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## **Motion Planning Using Adaptive Random Walks**

Stefano Carpin and Gianluigi Pillonetto

Abstract—We propose a novel single-shot motion-planning algorithm based on adaptive random walks. The proposed algorithm turns out to be simple to implement, and the solution it produces can be easily and efficiently optimized. Furthermore, the algorithm can incorporate adaptive components, so the developer is not required to specify all the parameters of the random distributions involved, and the algorithm itself can adapt to the environment it is moving in. Proofs of the theoretical soundness of the algorithm are provided, as well as implementation details. Numerical comparisons with well-known algorithms illustrate its effectiveness.

Index Terms—Adaptive sampling, holonomic path planning, stochastic processes.

### I. INTRODUCTION

In the last years, the problem of robot motion planning received even more attention as a consequence of two possibly synergistic events. Starting from the seminal work presented in [1], there has been a massive introduction of randomized motion planners (see, for example, [2]-[5]) which allowed effectively facing problems in high-dimensional spaces, not always solvable by formerly developed planners. Moreover, motion planners are currently used for a number of applications that are beyond traditional robotics (see [6]). In fact, the continuous flow of innovative tools for efficiently searching configuration state spaces confirms that as efficient planners are introduced, the frontier of possible applications is also being pushed further. For example, we cite problems like multirobot motion planning, digital characters motion generation, or protein folding and ligand binding. They exhibit problem instances with tenths or even hundreds of degrees of freedom (DOFs), and are currently tackled by using randomized motion planners.

Recently there has also been an impulse to study algorithms not only probabilistically complete, but also resolution complete, and to speculate on a possible turn back to efficient deterministic approaches [7], [8]. However, the randomized approach appears to be the current mainstream of research in motion planning.

In this context, we have developed a new motion planner based on random walks. In the past, random walks have been sporadically used and only in particular situations; for example, to escape local minima in potential field methods [9]. Instead, our approach relies entirely on this concept. Numerical results illustrate that difficult problems can be efficiently solved. The algorithm belongs to the class of single-shot motion-planning algorithms. This means that it does not preprocess the environment to subsequently solve many queries. Instead, it is well-suited for single queries. In addition, the planner is well-suited for the introduction of adaptive components in order to let it adjust its random components to better address the environment where it moves.

The paper is organized as follows. Section II formally introduces the problem and illustrates the algorithm. The theoretical soundness of the

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algorithm is discussed in Section III, while Section IV provides details about simulations and numerical results. Conclusions are offered in Section V.

### II. RANDOM-WALK ALGORITHM

The problem of robot motion planning is usually dealt with by using the configuration space approach [10]. This means that a search is conducted in the configuration space of the DOFs in order to plan a motion. Every point of this space corresponds to a placement of the robot in its physical workspace. As not all the points correspond to valid placements of the robot in its workspace, the configuration space is partitioned into two subsets, the free configuration space and the obstacle configuration space. This approach has proved itself to be highly effective, as it turns out that often in real applications, the configuration space is a smooth manifold [11]. Thus, it is possible to address a wide variety of problems using the same approach by designing algorithms which search for paths in these well-characterized spaces.

One instance of the robot motion-planning problem can be formulated as follows. Given an n-dimensional configuration space C, let  $C_{\mathrm{free}}$  be the subset of free configurations of C. Let  $x_{\mathrm{start}} \in C_{\mathrm{free}}$ , and let  $X_{\mathrm{goal}}$  be a target subset contained in  $C_{\mathrm{free}}$  whose Lebesgue measure is strictly positive. The task is to find a path connecting  $x_{\mathrm{start}}$  with  $X_{\mathrm{goal}}$ , i.e., a continuous function  $f:[0,1] \to C_{\mathrm{free}}$  such that  $f(0) = x_{\mathrm{start}}$  and  $f(1) \in X_{\mathrm{goal}}$ . It has to be outlined that this problem belongs to the PSPACE-hard complexity class [12]. This justifies the massive use of approximated and randomized algorithms.

Many randomized planners which have been developed so far draw independent samples from an a priori fixed probability density function whose support is the entire configuration space. The aim is then to build a graph or a tree which captures its connectivity [1], [13]. Instead, the algorithm we propose tries to find a solution by building a random walk growing from  $x_{\text{start}}$ . At every step, a new sample is generated in the neighborhood of the last point, in accordance with a Gaussian distribution. If the segment connecting them lies entirely in  $C_{\text{free}}$ , the point becomes the last point in the walk, otherwise it is discarded. The basic version of the algorithm is illustrated in Algorithm 1. The  $v_k$  used in line 5 indicates an n-dimensional Gaussian vector with zero-mean vector and covariance matrix  $\Sigma_k$ . We also denote with "Update" the function which makes the algorithm adaptive by computing  $\Sigma_k$  from a certain number of last-accepted samples, which in the algorithm is indicated as M. As the search of the path is carried out in a possibly high-dimensional configuration space whose topological shape is usually unknown, it can be nontrivial to fix the values of  $\Sigma_k$ . For this reason, the algorithm starts with an arbitrary covariance matrix  $\Sigma_0$ , and during the computation, a sequence of covariance matrices is generated through the "Update" function on the basis of the evolution of the computation. For example,  $\Sigma_k$  could be the covariance matrix of the last H accepted samples or a similar function of the recent history of the random walk being built. As it will be shown in Section III, the conditions required for the probabilistic convergence of the algorithm are very mild, so that a wide variety of update rules can be used while setting the values. Specific choices will be discussed in Section IV. The aim of the update step is to introduce adaptivity in the sampling distribution. In this way, the algorithm is able to modulate the sampling process in order to address the specific shape of the region of the environment that the random walk is currently exploring. One important aspect of the random-walk

<sup>1</sup>While using both uniform and Gaussian distributions yields effective planners, we will concentrate our discussion exclusively on Gaussian distributions. This is because in addition to exhibiting better performance and being more suited for adaptivity, it allows easy and elegant proofs of the probabilistic convergence of the algorithm (see Section III).

algorithm is that the time spent to generate a new sample does not depend on the size of the previously generated samples set, provided that the update of the covariance matrix can be done in constant time. This implies that the overall time spent to generate the samples is linear in the number of samples. This, for example, always holds if we process a fixed number of the samples accepted up to that point. This is different from most of the formerly developed algorithms, like probabilistic roadmaps (PRM) and rapidly exploring random trees (RRT), where the performance is worst than linear in the number of generated samples (see, however, [14] for a version of the RRT algorithm where the time needed to generate a new sample is logarithmic in the number of already generated samples). A natural technique to speed up the search process is to let the algorithm perform a bidirectional search. This means that two random walks are expanded, one growing from the start point and the other from the goal region, periodically verifying if it is possible to join them. In order to avoid the examination of all the samples generated, the connect trial is performed only between the last two generated points in the two walks. We have chosen this approach to keep the time needed to generate a new sample constant. As already observed, the expedient of bidirectional search allows a substantial gain in the performance (see [13]). Extensive experiments outlined that even this kind of bidirectional search improves the performance over the simple single-direction exploration. In addition to this, further opportunism can be introduced by letting the algorithm to periodically try to connect the last sample with the goal point (or the start point, if in a bidirectional search we consider the walk growing from the goal region). Another strategy we have designed to improve the performance of the algorithm consists of incorporating a greedy component. When the segment connecting  $x_k$  with new generated sample s does not entirely lie in  $C_{\text{free}}$ , this technique is activated by trying to expand the random walk along this segment. To be specific, a set of fine-grain equally spaced points  $\{s_1 = x_k, \dots, s_P = s\}$  lying on the aforementioned segment is generated. The greedy strategy then sets  $x_{k+1}$  to the point  $s_i$ , such that  $s_{i+1} \notin C_{\text{free}}$  and  $s_j \in C_{\text{free}} \ \forall \ 1 \leq j \leq i$  [13].

```
Algorithm 1 Basic Random-Walk-Based Motion Planner
1: k \leftarrow 0
2: x_k \leftarrow x_{\text{start}}
3: \Sigma_0 \leftarrow \Sigma_{init}
4: while NOT x_k \in X_{\text{goal}} do
5: Generate a new sample s \leftarrow x_k + v_k
6: if the segment connecting x_k and s lies entirely in C_{\text{free}}
     then
7:
         k \leftarrow k+1
         x_k \leftarrow s
8:
         \Sigma_k \leftarrow \operatorname{Update}(x_k, x_{k-1}, x_{k-2}, \dots, x_{k-H})
9:
10:
11:
           discard the sample s
```

When the algorithm terminates, i.e., when the last generated point is in  $X_{\rm goal}$ , or when the two chains can be connected, the sequence of segments connecting  $x_i$  with  $x_{i+1}$  (with  $i=0,1,\ldots$ ) indeed builds up a path which solves the problem. However, the quality of such a path is usually poor, because it includes a large number of useless motions. Indeed, the path obtained resembles a Brownian motion. A postprocessing stage is thus needed in order to smooth the generated trajectory. For this reason, we used a *divide and conquer* algorithm similar to binary search (see Algorithm 2, where the *pushback* operations used therein append the given point to the end of the list). See [15] for a similar approach. The smoothing procedure is repeatedly called until it is

not able to further smooth the trajectory. It is worth noting that due to the good performance of the algorithm itself, the smoothing step turns out to be extremely fast and few iterative steps are needed. This will be clearly illustrated in Section IV.

```
Algorithm 2 Solution Smoothing
1: SMOOTH(V, first, last, S)
2: INPUT: V vector of points to smooth
3: INPUT: first, last extremes of V to be optimized
4: INPUT/OUTPUT: S list with the smoothed sequence
5: if first = last then
6: S.pushback(V[first])
7: else if first = last - 1 then
8: S.pushback(V[first])
   S.pushback(V[last])
10: else if the segment connecting V[first] and V[last] lies
 entirely C_{\text{free}} then
11: S.pushback(V[first])
12: S.pushback(V[last])
13: else
     SMOOTH(V, first, (first + last)/2, S)
14:
     SMOOTH(V, (first + last)/2 + 1, last, S)
15:
```

### III. THEORETICAL FOUNDATIONS

In this section, we provide the formalism and the proofs about the probabilistic convergence of the random-walk algorithms introduced in the former section. Demonstrations will be given for what concerns the basic version, from which the convergence of bidirectional and greedy search naturally follows. See [16]–[18] for similar proofs in the context of PRMs. After introducing the terminology used, we will define the stochastic process that will be called adaptive random walk (ARW). Then, we will show that ARW allows moving between a certain kind of adjacent sets with a probability greater than a constant dependent only on the topology of the environment under consideration. Next, it will be shown that a path which connects the starting region with the goal region and that crosses a set more than once can be turned into a path which crosses the same set just once. These results will be then used to prove the probabilistic convergence of the new randomized motion-planning algorithm.

Let  $\Omega$  be a probability space with  $\omega \in \Omega$  [19], [20]. Let C be  $[0,1]^n$ , whose generic element is denoted x and equipped with the  $\sigma$ -algebra B(C) consisting of all the Borel sets in  $\Re^n$  (i.e., the open sets and their complements, as well as the countably infinite unions) contained in C. We will also denote with  $\mu(A)$  the Lebesgue measure of a generic set A in B(C). Moreover, we will denote  $N_{x,\Sigma}(y)$  the Gaussian probability density as function of y, having mean x and covariance matrix  $\Sigma$ .

Definition 1: We define  $B^+(C)$  as corresponding to the sets in B(C) whose measure is strictly positive with respect to  $\mu$ .

Hereby, we will denote with  $C_{\rm free}$  a fixed open set belonging to  $B^+(C)$ .

Definition 2: We call  $\{x_1, x_2\}$  an admissible couple if  $tx_1 + (1-t)x_2 \in C_{\text{free}}$  for  $t \in [0, 1]$ .

Let  $g: C \times C \to C$  such that

$$\begin{cases} g(x_1, x_2) = x_1 + x_2, & \text{if } \{x_1, x_1 + x_2\} \text{ is an admissible couple} \\ g(x_1, x_2) = x_1, & \text{otherwise.} \end{cases}$$

We define the discrete-time stochastic process  $\{X_k\}_{k=0,1,2,...}$  on  $(\Omega, \Gamma, \eta)$  and taking values on C, by the following recursive formulas:

$$\begin{cases} X_0(\omega) = x_{\text{start}}, & \text{with } x_{\text{start}} \in C_{\text{free}} \\ X_k(\omega) = g\left(X_{k-1}(\omega), v_k(\omega)\right), & \text{for } k = 1, 2, \dots \end{cases}$$
 (1)

where for every k the random vector  $v_k(\omega)$  is normal, zero-mean with covariance matrix  $\Sigma_k(\omega)$  (the dependence of a random variable on  $\omega$  will be hereby implicit). We call the stochastic process in (1) ARW.

Definition 3: We define the random vector  $H_k$  as corresponding to  $\{X_0, X_1, \ldots, X_k\}$ .

Assumption 4: For every k,  $v_k$  is independent from  $H_{k-1}$  once  $\Sigma_k$  is known. Moreover, every entry of  $\Sigma_k$  is a random variable which is measurable with respect to the  $\sigma$ -algebra generated by  $H_{k-1}$ , i.e., there exists for every k a function  $L_k: C^k \to \mathbb{R}^{n \times n}$  such that  $\Sigma_k = L_k(H_{k-1})$ .

We call  $L_k$  the learning rule of the random walk at instant k.

Assumption 5: For every k,  $\varepsilon_1 I_n \leq \Sigma_k \leq \varepsilon_2 I_n$ , where  $I_n$  is the  $n \times n$  identity matrix and  $\varepsilon_1$ ,  $\varepsilon_2$  are strictly positive numbers.<sup>2</sup>

Definition 6: Given  $H_{k-1}$  and assigned  $x \in C$  and  $A \in B(C)$ , we denote with  $P^r_{k,H(k-1)}(x,A)$  the r-step transition kernel of the chain, i.e., once known the history of the chain until instant k,  $P^r_{k,H(k-1)}(x,A)$  provides the probability that ARW takes value in A at instant k+r.

We next define the concept of visibility set [21].

Definition 7: Assigned a set q belonging to  $B^+(C)$  and contained in  $C_{\text{free}}$ , we denote with V(q) the maximal open set of points  $x \in C_{\text{free}}$  such that for every y belonging to q,  $\{x,y\}$  is an admissible couple.

Lemma 8: Let q be a set belonging to  $B^+(C)$  and contained in  $C_{\text{free}}$ . Then, given an arbitrary  $D \in B^+(C)$  contained in V(q), there exists a strictly positive m such that for every  $x \in V(q)$ , k and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^2(x,D) \ge m$$
.

*Proof*: By definition of V(q), we have that for every  $x \in V(q)$ , k, and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^{1}(x,q) \ge \int_{q} N_{x,\Sigma_{k+1}}(y)dy \ge l\mu(q) \doteq m_{1}$$

where l is a real number which uniformly bounds from below the function to be integrated on q (note that the existence of l comes from Assumption 5). Moreover, we also have that for every  $x \in q, k$ , and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^{1}(x,D) \ge \int_{D} N_{x,\Sigma_{k+1}}(y)dy \ge l\mu(D) \doteq m_{2}.$$

Combining the two inequalities, we have that for every  $x \in V(q)$ , k, and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^2(x,D) \ge m_1 m_2.$$

*Remark 9:* It is easy to note that when V(q) is convex, we could replace  $P^2_{k,H_{k-1}}(x,D) \geq m$  with  $P^1_{k,H_{k-1}}(x,D) \geq m$  in the statement of *Lemma 8*.

From this point on, if S is a subset of  $\mathbb{R}^N$ , we indicate with  $\overline{S}$  the closure of the set.

Definition 10: Let T and U be subsets of  $C_{\rm free}$ . We say that T and U are adjacent (and we write  $\operatorname{Adj}(T,U)$ ) when the following holds:

$$\overline{T} \cap \overline{U} \cap C_{\text{free}} \neq \emptyset.$$

Lemma 11: Let  $A_i$  and  $A_j$  belong to  $B^+(C)$  and to  $C_{\text{free}}$ , such that  $\operatorname{Adj}(V(A_i),V(A_j))$ . Then, there exists a strictly positive m, such that for every  $x \in V(A_i)$ , k, and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^{3}(x,V(A_{j})) \geq m.$$

<sup>2</sup>Where  $A \leq B$  means that B - A is positive semidefinite.

*Proof:* Since  $Adj(V(A_i), V(A_i))$ , there exists a point

$$x_0 \in \overline{V(A_i)} \cap \overline{V(A_j)} \cap C_{\text{free}}.$$

Then, since  $C_{\mathrm{free}}$  is open, there exists a ball  $B_{\varepsilon}(x_0)$  with radius  $\varepsilon$  and center  $x_0$  which is entirely in  $C_{\mathrm{free}}$ . We define the set  $D_1 = V(A_i) \cap B_{\varepsilon}(x_0)$ . Clearly, we have  $D_1 \in B^+(V(A_i))$ . Then, by applying *Lemma 8*, there exists a strictly positive constant  $m_1$  such that for every  $x \in V(A_i)$ , k, and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^2(x,D_1) \ge m_1.$$

Let  $D = V(A_j) \cap B_{\varepsilon}(x_0)$ . Again, by applying *Lemma* 8 and by considering that  $B_{\varepsilon}(x_0)$  is convex, there exists a strictly positive constant  $m_2$ , such that for every  $x \in D_1$ , k, and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^1(x,D) \ge m_2.$$

Composing the two inequalities, we have that for every  $x \in V(A_i), k$ , and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^3(x,D) \ge m_1 m_2$$

which completes the proof.

Definition 12: A triplet  $\{C_{\text{free}}, x_{\text{start}}, x_{\text{goal}}\}$  is a solvable instance of the motion-planning problem if the set

$$\begin{split} S(C_{\text{free}}, x_{\text{start}}, x_{\text{goal}}) &= \left\{ f: [0, 1] \to C_{\text{free}} \right. \\ &\quad \text{such that } f(0) = x_{\text{start}}, f(1) = x_{\text{goal}}, f \in C^0 \right\} \end{split}$$

is not empty.

Definition 13: Let  $\{C_{\text{free}}, x_{\text{start}}, x_{\text{goal}}\}$  be a solvable instance of the motion-planning problem, let  $f \in S(C_{\text{free}}, x_{\text{start}}, x_{\text{goal}})$ , and let A be a subset of  $C_{\text{free}}$ . We say that a solution f crosses Ar times if the set

$$t_A^f = \{t \in [0,1] \text{ such that } f(t) \in A\}$$

is the union of r disjoint nonempty intervals (either open, closed, or half open).

Lemma 14: Let  $\{C_{\text{free}}, x_{\text{start}}, x_{\text{goal}}\}$  be a solvable instance of the motion-planning problem. Let us suppose that there exists a finite sequence  $A_1, A_2, \ldots, A_k$ , where every  $A_i$  belongs to  $B^+(C)$  and is contained in  $C_{\text{free}}$ , such that

$$C_{\text{free}} = \bigcup_{i=1}^{k} V(A_i).$$

Then, there exists  $f \in S(C_{\text{free}}, x_{\text{start}}, x_{\text{goal}})$  that crosses at most once  $V(A_i)$  for  $i = 1, 2, \dots, k$ .

*Proof:* Let  $f^*$  be a solution of the solvable instance  $(C_{\mathrm{free}}, x_{\mathrm{start}}, x_{\mathrm{goal}})$  that crosses two times  $V(A_j)$ . Then, since  $V(A_j)$  is open

$$t_{V(A_j)}^{f^*} = \{t \in [0,1] \mid f^*(t) \in V(A_j)\} = (a_1, b_1) \cup (a_2, b_2).$$

Let  $p \in A_i$ . Starting from  $f^*$ , we define a new function  $f^{**}$  as follows:

$$f^{**}(t) = \begin{cases} f^*(t), & \text{if } t \le a_1 \\ f^*(a_1) \frac{b_2 + a_1 - 2t}{b_2 - a_1} + 2p\left(\frac{t - a_1}{b_2 - a_1}\right), & \text{if } a_1 < t < \frac{a_1 + b_2}{2} \\ p\frac{2b_2 - 2t}{b_2 - a_1} + f^*(b_2)\left(\frac{2t - a_1 - b_2}{b_2 - a_1}\right), & \text{if } \frac{a_1 + b_2}{2} \le t < b_2 \\ f^*(t), & \text{if } t > b_2. \end{cases}$$

By the definition of  $V(A_j)$ , it follows that also  $f^{**}$  is a solution of the instance  $(C_{\mathrm{free}}, x_{\mathrm{start}}, x_{\mathrm{goal}})$ . It is then evident that for each solution of  $(C_{\mathrm{free}}, x_{\mathrm{start}}, x_{\mathrm{goal}})$ , by repeatedly applying the former procedure, it is possible to build a new function which is also a solution, and indeed crosses at most once every set in the sequence  $V(A_1), \ldots, V(A_k)$ .

Theorem 15: Let  $C_{\text{free}}$  be connected and such that there exists a finite sequence  $A_1,A_2,\ldots,A_k$ , where every  $A_i$  that belongs to  $B^+(C)$  is contained in  $C_{\text{free}}$ , and

$$C_{\text{free}} = \bigcup_{i=1}^{k} V(A_i).$$

Then, for each  $x_{\text{start}} \in C_{\text{free}}$  and  $X_{\text{goal}}$  belonging to  $B^+(C)$  and to  $C_{\text{free}}$ , the algorithm ARW started in  $x_{\text{start}}$  will reach  $X_{\text{goal}}$  with probability 1.

*Proof*: Without loss of generality, we suppose that  $X_{\mathrm{goal}}$  is entirely included in one of the sets of the sequence  $V(A_1), V(A_2), \ldots, V(A_k)$ , denoted  $V(A_l)$ . In the light of *Lemmas* 8 and 11, there exists a strictly positive m, such that for every  $V(A_i)$  and  $V(A_k)$  with  $\mathrm{Adj}(V(A_i), V(A_k))$  and for every  $x \in V(A_i)$ , k, and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^{3}\left(x,V(A_{k})\right) \geq m$$

and for every  $x \in V(A_l)$ , k, and  $H_{k-1}$ 

$$P_{k,H_{k-1}}^2(x, X_{\text{goal}}) \ge m.$$

This, in addition to *Theorem 14*, allows us to conclude that there exist constants h and s, independent from  $x \in C_{\text{free}}, k$ , and  $H_{k-1}$  such that

$$\sum_{r=1}^{h} P_{k,H_{k-1}}^{r}(x, X_{\text{goal}}) \ge s > 0.$$

It follows that the probability that ARW has never entered the set  $X_{\rm goal}$  after Zh steps is less than or equal to  $(1-s)^Z$ . Clearly, when Z diverges, this probability goes to zero.

It is easy to observe that the last formula is similar to those found while analyzing the convergence of PRM or RRT. We outline that when dealing with the bidirectional search algorithm, as the two random walks are statistically independent, formerly introduced convergence results still hold. Moreover, for what concerns greedy search, all the lower-bound inequalities are still valid, and then again the convergence holds.

Remark 16: One of the most studied aspects of stochastic processes, in particular Markov chains, concerns their behavior in the long term (see, e.g., [22]). Investigations often concern the existence of a limiting distribution together with the way the process converges toward this distribution. To cite an important example, Markov chains are currently used for evaluating expectations of functions of interest under a target distribution [23]. In this context, one has to design the simulation so that it can be guaranteed that asymptotically the state of the process is exactly drawn from the desired distribution. Many of these theoretical aspects related to the above-defined ARW turn out to be very complex to study. This is due to the fact that we allow the transition kernel to adapt whenever new features of the configuration space are encountered during the run. As a consequence, it appears very hard to establish, e.g., if ARW possesses a limiting distribution and which form it could have. In fact, the adaptation occurs infinitely often, and the stationary distribution of the chain may be disturbed [24]. In the context of randomized motion planning, the interesting point is, however, that one has only to design chains able to quickly visit all the free

TABLE I							
TIME SPENT IN VARIOUS STEPS OF THE ARW ALGORITHM. DATA ARE AVERAGED OVER 200 TRIALS AND TIME IS EXPRESSED IN SECONDS							
Environment   Number of   Time to   Number of   Overall   Number							
Environment	Number of	Time	to	Number	of	Overall	Number

Environment	Number of	Time to	Number of	Overall	Number
	points in the	gener-	points in the	Planning	of d.o.f.
	random walk	ate the	$\operatorname{smoothed}$	Time	
		random	path		
		walk			
Car 1	182	0.0229	10.03	0.0313	3
Car 2	907.93	0.2844	16.79	0.3497	3
Car 3	132.35	0.0372	8.9	0.0556	3
Car 4	105.7	0.0088	7.17	0.0114	3
Wrench	315.81	0.5500	23.5	1.4738	6
Cage	436.06	0.7588	14.03	1.2225	6
Truck	1731.73	7.8384	19.57	10.9321	6
Coffemug	226.78	0.3161	8.65	0.4537	6
3drigid 1	18.83	0.0532	6.22	0.1246	6
3drigid 2	8077.51	7.03	21.46	7.6171	6
3drigid 3	1657.05	1.7477	12.14	2.0408	6
Multi1	14.32	0.0181	6.43	0.057	18
Multi2	931.35	0.9464	17.6	1.3102	12
Multi3	341.2	0.5778	22.31	1.55	48
Multi6	10703.06	3.6282	69.87	4.3298	6



Fig. 1. One of the environments provided with the MSL software. The truck has to be moved out of the cage.

configuration space, then prove their capability to reach the goal set with probability 1. This is exactly the purpose we have attained in this section.

### IV. SIMULATION DETAILS AND NUMERICAL RESULTS

The proposed algorithm with all its variants (greedy, nongreedy, bidirectional) has been implemented and integrated into the Motion Strategy Library (MSL) software developed at the University of Illinois, Urbana-Champaign [25], in order to compare it with different motion-planning algorithms over a set of standard problems. The MSL includes a wide range of versions of RRT-based motion planners, as well as the basic PRM motion planner. Problems included in the MSL involve simple planar motion-planning problems, as well as complicated three-dimensional (3-D) environments with possible multiple objects moving inside them (see Fig. 1 for an example). The planar problems involve 3 DOFs (x, y, and the orientation), while objects in the 3-D environments has 6 DOFs (three for the position, plus the yaw, pitch, and roll angles). MSL performs collision detection

using the Proximity Query Package (PQP) library developed at the University of North Carolina, Chapel Hill [26].

The adaptive rule we have used is the following (see [27]). Let

$$\overline{x}_k = \frac{1}{H} \sum_{i=k-H}^{k-1} x_i$$
 (2)

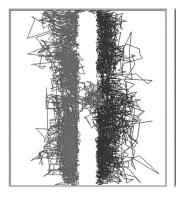
be the average of the last H accepted samples. Given a square p-dimensional matrix M, let  $m_{ij}$  be its generic element in position i, j. We define  $\operatorname{diag}(M)$  as follows:

$$\operatorname{diag}(M) = \begin{bmatrix} m_{11} & 0 & \cdots & \cdots & 0 \\ 0 & m_{22} & 0 & \cdots & 0 \\ \vdots & \cdots & \cdots & \vdots \\ 0 & \cdots & 0 & m_{p-1,p-1} & 0 \\ 0 & \cdots & \cdots & 0 & m_{pp} \end{bmatrix}$$
(3)

i.e., the matrix obtained from M by setting to 0 all the elements outside the main diagonal. The update rule for  $\Sigma_k$  is then

$$\Sigma_{k} = \max \left( \operatorname{diag} \left( \frac{1}{H} \left( \sum_{i=k-H}^{k-1} x_{i} x_{i}^{T} - H \overline{x}_{k} \overline{x}_{k}^{T} \right) \right), \Sigma_{\text{MIN}} \right)$$
(4)

where the function max returns a matrix whose generic element is the greatest of the corresponding elements in the two argument matrices, and  $\Sigma_{\text{MIN}}$  is a diagonal constant matrix with strictly positive elements on the main diagonal. It then follows that  $\Sigma_k$  is diagonal, too. The square roots of diagonal values of  $\Sigma_{\rm MIN}$  are set to one-sixth of the difference between the maximal and minimal values which can be assumed by the corresponding DOF.  $\Sigma_0$  is initially set to  $\Sigma_{\rm MIN}$ . As we are looking for local adaptivity, i.e., short-term memory, and for a fast update of the matrix  $\Sigma_k$ , history size is set to 10. We also have a posteriori assessed that this specific choice does not greatly influence the final results for values of H, say, less than 100. The computer used is a 2.0-Ghz Pentium IV with 512 Mbytes of RAM running Linux. All the subsequent numerical results refer to the bidirectional greedy planner, which turned out to be the most efficient. Table I reports the data relative to some the standard environments provided with the MSL. For lack of space, we can not give an in-depth description of every environment. However, we use the same names given in the MSL, so that the interested reader can refer to its documentation. Two aspects are evident. First, the time spent by the algorithm is linear in the size of the



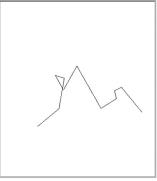


Fig. 2. Left figure illustrates a path generated using the random walk. Right figure illustrates the same path after the smoothing.

TABLE II
COMPARISON BETWEEN PRM AND ARW PLANNERS. TIME IS EXPRESSED IN
SECONDS AND DATA ARE AVERAGED OVER 200 TRIALS

Environment	PRM	ARW
Car 1	0.7447	0.0313
Car 2	1.3693	0.3497
Car 3	1.9961	0.0556
Car 4	0.3989	0.0114

generated random walk (compare columns  $T_1$  and  $N_1$ ). Second, even if the path produced by the random walk includes a great number of useless motions, significant improvements can be gained with the very fast postprocessing algorithm depicted in Section II. Fig. 2 illustrates an example of path smoothing. We then compare the random-walk algorithm performance with that concerning the basic PRM motion planner. For this aim, Table II compares the overall time spent by both algorithms to find the solution. We choose to display the comparison for just the four car-like environments, where the performance of the two is comparable. In more involved problems, there are usually two or three orders of magnitude of difference on the overall performance in favor of the ARW-based planner. It has, however, to be reminded that ARW is a single-shot algorithm, while PRM is oriented toward repeated queries over the same environment. Finally, we compare the performance of the random-walk planner with the RRTConCon algorithm provided in the MSL (see Table III). RRTConCon is a bidirectional single-shot greedy planner based on RRT [13]. Of course, comparing motion-planning algorithms experimentally using a restricted number of examples is hardly conclusive, in general. Moreover, both algorithms have been run without trying to find out the optimal values for the many parameters involved, and this could lead to different overall performance. What emerges, however, is that the performance of ARW appears interesting when dealing with single-shot problems.

## A. Importance of Adaptivity: Role of the History Size

We previously anticipated that the performance of the algorithm is not very sensitive to the history size. To prove this claim, we have run the algorithm with different history sizes and compared the results (see also [28]). We also ran the algorithm without adaptive components to verify the effectiveness of the proposed technique. In particular, due to its probabilistic nature, a maximum number of iterations is fixed. If the algorithm does not find a solution within that limit, this is considered a failure. Table IV illustrates the success ratio, i.e., the percentage of instances successfully solved within the given iterations bound. In all the cases, the initial matrix  $\Sigma_0$  was the same.  $\Sigma_0$  was fixed to be again a diagonal matrix whose elements were fixed to be the square of the difference between the maximal and minimal values, which can be assumed by the corresponding DOF. These values are fixed at will to

TABLE III
COMPARISON BETWEEN RRTCONCON AND RANDOM WALKS

Environment	RRT	ARW
Wrench	0.9824	1.4738
Cage	1.3908	1.2225
Truck	9.5486	10.9321
Coffemug	0.4825	0.4537
3drigid 1	0.0652	0.1246
3drigid 2	11.6726	7.6171
3drigid 3	14.0199	2.0408
Multi1	0.3842	0.057
Multi2	4.5345	1.3102
Multi3	84.2599	1.55
Multi6	6.3029	4.3298

be very large. It can be observed that without adaptivity, the algorithm suffers from drawing samples from a bad distribution. Instead, when adaptivity is included, the covariance matrix is updated and driven to more favorable values. This is indeed a big advantage, as it is then not necessary to invest a lot of time to tune the algorithm's parameters, since a rough initial choice can also be used. Table V instead compares the average time spent for different values of the history size. It can be observed that no significant variations emerge. The second important observation can be drawn from the execution times. We should remind readers that H = k means that the last k accepted samples are used in order to compute their variance. This means that at every iteration, it is necessary to process a  $k \times n$  data matrix. In spite of the increased size of such a matrix, we can observe that the time spent is more or less constant, and sometimes even decreases. Interestingly, this clearly indicates that the information acquired from the last k accepted samples is successfully used to quickly improve the sampling parameters. To make an example, when the variance of the increments through which the robot is currently moving is too large, so as to make many of the proposed movements fall outside the free configuration space, the algorithm is able to automatically reduce the variance of its steps, thus augmenting the probability of exploring a narrow space, e.g., a corridor, in a more effective and rapid way (see Fig. 3). An important point, which will deserve more attention in the future, is an accurate investigation in order to get clearer indications about the equilibrium point, i.e., to have indications about a range of values for H giving a minimum in the search times. This point is clearly nontrivial, since the expected time for finding the solution is, of course, a complex function of H which depends on the unknown environment where the robot moves. We also mention that we have repeated the simulations by computing the correlation between the last accepted samples, but it turned out that no significant speedup can be observed by using nondiagonal matrices. Of course, the design of more refined adaptivity rules could be important to further improve the performance of the proposed algorithm. For example, if knowledge about the C-space was available, it could be used in the choice of an adaptation rule for a more efficient path finding.

## V. CONCLUSIONS AND FUTURE WORK

We introduced the first random-walk-based adaptive motion planner. The algorithm builds a random walk with Gaussian increments over the configuration space. As the produced path can be very uneven, an efficient postprocessing step is used, yielding a fast smoothing of the produced path. A simple but effective technique for letting the algorithm adapt the random distribution parameters, in order to speed up the exploration process, has been devised. Thanks to this feature, it is possible to get a variable sampling resolution, which is indeed believed to be one of the most promising research directions. Rigorous proofs of the theoretical soundness of the algorithm have been provided under mild assumptions on the environment to explore and on the

Environment	No Adaptivity	H=5	H = 10	H=20	H = 30	H=40	H=50	H=100
Coffeemug	32	100	100	100	100	100	100	100
Cage	20	100	100	99	100	100	99	99
Wrench	10	61	61	70	59	60	65	68
Truck	0	91	88	88	89	94	90	89
3drigid1	100	100	100	100	100	100	100	100
3drigid2	0	95	83	80	90	86	88	80
3drigid3	28	100	100	100	100	99	100	100

TABLE V TIME PERFORMANCE OF THE ARW ALGORITHM FOR DIFFERENT VALUES OF THE HISTORY SIZE H. TIME IS MEASURED IN SECONDS. AVERAGE IS COMPUTED OVER 200 TRIALS FOR EACH BENCHMARK

Environment	H=5	H = 10	H=20	H = 30	H=40	H=50	H=100
Coffeemug	2.089	2.118	2.208	1.988	2.169	2.231	1.978
Cage	3.463	3.219	3.690	4.128	3.476	4.085	3.935
Wrench	8.816	7.954	8.809	7.603	7.857	8.243	8.063
Truck	38.277	35.605	30.095	38.317	36.251	34.542	31.422
3drigid1	0.355	0.400	0.389	0.369	0.392	0.400	0.388
3drigid2	14.908	14.660	14.484	14.287	14.893	14.901	13.515
3drigid3	3.564	3.413	3.065	3.035	3.173	2.946	3.106

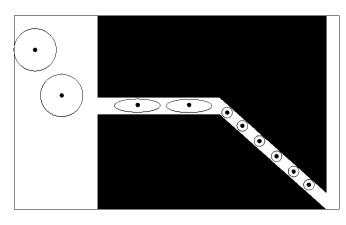


Fig. 3. When the robot moves in wide spaces (left), then a big covariance allows trying big steps. When it moves into the horizontal corridor, the covariance matrix indicates that big steps along the x axis are still possible, but not along the y direction. This is because proposed samples with big displacements along the y directions are refused, and then the corresponding matrix entry reduced. In the diagonal corridor, small steps in both directions will be taken, as samples with big steps in both directions will be rejected. On the other hand, if one allows for off-diagonal covariance matrix, then the covariance ellipsoid could align with the passage.

learning rule, which makes the algorithm adaptive. Numerical results clearly illustrate that for most of the investigated problems, the adaptive random-walk planner appears to be competitive. A direction for further development is the introduction of a bias to let the samples be generated not only according to the covariance matrix, but also taking into consideration unexplored regions. We are at the moment working on this extension. In the near future, we also plan to investigate the use of more refined adaptive techniques which could be useful while applying the algorithm in domains different from robot motion planning. For example, it appears particularly interesting to apply the algorithm to bioinformatics problems such as protein folding and ligand docking. In this context, we expect the adaptivity to be even more important, thanks to the continuous nature of the energetic levels concerning proteins instead of the Boolean nature of the configuration space characterizing robot motion planning. This will likely call for the study of more refined online learning techniques than those presented herein.

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