REASONING ABOUT UNCERTAINTY IN ROBOT MOTION PLANNING

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Anthony Lazanas
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I certify that I have read this thesis and that in my opinion it is fully adequate, in scope and in quality, as a dissertation for the degree of Doctor of Philosophy.

Jean-Claude Latombe
(Principal Advisor)

I certify that I have read this thesis and that in my opinion it is fully adequate, in scope and in quality, as a dissertation for the degree of Doctor of Philosophy.

Leonidas J. Guibas

I certify that I have read this thesis and that in my opinion it is fully adequate, in scope and in quality, as a dissertation for the degree of Doctor of Philosophy.

Oussama Khatib

Approved for the University Committee on Graduate Studies:

Dean of Graduate Studies

Abstract

Robot motion planning has been a central problem of robotics research since its early stages, yet no satisfactory solution has been found to date. The difficulty stems from the fact that the real world is too complex to be modelled accurately. The details that are omitted from a model create uncertainty. More detailed models have less uncertainty, but increase computational complexity; less detailed models may result in incorrect and/or incomplete planners. In this thesis, we investigate the effects of uncertainty on the difficulty of robot motion planning, and we study the tradeoff between physical and computational complexity.

Uncertainty makes the execution of motion plans non-deterministic. Sensing is used to help the robot identify its state, but measurement errors interfere with the process of state identification. Reasoning about potential (actual and perceived) states during plan execution is what makes planning with uncertainty a difficult problem. To keep the complexity of the problem polynomial, we have to represent all possible states with a limited number of equivalence classes. We present a model of a robot and its workspace, which yields a polynomial, correct, and complete motion planning algorithm. The key idea is the existence of regions in the workspace (landmark regions), where the robot has perfect position sensing and can navigate without error. Outside the landmark regions, the robot has no sensing at all and navigates with bounded control error. The proposed algorithm creates a distributed motion plan, by associating with every landmark region a motion command. This command is guaranteed to bring the robot into the goal, or into another region with a similar motion command attached to it.

Planning is performed using the preimage backchaining method. We extend

the standard definition of a "nondirectional preimage" (the set of all points of the workspace from where a robot can reliably achieve a goal set) to the case where a motion command depends on an arbitrary number of control parameters. We show that the resulting multi-dimensional preimage can be represented with a polynomial number of 2-D slices, each computed for a critical combination of values of the parameters. We present implemented algorithms for one parameter (the commanded direction of motion) and for two parameters (the commanded direction of motion and the directional uncertainty). The latter algorithm has many interesting applications, like planning in anisotropic uncertainty environments, finding probabilistic plans when no guaranteed plans exist, avoiding unexpected obstacles, etc.

Experimentation with the algorithm using a real mobile robot has been successful. By engineering the workspace, we have been able to satisfy all the assumptions of our planning model. As a result, the robot can operate for long periods of time with no failures. Since our planning algorithms are provably correct, experimental failures can be attributed to violations of the assumptions on which the algorithm depends. Recursive workspace engineering can be used until all assumptions are satisfied. Thus, experimentation and workspace engineering become essential tools for the creation of useful robot motion planning algorithms.

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Chapter 1

Introduction

In dealing with physical reality, which is too complex to be accurately modelled, one must decide on an appropriate modeling granularity level. The more details in the model, the bigger the complexity, and the harder the problems become. If too many details are omitted, then the class of solvable problems becomes uninteresting. Once a model of the real world has been selected, everything that has been omitted from the model becomes *uncertainty*. Dealing with uncertainty is the major problem of most systems aspiring to operate in the real world, and has been the most significant obstacle for the introduction of robots in everyday life. This work addresses the problem of dealing with uncertainty in robot motion planning. The general ideas presented, however, are applicable to any system that deals with the physical world.

1.1 About robots, motions, and planning

A *robot* is an autonomous device that is capable of changing the physical state of the world in which it operates, according to some guidelines specified by an external operator. The level of autonomy of a robot is determined by the level of details that it requires for its operation. For example, a robot at the human level of autonomy would accept commands expressed in ordinary human language; a superhuman robot would accept a single command, "Satisfy my wishes!," and would act accordingly. Since robots are neither human nor superhuman, guidelines must be expressed in an appropriately simple language.

As opposed to information, which requires very low energy and can be represented and propagated with electromagnetic interaction, a physical change requires a relatively high energy exchange. The simplest mechanism to deliver a burst of energy at a specific location is contact. Contact requires motion, so most operations of a robot involve some kind of motion. The study of robotic behavior is, to a large extent, the study of the motion of mechanical devices.

Every task, expressed in some input language that a robot accepts, must be translated into a collection of motions. Such a collection is called a *motion plan*. The procedure of deriving motion plans from more abstract task specifications is not an easy task for humans, so we require that robots perform it automatically. Two modules are usually involved in this process. First, a *task planner* translates commands expressed in a human-like language (e.g., "Get me a glass of water") into a set of geometric specifications (e.g., "Go to the refrigerator, grab the door, open the door, ..."). Then, a *motion planner* translates the high-level geometric specifications into a set of primitive motions, i.e., motions that can be easily effected with simple control of the motors that drive the robot.

The physical state of a robot and its environment can be described by a vector of parameters $\vec{\mathbf{p}}$. The geometric specifications of tasks can be described by a set of constraints $\mathcal{G}(\vec{\mathbf{p}})$. The initial state of the system can be described by another set of constraints $\mathcal{I}(\vec{\mathbf{p}})$. Motion planning amounts to the construction of an algorithm of primitive motions that transform $\mathcal{I}(\vec{\mathbf{p}})$ into $\mathcal{G}(\vec{\mathbf{p}})$, given some description of the effects of primitive motions on $\vec{\mathbf{p}}$. This formulation indicates that robot motion planning is just an instance of the more general problem of controlling the behavior of a system that obeys certain rules. Therefore, most of the ideas that we develop here have a much bigger domain of potential applications.

1.2 About uncertainty

In a perfect world, the result of primitive motions is deterministic. In the real world, however, nothing is deterministic. After all, the Second Coming may occur any time. Introducing uncertainty to the outcome of primitive motions complicates motion planning considerably. To begin with, a motion plan may not be a simple string of primitive motions any more. Since a primitive motion may lead the robot to a number of different states, which motion must be executed next may be a function of the resulting state. This observation introduces the need to sense the state of the system after the execution of each primitive motion. The sensing mechanism introduces its own errors: The identification of states may not be perfect. Because of that, a motion-planning algorithm must reason about the information that may be revealed after each primitive motion, as well as about the reliability of this information. As uncertainty grows, the number of states that may result from a primitive motion also grows, making the task of a motion planner more complex.

The deepest questions regarding uncertainty are philosophical in nature. Is it possible to predict the future given complete knowledge of the present? Is there such thing as complete knowledge? What constitutes knowledge? The first two questions are equivalent to the question regarding the existence of God, to which no satisfactory answer has be given to date. Adopting a utilitarian point of view, it is possible to give an answer to the third question. Knowledge is a collection of "words" in some representation language. Representation is the concept underlying knowledge. So the question regarding complete knowledge becomes a question about complete representation. Are there languages rich enough to represent the physical world? Even if we assume that such a language exists, we must use measurement in order to identify the actual state of the world in the space defined by the language. Measurement is subject to the limitations of measuring devices, and subject to Heisenberg's uncertainty principle. Finally, there is always the question whether the real world is a deterministic system or not, i.e., whether there exists some divine randomness.

Regardless of philosophical views, uncertainty is a major hindrance in the attempt to control physical systems. Its existence must be acknowledged and dealt with. Various approaches can be used in dealing with uncertainty.

- **Ignoring:** The simplest approach is to proceed as if uncertainty were not present and converge to the goal state with iterative process execution. The advantage of this method is that it is easy to design and execute. The disadvantages are (a) that it may often fail, (b) that, even when successful, it may take a big number of iterations, and (c) that it is hard to explain failure and derive useful conclusions from it.

- **Engineering:** Another approach is to attempt to eliminate as many sources of uncertainty as possible by simplifying the structure of the physical system and aligning it with the capabilities of the measuring devices. Care must be taken that the engineering of the system do not interfere with its ability to perform its assigned duties.
- **Reducing:** A third approach is to use a richer representation language. The more detailed the description of a system, the less uncertain its operation is. However, using a rich representation language increases computational complexity and may make problem solving hard.
- **Reasoning:** Uncertainty that has not been ignored, eliminated by engineering, or reduced by representation, becomes a parameter of the system model. To predict the behavior of the system we must explicitly reason about uncertainty.

1.3 Problem solving methodology

Problem solving in the real world has four phases. The first phase (*engineering*) is an attempt to reduce the physical complexity of a given problem, by means of physical alteration of the characteristics of the problem. Irrelevant details are eliminated, helpful features are added, and mechanisms are devised and implemented, so that the original problem becomes easier to solve. During the second phase (*modeling*), the simplified actual problem is transformed into a formal problem through the choice of a representation language and a solution language. The formal problem is the projection of the real problem onto the space defined by the representation and the solution language. The physical complexity of the real problem is transformed into the computational complexity of the formal problem. The choice of the representation and the solution language is a major factor in determining the computational complexity of the formal problem. During the third phase (*computation*), a mapping between the representation language space and the solution language space is defined

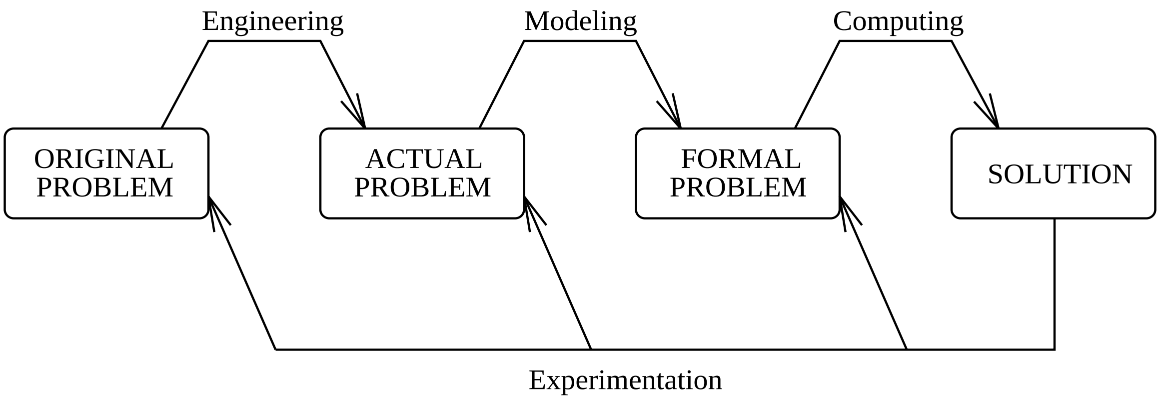

Figure 1.1: Problem solving phases

as the solution of the formal problem. An algorithm is a procedure that attempts to compute the above mapping. Finally, *experimentation* is used to evaluate the quality of a problem solver. Failures are attributed to deficiencies of engineering, modeling, or computation, which are accordingly adjusted to increase the effectiveness of the problem solver.

The quality of a problem solver is evaluated on three grounds: *correctness*, *completeness*, and *complexity*. Correctness measures the ability of the problem solver to produce solutions which are actual solutions of the original problem. Completeness measures the percentage of actual solutions which can be produced by the problem solver. Complexity is judged by how fast the solver generates the solutions, as a function of the size of the input problem.

Even though the above attributes can be associated with all phases of problem solving, they are only used to evaluate the computation phase. The reason is that it is easier to quantify the attributes using the description language of the formal problem. The computational complexity of a formal problem is defined as the lowest possible complexity of a correct and complete algorithm that solves the problem. The computational complexity of a formal problem depends on the physical complexity of the underlying physical problem, as well as on the representation and solution language.

The term "epistemic adequacy" has been used to describe how well a formal problem approximates a real one. The choice of the representation and the solution language determine the epistemic adequacy of modeling. Although high epistemic adequacy is desirable, it usually results in formal problems that cannot be solved with algorithms that are correct, complete, and of low complexity. In dealing with hard problems, an important decision to be made is how many physical details of the original problem we are willing to hide through engineering and modeling, in order to produce a formal problem with sufficiently low complexity.

1.4 The philosophy of the proposed approach

Robot motion planning in the real world is a difficult problem mainly because of the existence of uncertainty. Two major schools of thought have evolved. The "reactionists" ignore uncertainty, with the claim that it is too expensive to reason about it. Motion planning is performed as if no uncertainty were present. Robots are equipped with a set of predefined local behaviors to help them react to the eventual failures during plan execution. The "reasoners" model uncertainty (or some part of it) and devise algorithms that reason explicitly about it, so plans do not fail as often. However, the models that have been used to describe uncertainty result in intractable formal problems. To date, no universally acceptable solution has been given to the problem of robot motion planning with uncertainty. Algorithms are either too slow or unacceptably incorrect.

The role of engineering Both approaches ignore the value of problem engineering as a means to reduce the complexity of the motion-planning problem. In this work, we begin from the premise that we need motion-planning algorithms which are correct, complete, and have computational complexity that is a polynomial function of the size of the formal problem. We use workspace and robot engineering to produce a real problem that can be modeled into a formal problem, which in turn can be solved with algorithms having the above properties. In our approach, engineering is not confined to creating robots able to cope with a given environment, but it attempts to adapt the environment to the capabilities of robots (much in the same way that paved roads are constructed to help cars navigate). Therefore, our approach is geared towards applications in environments that can be controlled (at least to some extent), and not towards purely exploratory applications.

The role of experimentation By insisting on the correctness of algorithms, we are able to redefine the role of experimentation. When an algorithm is provably correct, experimental failures can be directly attributed to the violation of one or more modeling assumptions. Then, we can decide whether we can augment the used model (always retaining the desired properties of the algorithm), or whether further engineering is necessary. Thus, workspace engineering enters the problem-solving loop, as experimental feedback can be used to provide guidelines for engineering.

Modeling requirements The correctness of a motion planning algorithm means that the plans it produces are guaranteed to succeed. A motion plan can be guaranteed in the presence of uncertainty only if uncertainty is bounded. If anything can happen, no plan is guaranteed to succeed. Therefore, we represent stochastic quantities with distributions that are bounded over some domain. Since we are only interested in the correctness of a plan (and not in its stochastic attributes), we do not reason at all about the characteristics of the probability distributions; rather, we only use their bounded domains as a sufficient representation.

To solve the motion-planning with uncertainty problem in polynomial time, it is necessary that the planner only consider a small finite number of world states. We embed such a discrete structure into the robot workspace by introducing uniquely identifiable characteristics called *landmarks*. Each state is characterized by the perception of a unique landmark. All other information is shielded out of the planning model to preserve its polynomial complexity. Each landmark is associated with a region of the workspace (a landmark region) which allows a robot to navigate in it without loss of information. Uncertainty increases during navigation outside such regions and decreases when the robot enters one of them. Therefore, the role of the planner is to make sure that every motion of the robot will eventually terminate within some landmark region. Such a design allows the robot to navigate while keeping its uncertainty bounded. This approach has distinct similarities to the binary transmission of signals, where a limited representation language (0 and 1) is used to preserve the quality of data.

1.5 Overview of the thesis

In Chapter 2 we present a very general model of a system that is controllable through a set of primitive commands, and that is capable of affecting the physical state of the world in which it operates. We begin with a very simple example, and then we gradually build on it introducing many important issues and structures. We study planning under uncertainty, and we show how we can derive guaranteed plans by bounding uncertainty over some domain. We discuss perception and interpretation of perceived data, and we demonstrate how a planner must explicitly reason about information that may become available at execution time. The more information that is available, the more complicated the planning task is. Therefore, it is of utmost importance to select the right pieces of information: those pieces that enable a planner to solve most tasks in an efficient manner. Too much information may be unnecessary, too little information may be inadequate. We give several versions of planning and execution algorithms. Since it is difficult for a planner to produce plans that react to the sensed data at a very fast rate, we decouple the motion control loop from the plan-execution control loop and we introduce the notion of continuously executed commands and the termination condition. Finally, we present high-level definitions of the forward projection, the preimage, and the kernel, as tools for producing guaranteed plans under uncertainty.

In Chapter 3 we present the general problem of robot motion planning and, more specifically, robot motion planning under uncertainty. After an overview of the literature in the field, we introduce the notion of the configuration space and formally define the planning problem. Then we present the preimage backchaining planning technique and an implementable version in two dimensions, backprojection from a target kernel. We consider three kinds of sensing, position sensing, force sensing and sticking on obstacle edges, and we show that their combined use may have a nonlinear effect on the capabilities of a planner. Several actual planning examples, solved with the implemented planner are presented. We also present an algorithm to compute the maximal preimage of a convex set, and we show that a concave set may have no unique maximal preimage. By combining two-dimensional preimages computed for all possible commanded directions of motion, we create three-dimensional omnidirectional and nondirectional preimages; these are essential building blocks of complete polynomial planners under uncertainty.

In Chapter 4 we discuss in detail how the amount of information considered by a planner affects the complexity of the planning algorithm. Practical planners must be fast and, therefore, must select carefully what information to consider. Since the appropriate information may not always be available in the workspace, we propose the idea of workspace engineering for planning complexity reduction. We introduce the notion of landmarks as sources of information, and the notion of landmark regions as information fields in the workspace. By assuming a finite number of such fields, we make the first step towards polynomial planning. We give a formal definition of the problem of landmark-based navigation and a correct, complete, and polynomial algorithm for it. The algorithm is based on the fact that a nondirectional preimage can be represented with a data structure of polynomial complexity. This polynomial description consists of a collection of slices (two-dimensional preimages) computed at certain critical values of the commanded direction of motion d. The critical events that produce the critical values of d are listed and explained in detail. We present several examples of the algorithm. At the end of the chapter, we explain why the motion plans produced by our planning algorithm have many desirable properties. Specifically, in addition to having been produced by a correct, complete, and polynomial algorithm, the motion plans are optimal in the number of executed steps, distributed over the workspace, and robust (i.e., it is easy to reattach to a plan that has been disrupted by an unexpected event).

In Chapter 5 we discuss several shortcomings of the landmark-based navigation algorithm and we attempt to overcome them. We show that it is straightforward (a) to allow generalized polygonal objects (instead of simple disks) in the configuration space, (b) to consider compliant motions, and (c) to allow contact between landmark regions and C-obstacles. We also show that it is easy to precompute all possible plans in a specific workspace. The assumption that is the hardest (actually impossible) to satisfy is that sensing and control are perfect within landmark areas. To deal with this issue, we introduce the notion of generalized landmarks, where the robot is only required to be able to get reliably within a subregion of the landmark region. Planning with generalized landmarks is actually much easier than planning with our standard definition of landmarks. We then attempt to relax the constraint that all landmarks must be recognizable from each other, and present techniques that deal with confusable landmarks. We conclude that confusable landmarks introduce too many complications and, in general, lead to loss of polynomiality. Therefore, making landmarks individually recognizable is an important goal of workspace engineering. Finally, we give a graph representation of planning with uncertainty. We are able to represent with a graph all possible plans, including guaranteed plans, plans that always fail, and plans that may fail or succeed. An important family of plans are the probabilistically guaranteed plans (introduced by Erdmann in [34]), which are plans that are guaranteed to succeed, but the number of execution steps is not bounded. (The expected number of execution steps, though, is a finite number.) Many problems that do not accept guaranteed plans accept probabilistically guaranteed plans. We give a recursive algorithm to compute such plans within our landmark framework, but we make no attempt to analyze its complexity.

Up to this point we have assumed that the robot controls a single parameter, the commanded direction of motion d. In Chapter 6 we consider the situation where more than one parameters are changeable, either under the control of the robot, or not. In this case, the two-dimensional preimage depends on many parameters, and the notion of nondirectional preimage is generalized to the multi-parametric preimage. We show that the ideas of Chapter 5 can be easily extended to higher-dimensional spaces. As a specific case, we study planning where the directional uncertainty half-angle θ is controllable by the robot. We wish to produce guaranteed plans using the highest possible values of θ, since, presumably, some kind of cost must be paid in order to reduce uncertainty. We develop a correct, complete, polynomial, maximum-uncertainty one-step planner, as well as a multi-step one that maximizes the value of θ at each step. We show that these algorithms can be also used in the case where θ

is not controllable by the robot, for problems that do not accept guaranteed plans. In this case, the algorithms produce non-guaranteed plans that have the maximum likelihood of success. The algorithms have been implemented and their use is demonstrated with several examples. Because non-guaranteed plans may fail in undesirable ways, we explore ways to produce plans that either succeed or fail recognizably (Donald's EDR plans [23]). Finally, we present many practical problems that can be solved using the multi-parametric planning technique (planning with anisotropic directional uncertainty, unexpected obstacle avoidance, controllable sensor sensitivity, robot cooperation, and directional sensing).

In Chapter 7 we test our algorithms by experimenting with them on a real robot. We use NOMAD-200, a small omnidirectional cylindrical robot from Nomadic Technologies. We conduct two sets of experiments. In the first set we use prismatic landmarks standing in the workspace that are detected by the reflection of a plane of laser light on them. In the second set of experiments we use landmarks placed on the ceiling that are detected by a camera fixed on the robot and looking straight up. We describe in detail the computation of the induced landmark regions, and the detection, recognition and localization algorithms. In the first set of experiments, not all of our model assumptions are met and the execution of plans is not 100% reliable. In the second set of experiments, we use iterative workspace engineering (especially with respect to the location of landmarks) to eliminate as many sources of errors as possible. Since we cannot completely eliminate uncertainty in the landmark regions, we make use of the generalized landmarks algorithm. The result is a system that operates reliably until the robot runs out of power (a little less than an hour). This is possible because in our system errors do not deteriorate over time, but always remain within their assumed bounds. We conclude that, with relatively little effort and resources, it is possible to create a planning system that remains reliable over time.

In the last chapter, we summarize the conclusions from this work, we identify promising areas for future research, and we discuss questions raised by critics of our approach.

Chapter 2

Models of Perception and Action

Before we can solve problems in the physical world, we must first build an adequate representation of it. This representation, called a *world model*, is a projection of the physical world onto the reasoning space. The projection must maintain all important characteristics of the world (relative to the problem we wish to solve), but it should conceal all parameters that are irrelevant to the problem. Selecting an appropriate world model greatly affects the existence and the complexity of problem solving algorithms.

In this chapter we consider a very generic dynamic world model. An agent (robot), obedient to commands from a controlling system, operates in the world. The agent acts as the interface between the controlling system and the world, by providing the means to affect the world state, and also by providing sensory feedback. We begin with a very simple discrete model with no uncertainty. Then we consider uncertainty both about the present (initial state uncertainty, perceptual uncertainty) and the future (control uncertainty). We introduce several functions that facilitate planning with uncertainty, and we sketch a planning algorithm. Finally, we consider systems with two distinct control loops, the motion control loop and the plan execution control loop, and we introduce the notion of the "termination condition."

2.1 Action

Consider a robot $\mathcal{R}$ operating in a workspace $\mathcal{W}$. Let the vector s contain all the parameters that are needed to describe the combination of the robot and the workspace. We will refer to s as the *state* of the system. Let c denote any *primitive command* that the robot can execute. The function **act** describes the results of the execution of c, i.e., it describes how s changes in response to the execution of c. Assume for the moment that this is the only way in which s may change, and that the system is completely deterministic.

Let s_0 represent the initial state of the system. A *plan* is a sequence of commands $c_0, c_1, \ldots, c_{n-1}$, which, if executed, will cause the system to pass through states $s_0, s_1, \ldots, s_{n-1}$ and finally arrive at s_n, where $s_i = \mathbf{act}(c_{i-1}, s_{i-1})$, $i = 1$ to n. Plans are provided by a user[1] trying to achieve goals with the help of the robot. A goal is defined as a set $\mathbf{G}$ of desirable final states of the system. *Planning* is the procedure which takes as inputs s_0, $\mathbf{G}$ and **act** and returns a sequence of primitive commands, i.e., a plan. The trace of an execution of a plan is the trajectory of states $s_0, s_1, \ldots, s_n$ that the system traverses during this execution. An execution is successful if $s_n \in \mathbf{G}$.

As an example, consider a point robot that can move on a straight line. Let the set of primitive commands that are available to the user be $\{F, B\}$, and let the function **act** be defined as follows:

$$\mathbf{act}(c, s) = \begin{cases} s+1, & \text{if } c = F; \\ s-1, & \text{if } c = B. \end{cases}$$

If $s_0 = 0$ and $\mathbf{G} = [2.5, 5]$, then plans (F, F, F) and (F, F, F, F) both achieve the goal, the first one in three steps and the second one in four. Generally, plans with fewer steps are preferable, unless there are other considerations like robustness. For example, plan (F, F, F, F) has a margin of error equal to 1, and might be preferable to plan (F, F, F) which has a 0.5 margin of error.

[1] A person, another robot, or a software program.

2.2 Uncertainty

Planning in deterministic systems is straightforward, though not always easy. Unfortunately, real systems are never that well behaved. There is always a stochastic aspect in their behavior. We call this stochastic aspect *uncertainty.* While in fully deterministic systems a plan has a unique execution trace, this is not the case in stochastic systems. We call a plan *guaranteed* if its executions are always successful. The probability of success of a non-guaranteed plan is the percentage of its successful executions.

Uncertainty about the initial state: Let us first consider the case where the initial state is not precisely known at planning time. Thus, instead of a single value s_0, the planning procedure knows a probability density function $f_0(s)$ about the possible initial states of the system. We introduce a new function **ACT** that takes $f_i(s)$ as input (instead of s), and returns the updated distribution of s after the execution of c_i: $f_{i+1}(s) = \mathbf{ACT}(c_i, f_i(s))$. Now the success probability of a plan is equal to $\int_{\mathbf{G}} f_n(s)\, ds$.

Returning to our simple example, assume that s_0 is not known precisely, but it is known to be somewhere in [-0.6,0.4] with uniform probability. If $u[a,b]$ denotes the uniform distribution on [a,b], then one possible form of **ACT** is the following:

$$\mathbf{ACT}(c, u[a,b]) = \begin{cases} u[a+1, b+1], & \text{if } c = F; \\ u[a-1, b-1], & \text{if } c = B. \end{cases}$$

With the goal being [2.5,5] the plan (F, F, F), is not guaranteed. It maps $u[-0.6, 0.4]$ into $u[2.4, 3.4]$, so its probability of success is 90%. However, the plan (F, F, F, F) is guaranteed.

Control Uncertainty: As long as **act** is deterministic, the variance of $f_i(s)$ remains constant and equal to the variance of $f_0(s)$ throughout the execution of a plan. In reality, however, this is rarely true. The mapping **act** is stochastic, causing the variance of $f_i(s)$ to grow as i grows. We call this fact *control uncertainty.*

In our example, let us assume for simplicity that **ACT** retains the uniformity of the distribution, but stretches it a bit, like for example:

$$\mathbf{ACT}(c, u[a,b]) = \begin{cases} u[a+0.8, b+1.2], & \text{if } c = F; \\ u[a-1.2, b-0.8], & \text{if } c = B. \end{cases}$$

Plan (F, F, F) maps the initial distribution $u[-0.6, 0.4]$ into $u[1.8, 4]$; its probability of success is equal to 15/22. But now, plan (F, F, F, F) also becomes non-guaranteed, because it results in a final distribution of $u[2.6, 5.2]$ that is not fully included in the goal; its probability of success is equal to 24/26.

2.3 Perception

In the above example, one could note that the plan (F, F, F, F) is guaranteed to make the robot *reach* the goal, but not to make the robot *stop* in the goal. Indeed, after three steps the distribution of s is $u[1.8, 4]$. If $s \leq 3.8$ the robot will get (and stop) in the goal after the next step. If $3.8 < s \leq 4$ the robot is already in the goal, but it may leave it with its next step. A successful plan must not only make sure that the system achieves a goal state, but that it also remains in a goal state after the execution of the plan.

If the robot had a way to verify goal achievement, we could derive the following guaranteed plan: "**while** (*not in the goal*) **execute** F." This is an example of a generalized plan structure, where the plan ceases to be linear and becomes an algorithm that is parsed and executed by the robot. In order to facilitate decisions during the plan execution, the robot observes the state s and transforms it into a variable $\hat{s}$ that is stored internally or fed to the execution module directly. This transformation is called *perception* and is modelled with the function $\hat{s} = \mathbf{per}(s)$. However, decisions like goal achievement, need to be functions of the actual state of the world s. Therefore, we need to be able to obtain estimates of the actual state s from the perceived state $\hat{s}$. Let us call this mapping *interpretation*: $s = \mathbf{int}(\hat{s})$.

Perceptual Uncertainty: If the perception function is one-to-one, then interpretation is a one-to-one function, too. Each value of s corresponds to a unique value of

$\hat{s}$. Unfortunately, this is never the case in the real world. First, perception is a lossy procedure: we cannot possibly record all parameters of a real system. Secondly, **per** is not even deterministic because of measurement errors; it is a stochastic function. In general, both perception and interpretation are many-to-many mappings. We call this fact *perceptual uncertainty*. We can model perceptual uncertainty, with the introduction of functions **PER** and **INT** that accept probability distributions as arguments, and return probability distributions: $f_{per}(\hat{s}) = \mathbf{PER}(f(s))$ and $f_{int}(s) = \mathbf{INT}(f(\hat{s}))$.

Back to our example, let us assume that: $\mathbf{PER}(u[a,b]) = u[a-0.1, b+0.1]$ and $\mathbf{INT}(u[a,b]) = u[a-0.1, b+0.1]$. The straightforward condition "*not in the goal*" must now be transformed either into a condition on $\hat{s}$, or, equivalently, into a condition on all possible interpretations of $\hat{s}$.

Consider the plan "**while** ($\hat{s} < 2.6$) **execute** F." Figure 2.1 demonstrates the execution of this plan. Dark-shaded areas correspond to all possible actual locations of the robot (derived only by the knowledge of the initial state of the robot prior to the execution of a command, and the accuracy of the controller). White boxes represent all possible actual readings of the robot's sensor.

In the first line of the figure the initial situation is shown: The robot lies somewhere in $[-0.6, 0.4]$ and its sensor may read anything in the range $[-0.7, 0.5]$. Since all possible sensor readings are lower that the cutoff value of 2.6, the robot will necessarily take a step forward. In fact, the robot will have to take two more steps forward before its sensory reading may exceed 2.6.

The situation is depicted in the second line of Figure 2.1. After three steps, s may lie anywhere in $[1.8, 4]$ and $\hat{s}$ anywhere in $[1.7, 4.1]$. If $\hat{s} \geq 2.6$, the execution of the plan will stop. However, if $\hat{s} \in [1.7, 2.6)$, the execution will proceed for at least one more step. If this happens, using the interpretation function and our prior knowledge about s, we know that s must lie in $[1.8, 2.7)$, as shown in the third line of Figure 2.1.

The fourth line of the figure depicts the situation after such a fourth step: The range of values of s is $[2.6, 3.9)$, and the range of values of $\hat{s}$ is $[2.5, 4)$. Again, if $\hat{s} \geq 2.6$ the execution will halt, otherwise it will proceed for a fifth step; s will start in $[2.6, 2.7)$ (line 5 of Figure 2.1), and will end up in $[3.4, 3.9)$. The new $\hat{s}$ will lie in $[3.3, 4)$, so the execution will terminate here (line 6 of the figure). In all cases,

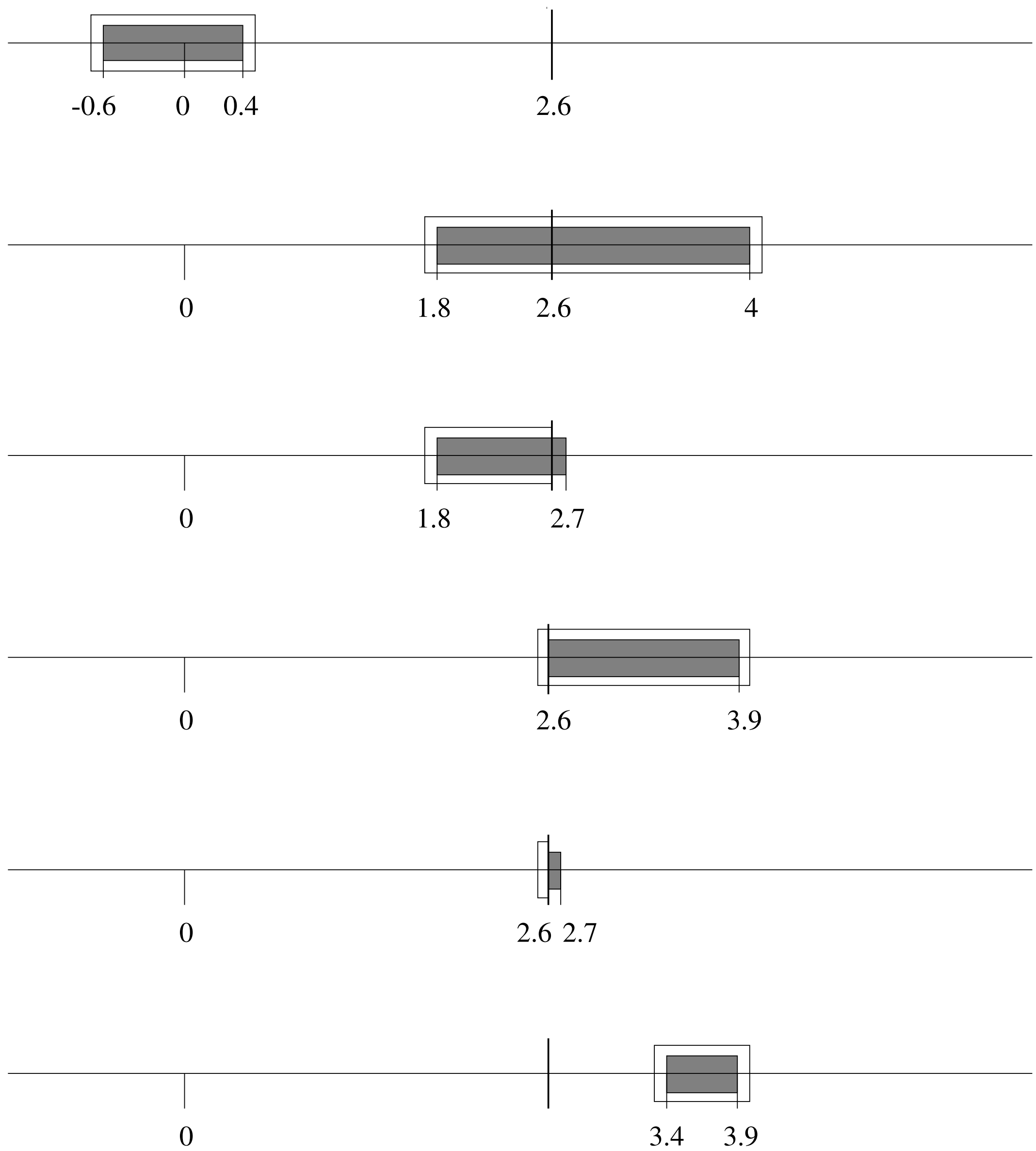

Figure 2.1: Planning with perceptual uncertainty

after the plan terminates, the system will be in a goal state $[2.5, 5]$, i.e., the plan is guaranteed.

Note that it is possible to terminate the execution of the plan after the fourth step, because we know that the actual location of the robot lies in $[2.6, 3.9]$, i.e., necessarily in the goal. However, since our termination condition is based solely on $\hat{s}$, it is not possible to do so. As we will see in the following chapters, it is much harder to derive plans that take into consideration the derived actual state of the robot.

Another interesting observation is that during plan execution the robot will have more accurate knowledge of its actual location s, because, after each step, the interpretation function will narrow down the possible values of s into a 0.2-wide interval. However, during the planning phase we need to take into account *all* possible situations that may result during plan execution, because we are interested in guaranteed plans. This observation can be extended to argue that while the planning module must use appropriate models for fast and guaranteed planning, the execution module may use different (yet compatible), more accurate models, so that the execution of plans becomes more efficient.

2.4 Deriving one-step plans

2.4.1 Algorithmic formulation of plan execution

Planning with uncertainty is quite complicated. In order to come up with a problem that we can solve rigorously, we will make a few simplifying assumptions. As a first step, we will not use probability density functions to represent beliefs about the state of the system. Instead, we will assume that beliefs are limited to the knowledge of a set, representing all values of s currently believed possible. We thus represent a density function $f(s)$ with the set $F = \{s \mid f(s) > 0\}$. According to this representation, $\mathbf{PER}(S)$ represents the set containing all values of $\hat{s}$ that can be observed, if s can take any value in S; $\mathbf{INT}(\hat{S})$ is the set containing all values of s that are compatible with a measurement $\hat{s} \in \hat{S}$; and $\mathbf{ACT}(c, S)$ is the set containing all possible values of s after the execution of command c in a system starting at any state $s \in S$. We also introduce the function $\mathbf{dec}(S, P)$ which returns the command prescribed by plan P when the state of the system can take any value $s \in S$. Now we can use the following algorithm to describe the execution of a plan P in a system which starts at an initial state known to be in a set S_0:

$\hat{s} \in \mathbf{PER}(S_0)$ is the perceived initial state of the system;
$\tilde{S} \leftarrow \mathbf{INT}(\{\hat{s}\}) \cap S_0$;
while $(\tilde{S} \not\subseteq \mathbf{G})$ **do** {
    $c \leftarrow \mathbf{dec}(\tilde{S}, P)$;
    execute c;
    $S \leftarrow \mathbf{ACT}(c, \tilde{S})$;
    $\hat{s} \in \mathbf{PER}(S)$ is the perceived new state of the system;
    $\tilde{S} \leftarrow \mathbf{INT}(\{\hat{s}\}) \cap S$;
}
return *success*;

2.4.2 The preimage and the kernel

Suppose we are interested in finding a *one-step* plan, i.e., a plan, all the executions of which will traverse the loop of the execution algorithm just once. The question whether a one-step guaranteed plan exists can be formulated as follows:

$$\forall \hat{s} \in \mathbf{PER}(S_0),\ \exists c:\ \mathbf{ACT}(c, \mathbf{INT}(\{\hat{s}\}) \cap S_0) \subseteq \mathbf{G}$$

Essentially, deriving a one-step plan entails associating a command c with all possible values of $\hat{s}$. This number is always finite (since $\hat{s}$ is the output of a physical instrument), but it may be very large. It would be a tedious and time consuming task to have the planner reason explicitly about each of these values separately.

Fortunately, there is a better way. Consider the *preimage* function (**PRE**), which operates inversely to **ACT**: For a given set of system states S and a command c, it returns the *maximal* set of system states S_P such that: $S_P = \mathbf{PRE}(c, S) \Rightarrow \mathbf{ACT}(c, S_P) \subseteq S$. In other words, the preimage of a set S of system states for a command c is another set of system states S_P with the following properties: a) If c is executed with the system starting in any state in S_P, the resulting state is guaranteed to be in S. b) There is no set with the above property that is not a subset of S_P. An interesting property of the preimage, which follows directly from its definition, is that the preimage of a subset of a bigger set is a subset of the preimage of the bigger set for the same command.

If computation of preimages is possible, then the following algorithm can be used to solve the one-step planning problem:

$S \leftarrow S_0$;
$\hat{S} \leftarrow \mathbf{PER}(S_0)$;
while $(\hat{S} \neq \emptyset)$ **do** {
  select a new command c that has not been already processed;
  if (*all commands have been processed*) **return** *failure*;
  $P \leftarrow \mathbf{PRE}(c, \mathbf{G})$;
  $S^+ \leftarrow S \cap P$;
  $S^- \leftarrow S \setminus P$;
  $K \leftarrow S^+ \setminus \mathbf{INT}(\mathbf{PER}(S^-))$;
  if $(K \neq \emptyset)$ **associate** c with each value $\hat{s} \in \mathbf{PER}(K)$;
  $S \leftarrow S \setminus K$;
  $\hat{S} \leftarrow \hat{S} \setminus \mathbf{PER}(K)$;
}
return *success*;

This algorithm attempts to cover $\mathbf{PER}(S_0)$ with a collection of sets $\mathbf{PER}(K^i)$, and associate a command c^i with each of these sets. If $\mathbf{PER}(K^i) \cap \mathbf{PER}(K^j) \neq \emptyset$ with $i \neq j$, then the states in this intersection have the following property: The command prescribed by the plan, if the system initially lies in one of these states, may vary in different plan executions.

The set $\mathbf{KER}(S^+, S) \equiv S^+ \setminus \mathbf{INT}(\mathbf{PER}(S \setminus S^+))$ is called the *kernel* of the set S^+, and it has a special significance in reasoning about perception under uncertainty. Suppose we know that the state of the system s lies in a set S, and we wish to derive a condition based on the perceived value of the system state $\hat{s}$ so that, if the condition is true, the state is guaranteed to lie in some prespecified set S^+. This condition is simply $\hat{s} \in \mathbf{PER}(\mathbf{KER}(S^+, S))$. The structure $\mathbf{INT}(\mathbf{PER}(S \setminus S^+))$ represents all the system states that may produce the same perceived value $\hat{s}$ with some system state not in S^+, i.e., all system states that can be *confused* with a state outside S^+. Subtracting these states from our target set S^+ leaves only those states in S^+ that cannot be confused with feasible states outside S^+.

2.4.3 An example

Let us examine how the above ideas apply to a particular example. Consider again the point robot that moves on a straight line. Let its initial position be known to be in $[-0.6, 0.4]$, and let the goal be [2.5,5]. This time, however, the command set that is available to the user is richer: $\{M_i, \mid i = 0, \pm 1, \pm 2, \ldots, \pm m\}$. The action function is defined as follows:

$$\mathbf{ACT}(M_i, [a, b]) = [a + 0.8 \times i,\ b + 1.2 \times i],$$

hence the preimage function is

$$\mathbf{PRE}(M_i, [a, b]) = \begin{cases} \emptyset, & \text{if } 0.4 \times i > (b - a); \\ [a - 0.8 \times i,\ b - 1.2 \times i], & \text{otherwise.} \end{cases}$$

Keeping the same perception and interpretation function as before, the possible initial perceived values are included in $[-0.7, 0.5]$, the perception set. The planning algorithm visits all commands in arbitrary order (say, in increasing order of i), and computes the preimage of the goal for this command and the kernel of the intersection of the preimage with the initial set. The first command with a non-empty kernel is M_3. The preimage of the goal for this command is $[0.1, 1.4]$, so S^+ is $[0.1, 0.4]$, S^- is $[-0.6, 0.1)$, and $\mathbf{INT}(\mathbf{PER}(S^-)) = [-0.8, 0.3)$. Now we can compute the kernel to be $[0.3, 0.4]$ and its perception set to be $[0.2, 0.5]$. Thus, whenever the robot perceives an initial value of $\hat{s}_0$ in $[0.2, 0.5]$, it should execute M_3. We shrink the initial set to $[-0.6, 0.3)$ and the perception set to $[-0.7, 0.2)$.

We move to the next command M_4. The backprojection of the goal for this command is $[-0.7, 0.2]$, S^+ is $[-0.6, 0.2]$ and S^- is $(0.2, 0.3)$. We compute the kernel to be $[-0.6, 0]$, and we associate M_4 with any value of $\hat{s}_0$ in $[-0.7, 0.1]$. The initial set is shrunk to $(0, 0.3)$ and the perception set to $(0.1, 0.2)$.

Unfortunately, no other command produces a backprojection with a non-empty intersection with the initial set, so the planner returns failure. Indeed, no guaranteed plan exists. Imagine that we get an initial reading $\hat{s}_0 \in (0.1, 0.2)$. According to the interpretation function, the actual position of the robot can be less than 0.1 or greater than 0.2, but we have no way to know it. If for this value of $\hat{s}_0$ we command M_3 and the actual position of the robot is less than 0.1, then the resulting actual position of the robot can be less than 2.5, i.e., outside the goal. If, on the other hand, we command M_4 and the actual position of the robot is greater than 0.2, the resulting actual position may be greater than 5, once again outside the goal. Hence, there is no safe command to associate with an initial measurement in the set $(0.1, 0.2)$, a fact that was discovered by our planner, too.

2.4.4 Conditional plans

Let us now consider the same example but with slightly smaller control uncertainty:

$$\mathbf{ACT}(M_i, [a, b]) = [a + 0.85 \times i,\ b + 1.15 \times i],$$

and

$$\mathbf{PRE}(M_i, [a, b]) = \begin{cases} \emptyset, & \text{if } 0.3 \times i > (b - a); \\ [a - 0.85 \times i,\ b - 1.15 \times i], & \text{otherwise.} \end{cases}$$

In this case, the kernel that corresponds to M_3 is found to be $[0.15, 0.4]$, and the kernel that corresponds to M_4 afterwards is $[-0.6, 0.15)$. Their perception sets cover all possible initial readings of $\hat{s}_0$, so we get the following guaranteed plan:

if $(\hat{s}_0 \in [0.25, 0.5])$ **execute** M_3;
else if $(\hat{s}_0 \in [-0.7, 0.05])$ **execute** M_4;
else if $(\hat{s}_0 \in (0.05, 0.25))$ **execute** M_3 **or** M_4;

The above is an example of a *conditional* plan, where the selected command depends on the initial perceived state, as opposed to a *linear* plan, where the choice of command is independent of the initial measurement. Conditional plans are much more powerful than linear plans, because they solve a richer class of planning problems. However, complete planners that return conditional plans are more complicated to build than their linear counterparts. Indeed, if we are seeking only linear plans, we must just find a command for which the preimage of the goal fully includes the initial set.

2.5 Multi-step plans

As we saw in the previous example, a guaranteed one-step plan may not exist. In this case the planner must attempt to find a multi-step plan. Such plans have the advantage of allowing perception actions between the execution of commands, thus diminishing the uncertainty about the state of the system. The planner must reason about what information might become available to the robot at execution time, and try to create a plan that allows the robot to gather as much of this information is necessary in order to reliably achieve the goal.

In the example of the previous section with the higher perceptual uncertainty, we were not able to produce a one-step guaranteed plan when $\hat{s}_0 \in (0.1, 0.2)$. However, the following two-step plan is guaranteed:

execute M_2
get measurement $\hat{s}$
if $(\hat{s} \in (1.5, 2))$ **execute** M_2
else execute M_1

To verify that this plan is guaranteed, notice that the possible initial positions of the robot lie in the set $(0, 0.3)$. The first M_2 command transforms this set into $(1.6, 2.5)$, and the measurement $\hat{s}$ can take any value in $(1.5, 2.6)$. According to the plan, if $\hat{s} \in (1.5, 2)$, another M_2 is executed. In this case the possible starting positions of the robot lie in the set $(1.6, 2.1)$, so its possible final positions lie in $(3.2, 4.5)$, i.e., in the goal. Otherwise, $\hat{s}$ is in $[2, 2.6)$, the robot starts from $[1.9, 2.5)$, executes M_1, and ends up in $[2.7, 3.7)$, once again in the goal.

The general problem of whether a guaranteed n-step plan exists can be formulated as follows:

$$\begin{array}{l}
\forall \hat{s}_0 \in \mathbf{PER}(S_0),\ \exists c_0 :\ S_1 = \mathbf{ACT}(c_0, \mathbf{INT}(\{\hat{s}_0\}) \cap S_0), \\
\forall \hat{s}_1 \in \mathbf{PER}(S_1),\ \exists c_1 :\ S_2 = \mathbf{ACT}(c_1, \mathbf{INT}(\{\hat{s}_1\}) \cap S_1), \\
\qquad\qquad\qquad\qquad\qquad\qquad \ldots \\
\forall \hat{s}_{n-1} \in \mathbf{PER}(S_{n-1}),\ \exists c_{n-1} :\ \mathbf{ACT}(c_{n-1}, \mathbf{INT}(\{\hat{s}_{n-1}\}) \cap S_{n-1}) \subseteq \mathbf{G}
\end{array}$$

Finding a multi-step plan is a much harder problem than its one-step counterpart. The most common technique for finding multi-step plans under uncertainty is *preimage backchaining*. If the preimages of the goal do not cover the initial set in an appropriate way, they are considered as intermediate goals, and the same algorithm is called upon them, and so forth, until the entire initial set has been covered. In the following chapters we will study several versions of this powerful technique.

2.6 Commands with termination condition

In the previous section we have assumed that the robot reassesses its behavior as fast as it can perceive its environment and interpret the sensory readings. This assumption places a tremendous burden on planners that attempt to fully exploit the capability of robots for such adaptive behavior. This often yields very long plans, and thus planning efficiency is seriously affected.

In practice, the action and the perception-interpretation loops do not operate at the same frequency; the action loop is usually much slower. A command consists of a *control statement* **cs** and a *termination condition* **tc**, and prescribes the following behavior for the robot: "**while**($\neg$**tc**) **obey cs**." In general, the termination condition depends on the most recent estimate of the state of the system. The robot begins to alter the state according to **cs**; at the same time it keeps perceiving the current state, updating its estimate about the true state of the system, and evaluating the termination condition continuously. When **tc** evaluates to *true*, the robot stops obeying **cs** and seeks another command from the plan. Putting the above description into an algorithm, we have:

$\hat{s} \in \mathbf{PER}(S_0)$ is the perceived initial state of the system;
$\tilde{S} \leftarrow \mathbf{INT}(\{\hat{s}\}) \cap S_0$;
while $(\tilde{S} \not\subseteq \mathbf{G})$ **do** {
    $(\mathbf{cs}, \mathbf{tc}) \leftarrow \mathbf{dec}(\tilde{S}, P)$;
    adjust the robot's behavior to conform with **cs**;
    while $(\neg\mathbf{tc}(\tilde{S}))$ **do** {
        $S \leftarrow \mathbf{ACT}(\mathbf{cs}, \tilde{S})$;
        $\hat{s} \in \mathbf{PER}(S)$ is the perceived new state of the system;
        $\tilde{S} \leftarrow \mathbf{INT}(\{\hat{s}\}) \cap S$;
    }
}
return *success*;

However general the above formulation may seem, real systems behave in more convoluted ways. In fact, in real systems the outcome of the **ACT** function may depend on itself, yielding the recursive equation $S \leftarrow \mathbf{ACT}(\mathbf{cs}, S, \tilde{S})$. The simplest such situation occurs when we consider time t (a member of s in the above). Usually, the outcome of **ACT** depends on the elapsed time, i.e., on the value of one member of s. Many workarounds can be proposed (like splitting s into vectors with different constraints and/or properties), but for the purposes of this discussion we will ignore time (or more precisely its effects on **ACT**), and we will retain the notation $\mathbf{ACT}(\mathbf{cs}, \tilde{S})$ to represent all states achievable by the system if **cs** is executed continuously with no termination condition.

In the above algorithm the execution module keeps updating its estimate about the current state of the system with each new value of $\hat{s}$ it obtains. Let h, the *sensing history*, denote the series of all observed values of $\hat{s}$, starting with the one that led to the selection of the particular command $(\mathbf{cs}, \mathbf{tc})$, and up to the most recent one. We can always obtain the most current estimate about the state of the system using only S_0 and h. Therefore, the termination condition depends only on h and S_0. This most general form of the termination condition (that makes use of all available information) is called a termination condition with *history* and *state*.

A plan which can be described with the above formulation, makes use of all available information. This fact contributes to the complexity of the planning problem significantly. In what follows, we will study a restricted class of termination conditions, namely, those that do not depend on the sensing history. Therefore, we will assume that the robot does not update its estimate about the state with each sensing action. The execution algorithm of such a system looks like follows:

$\hat{s} \in \mathbf{PER}(S_0)$ is the perceived initial state of the system;
$\tilde{S} \leftarrow \mathbf{INT}(\{\hat{s}\}) \cap S_0$;
while $(\tilde{S} \not\subseteq \mathbf{G})$ **do** {
  $(\mathbf{cs}, \mathbf{tc}) \leftarrow \mathbf{dec}(\tilde{S}, P)$;
  adjust the robot's behavior to conform with **cs**;
  $S \leftarrow \mathbf{ACT}(\mathbf{cs}, \tilde{S})$;
  while $(\neg\mathbf{tc}(\tilde{S}))$ **do** {
    $\hat{s} \in \mathbf{PER}(S)$ is the perceived new state of the system;
    $\tilde{S} \leftarrow \mathbf{INT}(\{\hat{s}\}) \cap S$;
  }
}
return *success*;

Since the set S remains constant throughout the evaluation of **tc**, we can essentially treat **tc** as a function of $\hat{S}$. Furthermore, **tc** is a boolean function, so we can represent it with two sets $\hat{S}_{\mathbf{tc}}$ and $\hat{S}_{\neg\mathbf{tc}}$, where $\hat{S}_{\mathbf{tc}} \cup \hat{S}_{\neg\mathbf{tc}} = \hat{\Omega}$. The first contains all values of $\hat{s}$ that cause **tc** to evaluate to *true*, whereas the second contains all values that cause **tc** to become *false*. Their union $\hat{\Omega}$ is the set of all possible measurements.

Consider again a one-step plan. Let us assume that for a particular execution of the plan the command $(\mathbf{cs}, \mathbf{tc})$ has been chosen. In order to guarantee that the execution will be successful, we have to make sure that at some point $\mathbf{tc}(\hat{S})$ will evaluate to *true*, and that the first time that this happens the system will be in a goal state. The function **ACT** is not adequate to facilitate reasoning like the above. Because **cs** is executed continuously, we need a function that can return possible *trajectories* of s instead of possible values of it. We call this function the *forward* *projection* (**FP**).

Definition 2.1 (Forward Projection) *The forward projection* $\mathbf{FP}(\mathbf{cs}, S)$ *of a set of system states* S *for control statement* **cs** *is the set of possible trajectories of the system state* s*, that occur when the system, initially in some state in* S*, changes in response to the continuous execution of* **cs**.

Note that the set $\mathbf{ACT}(\mathbf{cs}, S)$ consists of all the points of the trajectories included in $\mathbf{FP}(\mathbf{cs}, S)$. Similarly, instead of the function **PRE** we use the *backprojection* function **BP**, defined as follows:

Definition 2.2 (Backprojection) *The backprojection* $\mathbf{BP}(\mathbf{cs}, S)$ *of a set of system states* S *consists of all system states, such that if a system initially being in one of these states starts changing by obeying* **cs**, *all possible trajectories of its state will pass through* S.

We can now use these definitions to derive a one-step planning algorithm for such a system. For each possible observable initial state of the system we need to compute a pair $(\mathbf{cs}, \mathbf{tc})$, such that all possible trajectories of the system state that are compatible with **cs** pass through the goal, i.e., we have to make sure that the goal is *reachable*. Reachability, though, is not a sufficient condition for the correctness of a plan. We also need to make sure that the system will stop following **cs** while it is still in the goal, i.e., that along each possible trajectory there exists a point for which **tc** will evaluate to *true*, and that the first time that this happens, this point will lie in the goal. This is the problem of goal *recognizability*.

The set $\mathbf{INT}(\hat{S}_{\mathbf{tc}})$ contains all the states that may cause **tc** to evaluate to *true*. Also, the set $\mathbf{K}_{\mathbf{tc}} = \mathbf{INT}(\hat{S}_{\mathbf{tc}}) \setminus \mathbf{INT}(\hat{S}_{\neg\mathbf{tc}})$ contains all the states that always cause **tc** to evaluate to *true*. As it can be verified, $\mathbf{K}_{\mathbf{tc}} = \mathbf{KER}(\mathbf{INT}(\hat{S}_{\mathbf{tc}}), \Omega)$, where Ω is the set of all possible states of s, hence $\mathbf{K}_{\mathbf{tc}}$ is called the *kernel* of **tc**.

Definition 2.3 (Kernel of a termination condition) *The kernel of a termination condition consists of all system states which* always *give rise to perceived states that cause the termination condition to evaluate to true.*

Clearly, all possible trajectories of a one-step guaranteed plan must lead the system into the intersection of the goal with the kernel of the termination condition, without first passing through a point that belongs to $\mathbf{INT}(\hat{S}_{\mathbf{tc}})$ but not to the goal. Thus, a planner should go over all possible combinations of **cs** and **tc**, and for each such pair compute the backprojection of $\mathbf{K_{tc}} \cap \mathbf{G}$ treating the set $\mathbf{INT}(\hat{S}_{\mathbf{tc}}) \setminus \mathbf{G}$ as an artificial obstacle to be avoided. If several of these backprojections can cover the initial set S_0 in the way described in Section 2.4, then we got a one-step guaranteed plan.

2.7 Selecting a termination condition

An efficient planner cannot afford to examine all possible termination conditions in order to determine the best one. Instead, planners must use heuristic algorithms to select a "nice" termination condition that depends only on the selected control statement.

The properties of a "nice" termination condition can be derived by our desire to obtain as big a backprojection as possible. The size of the intersection of its kernel with the goal is not always a reliable heuristic. What is of real importance is the size of the boundary of this intersection that is not blocked by obstacles (including the artificial obstacles induced by the termination condition). For example, if we use a **tc** that is always true, the intersection of its kernel with the goal is the entire goal, but it is totally surrounded by the artificial obstacle, which includes all non-goal system states. Thus, there exists no control statement that can lead the system into the goal from any state outside the goal. The backprojection for this termination condition is empty, although the size of the intersection of its kernel with the goal is the maximum possible. In other words, the power of the termination condition depends on its ability to discriminate between points in the goal and points outside it. A typical heuristic creates termination conditions which have the maximal kernel without inducing any artificial obstacles.

Let us now, for one final time, visit our example, to demonstrate the above ideas. We use again the command set (F, B), only this time we assume that these commands are executed continuously (i.e., they are control statements) until some termination condition becomes true. We assume that we know nothing about the speed of motion of the robot, thus the function **ACT** is defined as follows:

$$\mathbf{ACT}(c,[a,b]) = \begin{cases} [a,+\infty), & \text{if } c = F; \\ (-\infty,b], & \text{if } c = B. \end{cases}$$

Since the robot moves on a line, it will eventually get into any non-empty set that lies in the direction of its motion, so the backprojection function has the following description:

$$\mathbf{BP}(c,[a,b]) = \begin{cases} (-\infty,b], & \text{if } c = F; \\ [a,+\infty), & \text{if } c = B. \end{cases}$$

Consider the same one-step planning problem, with $S_0 = [-0.6, 0.4]$ and $\mathbf{G} = [2.5, 5]$. How can we select a "nice" termination condition? In this case where the robot moves on a line, the only consideration is not to create an artificial obstacle between the initial set and the goal. Thus "$\hat{s} \geq 2.6$" is a good termination condition: It yields $\hat{S}_{\mathbf{tc}} = [2.6, +\infty)$, $\hat{S}_{\neg\mathbf{tc}} = (-\infty, 2.6)$, $\mathbf{INT}(\hat{S}_{\mathbf{tc}}) = [2.5, +\infty)$, $\mathbf{INT}(\hat{S}_{\neg\mathbf{tc}}) = (-\infty, 2.7)$, and $K_{tc} = [2.7, +\infty)$. The intersection of the kernel with the goal is $[2.7, 5]$, and the artificial obstacle $(5, +\infty)$ is safely behind the goal, allowing the preimage of $[2.7, 5]$ for control statement F to fully enclose the initial set. In this case the plan "**while** $(\hat{s} < 2.6)$ **obey** F" is guaranteed.

If, however, we use the termination condition "$\hat{s} \geq 2.5$," it will create the artificial obstacle $[2.4, 2.5) \cup (5, +\infty)$ which fully encloses the goal: The robot cannot enter the goal without getting in danger to stop before it reaches it.

Chapter 3

Robot Motion Planning with Uncertainty

The difference between a robot and a computer is the ability of the former to exert forces upon the physical world, thus altering its configuration. Therefore, any task that is performed by a robot (but cannot be performed by a computer) must involve physical motion of the robot or components of its environment. Motion planning is the foundation upon which most robotic reasoning systems must be built. Unfortunately, motion planning is a very hard problem, and it is not surprising that a lot of effort has been directed by researchers towards this area. A comprehensive analysis of the various motion planning techniques can be found in [53].

Of all the parameters that describe the system of the robot and its environment, we separate those that can be changed, and we call them the *configuration* q. The number of dimensions of the configuration is referred to as the dimensionality of a motion planning problem. The complexity of motion planning depends to a great extent on the dimensionality of the problem.

Since no more than one physical object can occupy the same physical space, not all values of q are possible. We describe this fact with a set of constraints $\{\mathcal{B}_1(q), \ldots, \mathcal{B}_{n_b}(q)\}$, each corresponding to the existence of an *obstacle* in the workspace. If a particular value of q results in a collision (i.e., violates one of the constraints) it is called *infeasible*; otherwise it is called *feasible*. In Chapter 2 we implicitly assume that the existence of obstacles is incorporated into the function **act**, which never returns an infeasible system configuration. An alternative approach is to define **act** in the absence of obstacles, and reason about them explicitly.

In this chapter we make the ideas presented in the previous chapter concrete, by presenting a practical algorithm for planning the motion of a robot under control and sensing uncertainty. The algorithm uses the method of "preimage backchaining." We consider three sensing schemes (sticking, position sensing, force sensing) and their combination. Several planning problems, solved with the implemented algorithm, are shown. Finally, we discuss the notion of maximal preimages and show that they do not exist.

3.1 Background

3.1.1 The path-planning problem

Research in motion planning began in the mid-seventies. Some early results can be found in [92, 66, 65]. It was soon understood that the pure topological and geometrical aspects of motion planning constituted an interesting problem of its own, and a lot of effort was directed towards its solution. This problem, the *path-planning problem*, is the search for a path of feasible configurations (a collision-free path) starting at the initial state and ending at a goal state. The robot is assumed to be able to follow the path without error. The archetype of such problems is the so called *piano-movers problem*. When the dimensionality is not fixed, the path-planning problem has been shown to be PSPACE-hard [85]. Other theoretical results regarding lower and upper bounds of multiple variants of the problem can be found in [88]. When the dimensionality is fixed, algorithms have been proposed, which are polynomial in the number of algebraic surfaces bounding the workspace and the obstacles and their maximal degree [87, 13].

These results are of theoretical interest, but have little practical use in systems with high dimensionality, even when uncertainty is not an issue. Several other planning techniques have been proposed in an attempt to find more practical algorithms [11, 45, 57, 68]. Most of them trade *completeness* for *efficiency*, i.e., they search a smaller space with the hope that it contains at least one solution. The *artificial potential field* method, introduced by Khatib [50], has been widely adopted and has resulted in many practical motion planning algorithms, like the one described in [38]. The idea is that obstacles exert repulsive forces to the robot, whereas the goal attracts it. These forces generate a potential field in the workspace having a global minimum at the goal. The robot moves following the gradient of the potential field until it reaches a minimum; if the minimum is local, the robot tries to escape, otherwise the goal has been achieved. Detecting and escaping local minima is the major problem of potential-field-based methods. Recently, the introduction of randomized and parallel algorithms has resulted in very effective techniques to address these problems. Barraquand and Latombe [5] described such a planner that is able to solve complicated motion planning problems with more than 10 degrees of freedom in a few seconds.

3.1.2 Dealing with uncertainty

Few of the algorithms described above can be used independently in real systems. The reason is that they do not take into account uncertainty. Reasoning about uncertainty complicates the motion-planning problem significantly. Moreover, since we cannot anticipate anything that may happen, we have to define precisely the extent to which uncertainty will be dealt with during the planning phase. This issue has been the subject of debate among roboticists and AI researchers, and has led to the formation of two major schools of thought.

The *reactionists* believe that planning should not be concerned with uncertainty. If unexpected situations arise during the execution of a plan, the robot should attempt to react to them following some prespecified *behavior*. If the execution of the plan cannot continue, the planner is asked to provide a new plan. The arguments of this school are (a) that it is not worth to spent too much planning time preplanning about low probability events, and (b) that even if we can anticipate some uncertainty, there are so many unexpected events, that the plan will fail in one way or another; if the robot must deal with such failures, it will also be able to deal with failures that could have been anticipated. Several systems using the above "reactive planning" method have been proposed and implemented with various degrees of success [10, 39, 1, 73]. The major disadvantage of this method is the inability to model the behavior of the robot in a meaningful way. Success or failure cannot be explained, so results cannot be generalized or transferred to a different system. Furthermore, uncertainty exists both at planning time and at execution time. One example of execution time uncertainty is perceptual errors that may prevent pure reactive schemes from completely eliminating uncertainty by reading sensory inputs. Perceptual uncertainty may also hinder the robot to correctly identify its own state, and the state of the workspace. False state identification may trigger inappropriate robot behavior. It is necessary for the robot to reason in advance about the knowledge that may become available at execution time, and make sure that this knowledge will be sufficient to lead it towards the goal or recognize failure.

For these reasons, the other school proposes that a planner should anticipate "as much uncertainty as possible." Typically, the term "as much uncertainty as possible" means that the considered stochastic parameters have distributions with relatively small variance (i.e., we do have some knowledge about these parameters), and that an algorithm that solves the planning problem exists. Unfortunately, most planning-with-uncertainty algorithms that have been proposed up to now require exponential running time. Our view is that it is only worthwhile to reason about uncertainty if the planning algorithm is fast enough to be practical. Moreover, we believe that reactive behaviors (like unexpected obstacle avoidance) should be an integral part of a robust robotic system, because there is no way that a planner can anticipate *everything* that may happen *all* the time.

3.1.3 Planning with uncertainty

Four major approaches to planning with uncertainty have been proposed so far. The first, called *skeleton refinement* was introduced independently by Lozano-Pérez [66] and Taylor [92]. It consists of (a) retrieving a plan skeleton appropriate to the task at hand and taking that as an initial plan, and (b) iteratively modifying the skeleton by inserting complements (typically sensor readings). Complements are decided after the correctness of the skeleton is checked, either by propagating uncertainty through the steps of the plan skeleton [92] or by simulating several possible executions [66]. Subsequent contributions to the approach have been brought by Brooks [9], who developed a symbolic computation technique for propagating uncertainty forward and backward through plan skeletons, and by Pertin-Troccaz and Puget [81], who proposed techniques for verifying the correctness of a plan and amending incorrect plans.

The second approach to motion planning with uncertainty was proposed by Dufay and Latombe [29], and is known as the *inductive learning* approach. It consists of assembling input partial strategies into a global one. First, during a training phase, the system uses the partial strategies to make on-line decisions and execute several instances of the task at hand. Then, during an induction phase, the system combines the execution traces generated during the training phase, and generalizes them into a global strategy. In fact, the training phase and the induction phase are interweaved. The generation of a strategy for the task ends when new executions do not modify the current strategy. A system based on these principles has been implemented, and experimented successfully on several part mating tasks.

A third method, called *iterative removal of contacts* creates plans using compliant motions on obstacle faces. Compliant motions have the advantage that they completely eliminate uncertainty along the normal to the contact surface. Several variants of this method have been proposed [58, 44, 20].

The fourth approach is called *preimage backchaining* and it will be presented in detail in the following.

3.1.4 Preimage backchaining

The first three approaches essentially operate in two phases: first, a motion plan is generated assuming no uncertainty, and then it is transformed to deal with uncertainty. Instead, preimage backchaining takes uncertainty into account throughout the whole planning process. In principle, it can tackle problems where uncertainty shapes the structure of a plan to the extent that the plan cannot be generated by transforming an initial one produced under the no-uncertainty assumption. This method was first introduced by Lozano-Pérez, Mason and Taylor [69] and has since been accepted as the primary method for the creation of guaranteed plans despite the presence of uncertainty.

The *preimage* of a goal region, for a given motion command $\mathbf{M}$, is the set of all robot configurations such that if the robot starts executing the command from any one of these configurations, it is guaranteed to reach the goal and stop in it. Preimage backchaining consists of constructing a sequence of motion commands $\mathbf{M}_i$ ($i = 1, \ldots, n$), such that, if $\mathcal{P}_n$ is the preimage of the goal for $\mathbf{M}_n$, $\mathcal{P}_{n-1}$ the preimage of $\mathcal{P}_n$ for $\mathbf{M}_{n-1}$, and so on, then $\mathcal{P}_1$ contains all possible initial configurations of the robot.

One source of difficulty in computing preimages is the interaction between goal reachability and goal recognizability. The robot must both reach the goal (despite initial position and control uncertainty) and stop in the goal (despite sensing uncertainty). Goal recognition often depends on the region from where the command is executed. This region, which is precisely the preimage of the goal for that command, also depends on the way the goal is recognized. This recursive dependence was noted in [69]. Despite this difficulty, Canny [14] described a complete planner with very few restrictive assumptions. This planner generates a plan consisting of r motion commands by reducing the input problem to deciding the satisfiability of a semi-algebraic formula with $2r$ alternating existential and universal quantifiers. Such a decision takes double exponential time in r. Even worse, the smallest r for which a plan may exist grows with the complexity of the environment. Actually, various forms of the above motion planning problem have been proven intrinsically hard [15, 77, 14].

At the expense of completeness, Erdmann [32] suggested that goal reachability and recognizability be treated separately by identifying a subset of the goal, called a *kernel*, such that when this subset is attained, goal achievement can be recognized independently of the way it has been achieved. He defined the backprojection of a region $\mathcal{T}$, for a motion command $\mathbf{M}$, as the set of all points such that, if the robot executes $\mathbf{M}$ starting at any one of these points, it is guaranteed to reach $\mathcal{T}$. He proposed an $O(n \log n)$ algorithm to compute backprojections in the plane when the obstacles are polygons bounded by n edges. An implemented planner based on this approach is described in [55].

3.2 The configuration space

We have previously defined the *configuration* q of a robot and its environment as the values of the changeable parameters of the system. Let us now narrow this definition for pure robot motion planning problems, i.e., for problems where the robot is the only movable object in a workspace $\mathcal{W}$, populated by stationary obstacles $\mathcal{B}_i$, $i \in [1, n_b]$. A *configuration* of the robot is a specification of the position of every point of the robot with respect to a coordinate system embedded in $\mathcal{W}$ [2]. The *configuration space* denoted by $\mathcal{C}$, is the set of all possible configurations.

Each obstacle $\mathcal{B}_i$ is mapped to the subset $\mathcal{CB}_i$ of configurations for which the robot collides with $\mathcal{B}_i$:

$$\mathcal{CB}_i = \{q \in \mathcal{C} \mid \mathcal{R}(q) \cap \mathcal{B}_i \neq \emptyset\},$$

where $\mathcal{R}(q)$ denotes the subset of $\mathcal{W}$ occupied by the robot at configuration q. The region $\mathcal{CB}_i$ is called a *C-obstacle*.

In general, $\mathcal{C}$ is a non-Euclidean manifold. For instance, if the robot is a rigid planar object moving freely in $\mathcal{W} = \Re^2$, then $\mathcal{C} = \Re^2 \times S^1$, where S^1 denotes the unit circle. If the robot is a rigid three-dimensional object moving freely in $\mathcal{W} = \Re^3$, $\mathcal{C} = \Re^3 \times SO(3)$, where $SO(3)$ denotes the Special Orthogonal Group of orthonormal matrices with determinant $+1$ [2].

For the rest of the chapter, we assume that the robot is a two-dimensional polygonal object that can only translate in the plane, e.g., an omnidirectional mobile robot that cannot rotate. A configuration is represented by $q = (x, y) \in \Re^2$, where x and y are the coordinates of a specific point of the robot (the *reference point*) with respect to the coordinate system embedded in $\mathcal{W}$. Hence, both $\mathcal{W}$ and $\mathcal{C}$ are copies of $\Re^2$. Both the robot and the obstacles are modeled as polygonal regions, the robot as a simple polygon, and each obstacle as a region whose boundary is a simple polygonal curve

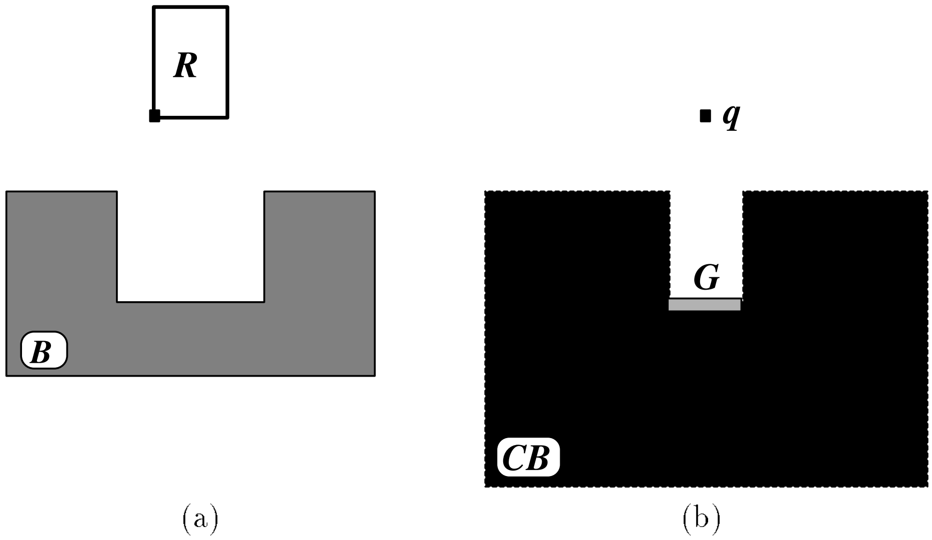

Figure 3.1: Peg-Into-Hole Task

that may, or may not be a closed-loop curve (allowing for non-bounded obstacles). Under these assumptions, each C-obstacle $\mathcal{CB}_i$ is also a polygonal region.

Figure 3.1 illustrates the above concepts. A simple setting in the workspace is depicted in Figure 3.1(a). The robot $\mathcal{R}$ is a rectangle and there is a single polygonal obstacle $\mathcal{B}$ with a rectangular depression. The goal of the task is to insert $\mathcal{R}$ in $\mathcal{B}$'s depression ("peg-into-hole" task). Figure 3.1(b) shows the mapping of this setting in $\mathcal{R}$'s configuration space. The robot $\mathcal{R}$ maps to the point denoted by q. The obstacle $\mathcal{B}$ maps to the C-obstacle $\mathcal{CB}$. The width of the rectangular depression in $\mathcal{CB}$ is equal to the difference between the width of the depression in $\mathcal{B}$ and the width of $\mathcal{R}$. The goal of the peg-into-hole task is to move q (the configuration of $\mathcal{R}$) to any location in the edge $\mathcal{G}$ at the bottom of $\mathcal{CB}$'s depression.

The union of the obstacles, $\bigcup_{i=1}^{n_b} \mathcal{B}_i$, is called the *obstacle region* and is denoted by $\mathcal{B}$. The union of the C-obstacles, $\bigcup_{i=1}^{n_b} \mathcal{CB}_i$, is called the *C-obstacle region* and is denoted by $\mathcal{CB}$. The complement of the C-obstacle region in $\mathcal{C}$ is called the *free space* and is denoted by $\mathcal{C}_{free}$. The subset of configurations q where $\mathcal{R}(q)$ intersects with the obstacle region without overlapping its interior is called the *contact space* and is denoted by $\mathcal{C}_{contact}$. The union of the free space $\mathcal{C}_{free}$ and the contact space $\mathcal{C}_{contact}$ is called the *feasible space* and is denoted by $\mathcal{C}_{feasible}$. A *feasible path* between two configurations q_1 and q_2 in $\mathcal{C}_{feasible}$ is a continuous map $\tau : [0,1] \rightarrow \mathcal{C}_{feasible}$ such that $\tau(0) = q_1$ and $\tau(1) = q_2$.

The set $\mathcal{C}_{free}$ is an open subset of $\mathcal{C}$, whose boundary $\partial\mathcal{C}_{free}$ is a set of polygonal curves. $\mathcal{C}_{contact}$ is also made of polygonal curves. It can be shown that $\partial\mathcal{C}_{free} \subseteq \mathcal{C}_{contact}$ and $cl(\mathcal{C}_{free}) \subseteq \mathcal{C}_{feasible}$ [48]. In Figure 3.1, the strict inclusion of $\partial\mathcal{C}_{free}$ in $\mathcal{C}_{contact}$ occurs when $\mathcal{R}$'s width is exactly equal to the width of $\mathcal{B}$'s depression. Then the depression in $\mathcal{CB}$ degenerates to a line segment contained in $\mathcal{C}_{contact}$, but not in $\partial\mathcal{C}_{free}$. In order to avoid such pathological cases, we assume that each maximally connected subset of the C-obstacle region $\mathcal{CB}$ is homeomorphic[1] to a closed disc or a closed half-space, i.e., is a manifold with boundary. This assumption entails that $\mathcal{C}_{contact} = \partial\mathcal{C}_{free}$ and $\mathcal{C}_{feasible} = cl(\mathcal{C}_{free})$; hence, a contact between $\mathcal{R}$ and the obstacle region can be broken by an arbitrarily small displacement of $\mathcal{R}$. It also implies that every vertex in the contact space $\mathcal{C}_{contact}$ is the extremity of only two edges; thus, $\mathcal{C}_{contact}$ consists of disjoint simple polygonal lines. An edge (resp. a vertex) in $\mathcal{C}_{contact}$ is called a *contact edge* (resp. a *contact vertex*). The *outgoing normal* of a contact edge E is the unit vector $\vec{\nu}(E)$ normal to E pointing toward $\mathcal{C}_{free}$. The vector $-\vec{\nu}(E)$ is the *ingoing normal* of E.

Let $n_\mathcal{R}$ be the number of edges of $\mathcal{R}$, and $n_\mathcal{B}$ be the number of edges of the obstacle region. $\mathcal{C}_{contact}$ contains $O(n_\mathcal{R}^2 n_\mathcal{B}^2)$ edges which can be computed in $O(n_\mathcal{R}^2 n_\mathcal{B}^2 \log n_\mathcal{R} n_\mathcal{B})$ time [89].

3.3 Control, sensing and uncertainty

3.3.1 Structure of a motion plan

According to the model of Chapter 2, the execution of a motion plan occurs in *steps*. During each step the execution module selects a *control statement* **cs** which prescribes a particular behavior to be followed by the robot, and a *termination condition* **tc** which is a boolean function of the sensory readings. (Time is considered to be one of the possible sensory readings.) The pair $(\mathbf{cs}, \mathbf{tc})$ constitutes a *motion command.*

[1] Two topological sets E and F are *homeomorphic* iff there exists a bijection $f : E \rightarrow F$ such that both f and f^{-1} are continuous.

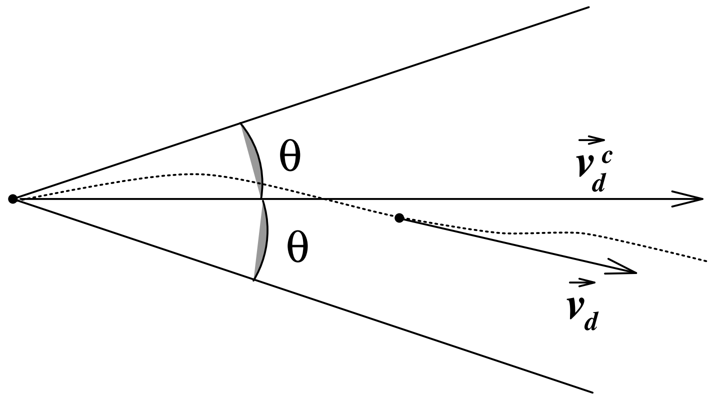

Figure 3.2: The control uncertainty cone

The robot moves obeying **cs**, until **tc** evaluates to true, at which point it stops instantaneously.

The way in which the execution module decides what motion command to execute next depends on the structure of the motion plan. *Linear plans* specify a sequence of motion commands that are to be executed one after the other. *Conditional plans* specify several possible motion commands at each execution step. The robot selects one of them to execute next, based on its beliefs about its own state and the state of the world. These beliefs depend mainly on sensory readings, but may also depend on sensing history, and other prior information. Finally, *randomized plans* use a stochastic function in order to select a motion command at each step. In this chapter we consider only linear plans.

3.3.2 Control

Several schemes have been proposed to control the motion of a robot. Here, we assume that the only control parameter is the *commanded velocity* of motion $\vec{\mathbf{v}}_d^c$. The robot is commanded to move along $\vec{\mathbf{v}}_d^c$, but it is not always possible to do so because of control uncertainty. At any instant during the motion in free space, the *actual velocity* of the robot is a vector $\vec{\mathbf{v}}_d$, such that the magnitude of the angle between $\vec{\mathbf{v}}_d^c$ and $\vec{\mathbf{v}}_d$ is less than a fixed angle $\theta < \pi/2$. In other words, $\vec{\mathbf{v}}_d$ lies in a cone of angle 2θ whose axis points along the d direction. This cone is called the *control uncertainty*

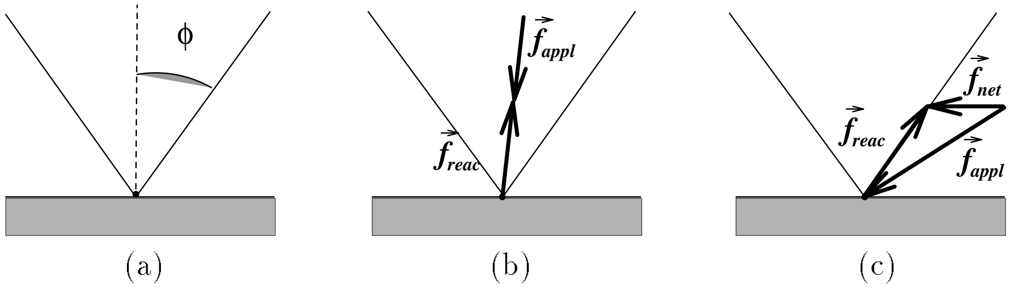

Figure 3.3: Reaction Force

cone. The modulus of the velocity of the robot is unknown to the planner, but is assumed to have some reasonable (positive) value. For simplification, we assume that there is no uncertainty in the velocity modulus, i.e., $\|\vec{\mathbf{v}}_d\| = \|\vec{\mathbf{v}}_d^c\|$. This assumption has no significant impact on the methods described below and can be easily relaxed. For further simplification (and without loss of generality), we assume that $\vec{\mathbf{v}}_d^c$ (hence, $\vec{\mathbf{v}}_d$) is a unit vector, so that the direction d suffices to characterize the vector $\vec{\mathbf{v}}_d^c$ (and the range of vectors $\vec{\mathbf{v}}_d$ compatible with $\vec{\mathbf{v}}_d^c$). During motion in free space, $\vec{\mathbf{v}}_d$ may vary arbitrarily between the two extreme orientations determined by $\vec{\mathbf{v}}_d^c$ and θ. Thus, if the robot is in the free space, it moves along a trajectory whose tangent at any configuration is contained in the control uncertainty cone anchored at this configuration (see Figure 3.2).

The behavior of the robot when it comes in contact with obstacles depends on the particular control scheme we use. Sometimes, contact with obstacles is not allowed. In this case, $\mathcal{C}_{feasible} = \mathcal{C}_{free}$. In most cases, though, compliant motions are allowed, because they eliminate the positional uncertainty along the normal to the contact surface and allow the robot to cover long distances with no uncertainty penalty along one axis.

The control scheme described above can be extended to deal with compliant motion [71, 83, 96]. It is called the *generalized damper compliance model*.

According to this model, the robot exerts a force on its environment that is proportional to the actual commanded velocity $\vec{\mathbf{v}}_d$, say $\vec{\mathbf{f}}_{appl} = B\vec{\mathbf{v}}_d$. When the robot is in free space, no reaction force is applied to the robot, and the force exerted by the robot is entirely used to create motion. When the robot is in contact space and pushes on an obstacle, the obstacle generates a reaction force. If the vector $-\vec{\mathbf{v}}_d$ lies in the friction cone at the contact configuration, the reaction force is $-B\vec{\mathbf{v}}_d$, i.e., it completely cancels the force exerted by the robot, and there is no motion (Figure 3.3(b)). If the vector $-\vec{\mathbf{v}}_d$ lies outside the friction cone, the reaction force only partially cancels the force exerted by the robot and the net resulting force makes the robot's configuration slide in contact space (Figure 3.3(c)).

Let q be the actual configuration of the robot at some instant t. We denote by $\vec{\mathbf{f}}_{reac}(q, \vec{\mathbf{v}}_d)$ the vector in configuration space representing the reaction force applied at the robot's configuration when the actual commanded velocity is $\vec{\mathbf{v}}_d$. $\vec{\mathbf{f}}_{reac}(q, \vec{\mathbf{v}}_d)$ is exhaustively defined as follows:

- if $q \in \mathcal{C}_{free}$, then $\vec{\mathbf{f}}_{reac}(q, \vec{\mathbf{v}}_d) = 0$;
- if $q \in \mathcal{C}_{contact}$, q is in the interior of a contact edge E, and $\vec{\mathbf{v}}_d$ projects positively on the outgoing normal of E, then $\vec{\mathbf{f}}_{reac}(q, \vec{\mathbf{v}}_d) = 0$;
- if $q \in \mathcal{C}_{contact}$, q is in the interior of a contact edge E, and $-\vec{\mathbf{v}}_d$ lies in the friction cone at the contact point, then $\vec{\mathbf{f}}_{reac}(q, \vec{\mathbf{v}}_d) = -B\vec{\mathbf{v}}_d$ (see Figure 3.3(b));
- if $q \in \mathcal{C}_{contact}$, q is in the interior of a contact edge E, and $-\vec{\mathbf{v}}_d$ projects positively on the outgoing normal of E but lies outside the friction cone at the contact configuration, then $\vec{\mathbf{f}}_{reac}(q, \vec{\mathbf{v}}_d)$ is equal to the projection of $-B\vec{\mathbf{v}}_d$ onto the nearest side of the friction cone at q (the projection is taken normal to the axis of the cone, see Figure 3.3(c));
- if $q \in \mathcal{C}_{contact}$, q is at a contact vertex, then the reaction force is a positive linear combination of the reaction forces that can be exerted by the edges abutting the vertex.

According to the above definition, if $\vec{\mathbf{v}}_d^c$ and θ are such that all directions in the control uncertainty cone project positively on the outgoing normal of a surface, the robot is guaranteed to move away from the surface. If the control uncertainty cone is completely included in the inverted friction cone, then the robot is guaranteed to stick on the surface. If all directions in the control uncertainty cone project positively on the ingoing normal of the surface but none of them lies in the inverted friction cone, then the robot is guaranteed to slide along the surface. In any other case the behavior of the robot cannot be predicted at planning time.

3.3.3 Sensing

Real robots are equipped with a variety of sensing equipment that provide measurements of the actual configuration of the robot and the workspace. These measurements contain stochastic errors, whose distribution is called *sensing uncertainty*. Clearly, the more information the robot has, the more complex tasks it is able to perform. On the other hand, though, reasoning about sensing information at planning time is a significant source of complexity for planners since the actual sensed values are not known at planning time. It is therefore important to distinguish the information that is more likely to have a significant impact on the structure of a plan, and ignore other potential sources of less important information. As a result, researchers have been particularly frugal when selecting the sensing inputs that a planner-with-uncertainty considers.

Sticking This is the simplest form of sensing, and for this reason robots that use it are often called *sensorless*. The only thing that the robot can sense is its own motion as a binary variable (1 if it is in motion, 0 if it is stopped). The robot moves following the generalized damper model, using a termination condition that stops its motion automatically (sticking on obstacle edges due to friction). Nevertheless, the robot must realize that motion has stopped, in order to proceed with the execution of the next command of the plan. It is assumed that the robot never makes a mistake in deciding whether it moves or not, so no sensing uncertainty exists in this case.

Force sensing A force sensor measures the reaction force exerted on the robot. The output of the sensor is a vector denoted by $\vec{\hat{\mathbf{f}}}$. For simplification, we assume that the modulus of the measured force is accurate. On the other hand, the magnitude of the angle between $\vec{\mathbf{f}}$ and $\vec{\hat{\mathbf{f}}}$ (if they are non-zero vectors) is less than a fixed angle $\varepsilon < \pi/2$. Hence, $\vec{\mathbf{f}}$ lies in an open cone of angle 2ε whose axis points along $\vec{\hat{\mathbf{f}}}$.

Position sensing A position sensor measures the current configuration of the robot. The sensed configuration is denoted by $\hat{q}$, while the current actual configuration is denoted by q. The uncertainty in the measurement is modeled as an open disc $\mathcal{U}_q(\hat{q}) \subset \Re^2$ of fixed radius ρ centered at the sensed configuration $\hat{q}$. This means that if the position sensor returns $\hat{q}$, the current actual configuration q may be anywhere in $\mathcal{U}_q(\hat{q}) \cap \mathcal{C}_{feasible}$. Reciprocally, if the actual configuration is known to be q, the sensed configuration $\hat{q}$ should necessary lie in $\mathcal{U}_q(q)$.[2]

3.4 Preimage backchaining in two dimensions

3.4.1 Statement of the problem

Let $\mathcal{I}$ be a subset of $\mathcal{C}_{feasible}$, in which it is known at planning time that the robot's configuration will be when the execution of the motion plan starts. $\mathcal{I}$ is called the *initial region*. Let $\mathcal{G}$ be another subset of $\mathcal{C}_{feasible}$ called the *goal region*. Given a robot that can be modelled as described in the previous sections, we want the planner to generate a sequence of motion commands (a linear motion plan), whose execution makes the robot move from its actual initial configuration in $\mathcal{I}$ to a final configuration in $\mathcal{G}$.

3.4.2 Preimage backchaining

The plan is generated by working backwards from the goal (backchaining). During each step, we identify a target set $\mathcal{T}$ (initially the goal), we select a motion command $(\vec{\mathbf{v}}_d^c, \mathbf{tc})$, and we compute the set of configurations $\mathcal{P}(d, \mathbf{tc}; \mathcal{T})$ from where the motion command is guaranteed to make the robot reach the target set (reachability) and stop in it (recognizability). $\mathcal{P}(d, \mathbf{tc}; \mathcal{T})$ is the preimage of $\mathcal{T}$ as defined in Chapter 2. If the initial region is covered entirely by the preimage, then the motion command $(\vec{\mathbf{v}}_d^c, \mathbf{tc})$ is guaranteed to bring the robot into $\mathcal{T}$ and make it stop in there. If this is

[2] This is not always true. It happens to be so in this case because the perception function $\mathcal{U}_q(\hat{q})$ is symmetric with respect to $\hat{q}$ (i.e., $x \in \mathcal{U}_q(\hat{q}) \equiv 2\hat{q} - x \in \mathcal{U}_q(\hat{q})$). If this weren't so, the interpretation function $\mathcal{U}_q^{-1}(q)$ would be the symmetric set of $\mathcal{U}_q(q)$ with respect to q.

the case, the algorithm terminates; otherwise, $\mathcal{P}$ becomes the target set of the next backchaining step. The number of motion commands in a plan is equal to the number of backchaining steps.

The computation of preimages combines the notions of reachability and recognizability. Reachability relates to the fact that any trajectory obtained by executing a motion commanded along $\vec{\mathbf{v}}_d^c$ from a preimage of $\mathcal{T}$ should reach $\mathcal{T}$. Given a starting configuration, $\vec{\mathbf{v}}_d^c$ only determines a *commanded trajectory* τ_d^c. Due to errors in control, any execution produces an *actual trajectory* τ_d that is slightly different. The planner must be certain that all the possible actual trajectories consistent with both $\vec{\mathbf{v}}_d^c$ and control uncertainty (i.e., θ) will reach $\mathcal{T}$ at some instant.

Reaching $\mathcal{T}$, however, is not sufficient since the termination condition could just let the robot traverse the target set without stopping in it, or, instead, it could halt the motion prematurely before $\mathcal{T}$ has been reached. In fact, the planner must be certain that the termination condition **tc** will stop the robot in $\mathcal{T}$ (recognizability). The function **tc** is an observer of the actual trajectory τ_d being executed. Since sensing is imperfect, **tc** perceives τ_d as an *observed trajectory* $\hat{\tau}_d$ which is most likely to be neither the commanded one, nor the actual one. Thus, the problem for the planner is to be sure that, for every possible observed trajectory $\hat{\tau}_d$, (a) **tc** becomes `true` at some instant, and (b) at the instant t_f when **tc** first becomes `true`, all the actual trajectories τ_d consistent with $\hat{\tau}_d$ (i.e., the trajectories which may be observed as $\hat{\tau}_d$ given the uncertainty in sensing) have reached the target set, namely, $\tau_d(t_f) \in \mathcal{T}$.

Preimage computation has been investigated in depth in [32, 52]. For a given commanded velocity $\vec{\mathbf{v}}_d^c$, the ideal method would compute the maximal preimage of $\mathcal{T}$ over all possible termination conditions, i.e., a preimage that is not contained in any other preimage[3] of $\mathcal{T}$ for $\vec{\mathbf{v}}_d^c$. The method would also return the termination condition for the maximal preimage. Intuitively, a large preimage has more chance to include the initial region $\mathcal{I}$ than a small one; in addition, if it is considered recursively as an intermediate goal, a large preimage has more chance to admit larger preimages than a small one. Thus, considering larger preimages may reduce the size of the search

[3]By definition, any subset of a preimage is also a preimage of the same set for the same motion command.

graph; it may also produce "simpler" strategies (strategies made up of less motion commands). Nevertheless, Erdmann [32] showed that: (a) a maximal preimage does not always exist; (b) if one does exist, it may not be unique; and (c) if a unique maximal preimage exists, it may depend in a very subtle fashion on sensing history, the elapsed time since the beginning of the motion, and the knowledge embedded in the termination condition.

Most of the difficulties related to maximal preimages are due to the subtle interdependence between reachability and recognizability. Indeed, even after a particular termination condition has been selected, the region of the workspace where this condition always evaluates to true (the kernel of the termination condition as defined in Chapter 2), which is needed in order to compute the preimage, may depend on the region from where this command is executed, which is precisely the preimage of this region.

3.4.3 Backprojection from target kernel

One way of breaking this recursive dependence is to use restricted forms of the termination condition so that these two issues can be treated separately. The basic idea is to identify a subset $\mathcal{K}$ of the target set $\mathcal{T}$, in which achievement of $\mathcal{T}$ can be recognized independently of the starting region. A termination condition is constructed, so that it always evaluates to `true` when the robot is in $\mathcal{K}$, i.e., $\mathcal{K}$ is the kernel of this termination condition according to definition 2.3. In this case, though, we first compute $\mathcal{K} = \chi(d; \mathcal{T})$ as a function of the target set and possibly the commanded direction of motion, and we construct the termination condition that corresponds to $\mathcal{K}$ afterwards. It now remains to compute the region $\mathcal{BP}(d; \mathcal{K})$, from where the robot is guaranteed to attain the kernel if the commanded velocity is $\vec{\mathbf{v}}_d^c$. In Chapter 2 we call this region the backprojection of $\mathcal{K}$.

For every possible termination condition, the backprojection of a set is bigger than (or equal to) any preimage of the set for the same commanded direction of motion:

$$\forall d\, \forall \mathbf{tc}\ \ \mathcal{P}(d, \mathbf{tc}; \mathcal{T}) \subseteq \mathcal{BP}(d; \mathcal{T}).$$

For example, look at Figure 3.4. Assume that the robot can only sense sticking on

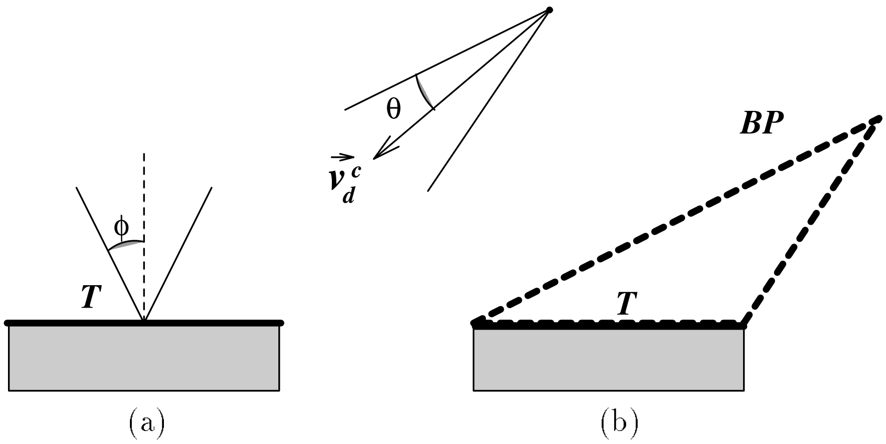

Figure 3.4: Preimage and backprojection

obstacle edges. Its target set $\mathcal{T}$ is the top edge of a rectangular box. For choices of $\vec{\mathbf{v}}_d^c$ such that the control uncertainty cone is not contained in the inverted friction cone, the preimage of $\mathcal{T}$ is the empty set since the robot cannot stick on the target edge. On the other hand, the backprojection of $\mathcal{T}$ is outlined with the thick dashed line in Figure 3.4(b).

Conversely, there exists a termination condition $\mathbf{tc}$, for which the preimage of a set is bigger than (or equal to) the backprojection of the kernel of the set for the same commanded direction of motion:

$$\forall d\, \exists \mathbf{tc}\;\; \mathcal{P}(d, \mathbf{tc}; \mathcal{T}) \supseteq \mathcal{BP}(d; \chi(d; \mathcal{T})).$$

This follows directly from the definition of the kernel. Whenever the termination condition $\hat{q} \in \mathcal{U}_q(\chi(d, \mathcal{T}))$ is true, we know that $q \in \mathcal{T}$. The preimage of $\mathcal{T}$ for this termination condition is equal to the backprojection of the kernel. Other termination conditions may produce bigger preimages.

3.5 Kernel computation

The computation of the kernel of a target set depends on the particular sensing readings that are being provided to the termination condition. In this section we describe kernel computation for all the sensing schemes presented in Section 3.3 (sticking, force sensing, position sensing), as well as for the combination of all three together.

3.5.1 Sticking

When sticking on obstacles is the only thing that can be sensed, there is no sensing uncertainty: once the robot sticks, motion terminates. Whether the robot will stick or not on an obstacle edge depends on its actual direction of motion at execution time. Since we are interested in producing guaranteed plans, the planner must identify those edges and vertices where the robot is guaranteed to stick (the *sticking edges/vertices*). Therefore, the planner must make sure that all actual velocity vectors that are compatible with the commanded velocity $\vec{\mathbf{v}}_d^c$ make the robot stick on such an edge or vertex.

The robot is guaranteed to stick on an edge, if the inverted control uncertainty cone lies entirely inside the friction cone of the edge. However, when the robot hits a vertex the situation is more complicated. We previously assumed that a vertex can exert a reaction force that is a positive linear combination of the forces exerted by the two edges forming the vertex. Consequently, the friction cone of a vertex consists of the friction cones of the two edges and the region between them. If the inverted velocity vector of the robot is included in the vertex friction cone, then the robot sticks at the vertex. However, small perturbations in the position of the robot may cause it to move on one of the two edges that form the vertex, and then move according to the reaction force that is exerted by this edge only. If this force pushes the robot back towards the vertex on both edges or makes it stick on the edge, then sticking at the vertex is stable; otherwise, the robot will move away from the vertex. We are only interested in stable sticking. The *sticking cone* of a vertex consists of the inverted velocity directions of the robot that cause stable sticking at the vertex.

Consider the friction cones (anchored at the vertex) of both edges. The region between each edge and its friction cone contains all inverted velocity directions that push the robot towards the vertex. The union of this region with the friction cone creates a bigger cone that contains all inverted commanded velocity directions that cause the robot either to stick at the vertex or to slide on the edge towards the vertex. The sticking cone of the vertex is the intersection of these cones of the two edges. In Figure 3.5 we show the sticking cones of three vertices (the region outlined with thick black rays): (a) For a concave vertex, the sticking cone is always equal to the friction

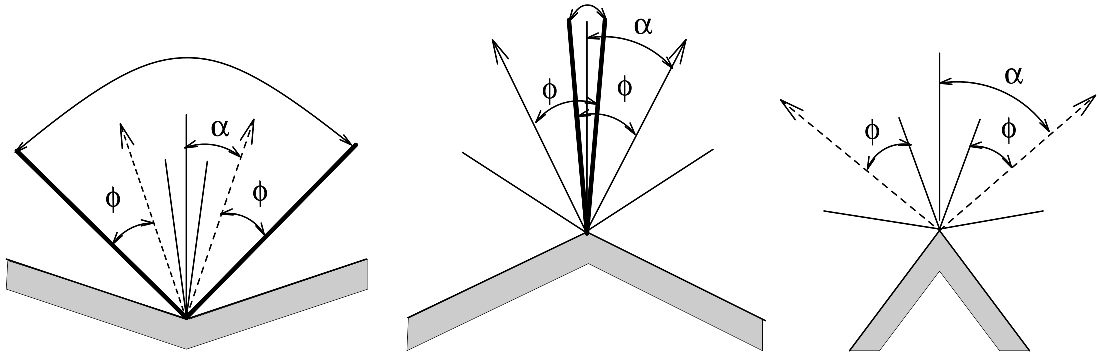

Figure 3.5: The sticking cone of a vertex

cone; (b) for a convex vertex, the sticking cone is the intersection of the friction cones of the two edges; but (c) when this intersection is empty, the sticking cone is null.

Given a commanded velocity $\vec{\mathbf{v}}_d^c$ and a target set $\mathcal{T}$, the computation of the kernel is done by (a) identifying all edges and vertices in contact space that belong to $\mathcal{T}$, (b) selecting those edges whose friction cone includes the inverted control uncertainty cone for $\vec{\mathbf{v}}_d^c$, and (c) selecting those vertices whose sticking cone includes the inverted control uncertainty cone for $\vec{\mathbf{v}}_d^c$. More specifically, the kernel consists of every edge E in $\mathcal{T} \cap \mathcal{C}_{contact}$ such that $|angle(-\vec{\mathbf{v}}_d^c, \vec{\nu}(E))| \leq \phi - \theta$, and every vertex V in $\mathcal{T} \cap \mathcal{C}_{contact}$ such that, if V is concave, $|angle(-\vec{\mathbf{v}}_d^c, \vec{\nu}(V))| \leq \phi + \alpha - \theta$, and if V is convex, $|angle(-\vec{\mathbf{v}}_d^c, \vec{\nu}(V))| \leq \phi - \alpha - \theta$. Here $\vec{\nu}(V)$ is the bisector of the angle of the normals of the two edges, and α is half the angle between these normals.

Notice that guaranteed sticking on edges is possible only if $\phi > \theta$, i.e., the friction cone is larger than the control uncertainty cone. This is not required for sticking at concave vertices. In fact, sticking at a concave vertex is easier and more stable than sticking along a contact edge, since the vertex may be reached by sliding along the two abutting edges.

3.5.2 Force sensing

The force sensor measures the reaction force exerted by an obstacle edge on the robot. Let $\mathbf{F}_{reac}(q, d)$ denote the set of all possible reaction forces at a configuration

q, when the commanded velocity is $\vec{\mathbf{v}}_d^c$. If, for example, the inverted velocity cone lies completely outside the friction cone, and it projects positively on the outgoing normal on the surface, then $\mathbf{F}_{reac}(q, d)$ consists of a single vector along one extreme ray of the friction cone. Let $\hat{\mathbf{F}}_{reac}(q, d)$ denote the set of all possible sensed forces at a configuration q, when the commanded velocity is $\vec{\mathbf{v}}_d^c$. $\hat{\mathbf{F}}_{reac}(q, d)$ is derived from $\mathbf{F}_{reac}(q, d)$ as follows:

- If $0 \in \mathbf{F}_{reac}(q, d)$, then $0 \in \hat{\mathbf{F}}_{reac}(q, d)$;
- If $\mathbf{F}_{reac}(q, d)$ contains a non-zero vector $\vec{\mathbf{f}}$, then all vectors $\hat{\vec{\mathbf{f}}}$ such that $\|\hat{\vec{\mathbf{f}}}\| = \|\vec{\mathbf{f}}\|$ and $|angle(\vec{\mathbf{f}}, \hat{\vec{\mathbf{f}}})| < \varepsilon$ are in $\hat{\mathbf{F}}_{reac}(q, d)$.

Computing the set $\hat{\mathbf{F}}$ for each edge and its intersection with other such sets is straightforward and takes constant time.

Force sensing is not able to distinguish points in the same edge, it can only recognize that the robot is in contact with some edge whose orientation is compatible with the sensed reaction force. If the reaction force on an edge may be null, this edge cannot be distinguished from the free space. Thus, if only force sensing is available, the kernel of a target set for a given commanded velocity consists of all contact edges that are entirely in the target and can neither generate a force measurement that could have been generated by an edge that does not belong to the target set, nor a null measurement.

3.5.3 Position sensing

In the case of position sensing, the sensory readings at a particular configuration do not depend either on the commanded or on the actual velocity. If the execution module makes use of only position sensing readings to decide whether the robot is in the goal or not, the computation of the kernel is conceptually simple: The kernel consists of all points of the target set $\mathcal{T}$ that cannot be confused with any point that does not belong to $\mathcal{T}$.

Two points of the configuration space are confusable if they may produce the same sensory reading. According to our position sensing uncertainty model, the distance

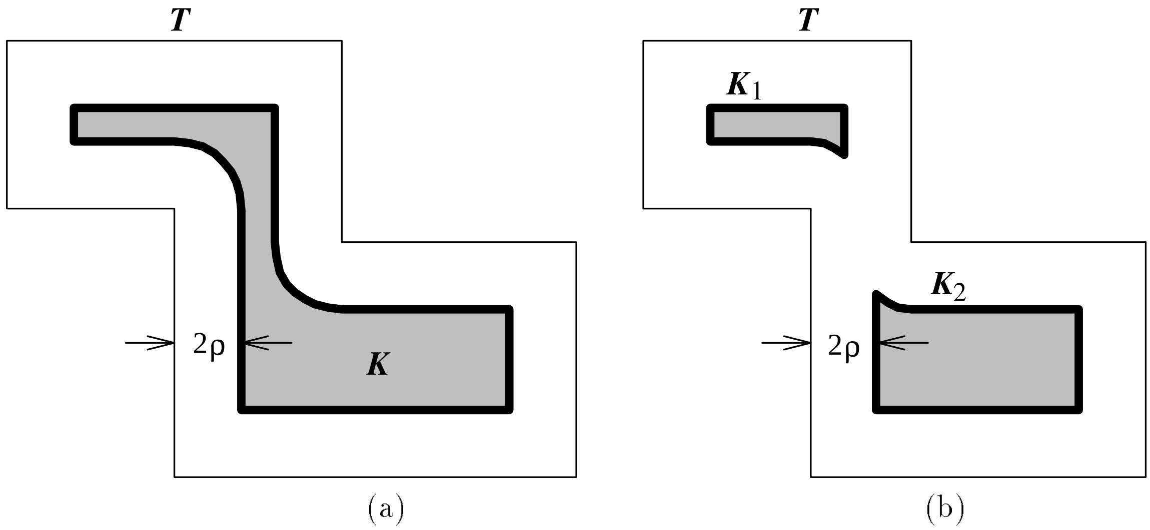

Figure 3.6: Computing the kernel in free space

between a point in the configuration space and the sensed configuration, when the robot lies at this point, is less than ρ, the position sensing uncertainty. Therefore, only points that are closer than 2ρ may produce the same sensory reading (and thus be confusable). In order to identify the points that are confusable with feasible configurations but do not belong to $\mathcal{T}$, we grow $\mathcal{C}_{feasible} \setminus \mathcal{T}$ by 2ρ. All the points of $\mathcal{T}$ that are not in this set constitute the kernel of $\mathcal{T}$.

In general, a target set $\mathcal{T}$ consists of polygonal regions, edges and points in $\mathcal{C}_{feasible}$. Edges and points are treated as degenerate polygons, so no special algorithms are needed to handle them. If a target region $\mathcal{T}$ lies in the free space without intersecting any other target regions, then its kernel $\mathcal{K}$ is found by shifting its edges inwards by 2ρ and drawing a circle of radius 2ρ at concave vertices. These lines are connected to form the kernel of the region (Figure 3.6). Of course, if $\mathcal{T}$ is thinner than 4ρ its kernel is the empty set. Also, the kernel of a connected set may be not connected, as is evident in the example of Figure 3.6(b).

If two or more target regions in the free space touch or overlap, we need to compute their union as a single set before we call the above algorithm. As an example, consider two adjacent rectangular regions. Their common edge should not be shifted inwards, as long as the robot does not need to know which of the two regions achieves. Indeed,

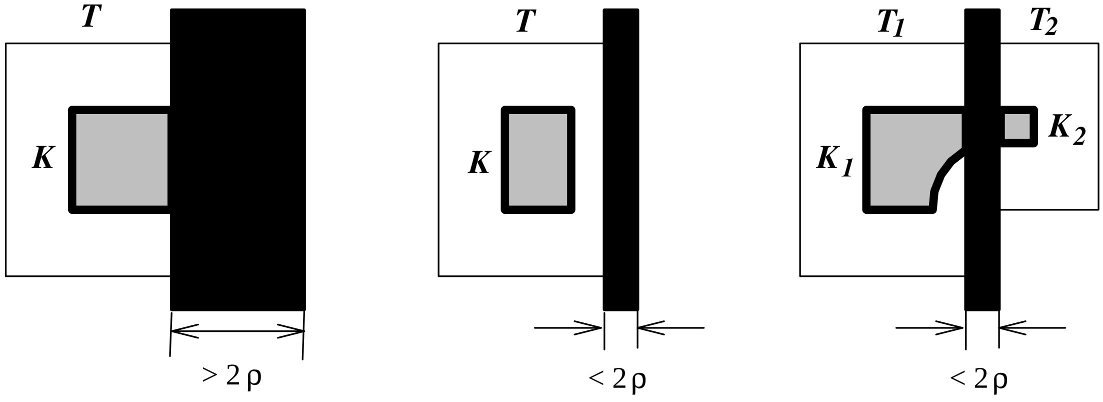

Figure 3.7: Computing the kernel in contact with obstacles

the points of the two regions whose distance from the common edge is smaller that 2ρ cannot be distinguished from points in the other region, but they are not confusable with points which lie outside both regions.

Another interesting case occurs when a target region is in contact with some obstacle. The interior points of the obstacle are not feasible, therefore we do not mind if they are confusable with points in the kernel. The general idea is that the edges of a target region that are in contact with obstacles should not be shifted inwards (Figure 3.7(a)). However, if the obstacle is too thin (its thickness is smaller than 2ρ), there are free space points that are confusable with points of the target set near the contact edge. In such a case, we compute the kernel by shifting inwards all target and obstacle edges that are not in contact with each other (Figure 3.7(b)). The same technique can be used to handle the case where another target region lies on the other side of the obstacle, and it is in contact with the obstacle (Figure 3.7(c)).

To formalize these ideas, we introduce the notion of the *cylsphere*. The cylsphere of an edge is the set of points whose distance from the edge is less than a real number r: $cylsphere(E, r) = \{q \in \mathcal{C} \mid dist(q, E) < r\}$. In order to compute the kernel $\mathcal{K}$ of a collection of target polygonal regions $\mathcal{T}_i$, we perform the following steps:

1. $\mathcal{K} = \emptyset$.

2. Select a target region $\mathcal{T}_i$.

3. Construct the maximal connected polygonal region S as follows: $\mathcal{T}_i$ is first included in S; then, every connected component of $\mathcal{CB}$ and target polygon $\mathcal{T}_j$, $j \neq i$, that shares an edge with a region already in S is iteratively[4] included in S, and it is marked as visited.

4. $S' \leftarrow S$.

5. For every edge E in the boundary of S, $S' \leftarrow S' \setminus cylsphere(E, 2\rho)$.

6. For all target regions $\mathcal{T}_j \in S$, add $\mathcal{T}_j \cap S'$ to the kernel $\mathcal{K}$.

7. If all target regions have been visited, return $\mathcal{K}$; otherwise, go back to step **2**.

Let $O(n)$ be the combined complexity of $\mathcal{C}_{contact}$ and $\mathcal{T}$. The region constructed at step **3** has $O(n)$ edges. Using a sweep-line algorithm, step **3** can be performed in $O((n+c)\log n)$ time, where $c \in O(n^2)$ is the number of intersections among target region edges. Step **5** can be performed in the same bound, only in this case c is the number of intersections among cylspheres. $\mathcal{T}_j \cap S'$ can be obtained from the sweep algorithm with the same time complexity.

Once the kernel has been computed, the remaining task is to extract a termination condition that is guaranteed to be true when the robot is in the kernel and false when the robot is outside the target set (so that the robot does not stop before reaching the target set). If $\hat{q}$ is the sensed configuration of the robot, the termination condition that corresponds to kernel $\mathcal{K}$ is $\mathbf{tc}(\hat{q}) \equiv dist(\hat{q}, \mathcal{K}) < \rho$. Equivalently, the robot terminates its motion whenever $\hat{q}$ lies within a set that we get by isotropically pushing out the boundary of the kernel by ρ. By construction of the kernel, all actual configurations that are compatible with sensory readings in this set lie in the target set. Figure 3.8 shows an example with a single target region $\mathcal{T}$ in contact with a C-obstacle $\mathcal{CB}$. The kernel $\mathcal{K}$ is the inner darker region; the set that contains the sensory readings that make the termination condition evaluate to true is outlined with the thin dashed line; and the set of configurations where the robot may stop is outlined with the thicker dashed line. Note that this set is a proper subset of the target set.

[4] In other words, S is the transitive closure of all connected components of $\mathcal{CB}$ and target polygons touching $\mathcal{T}_i$.

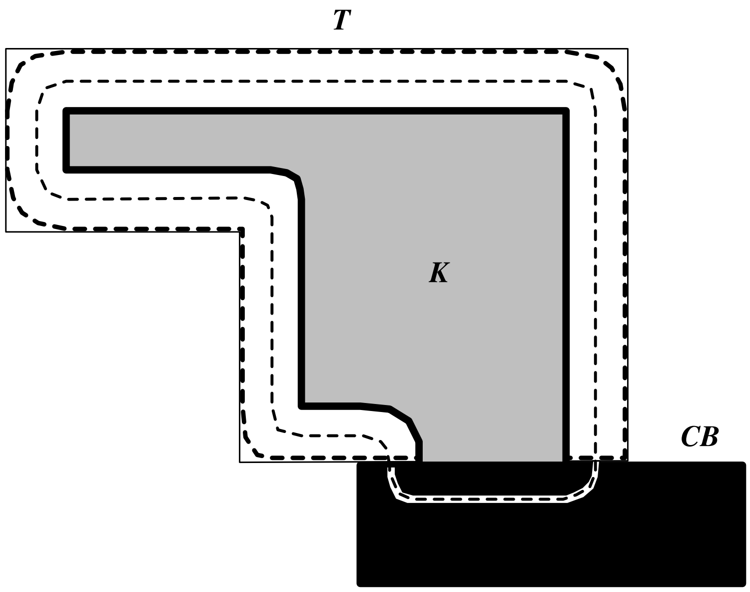

Figure 3.8: The kernel and the termination condition

3.5.4 Combination of sensing

Usually, robots have both position and force sensors, and can detect sticking, too. We now investigate how to compute the kernel for the combination of all three sensing schemes.

Sticking does not entail recognizing configurations, thus it does not interact with the two other kinds of sensing; it merely contributes a number of sticking edges to the kernel.

Position and force sensing do interact and create kernels that are bigger than the union of the kernels computed individually for each sensor. Such points, that are in neither individual kernel, but are in the combined kernel, must be in contact space so that both sensors can be used. Indeed, points in contact space can now be recognized using both position and force sensing, if they are either more than 2ρ apart, or if they cannot generate the same reaction force. This computation can be done as follows:

For each contact edge $E \in \mathcal{T}$:

1. If $\hat{\mathbf{F}}_{reac}(E, d)$ contains the null vector, then the edge is discarded.

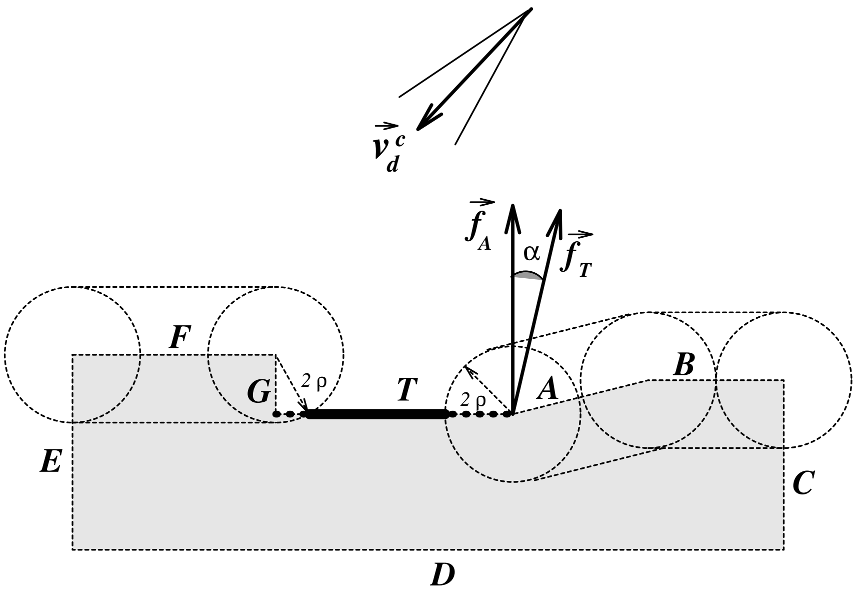

Figure 3.9: The kernel of a target in contact space

2. Otherwise, $S \leftarrow E$.

3. For every contact edge $C \notin \mathcal{T}$:
If $\hat{\mathbf{F}}_{reac}(E, d) \cap \hat{\mathbf{F}}_{reac}(C, d) \neq \emptyset$, then $S \leftarrow S \setminus cylsphere(E, 2\rho)$.

4. S is added to the kernel.

Consider the example of Figure 3.9. The goal is a single edge T of a C-obstacle that is depicted in light grey for clarity. The commanded velocity is such that the inverted control uncertainty cone is outside the friction cones of edges F, T, A and B, and projects positively on the outgoing normal of these edges (i.e., the robot is guaranteed to slide on these edges). The robot is also guaranteed to stick on C and G, and move away from D and E.

The sticking termination condition yields no kernel edges, because no sticking is possible on T. Using the above algorithm, we visit all edges in contact space, and if the reaction forces they may produce are confusable with the reaction force that may be produced by T, we subtract their cylspheres from T. The only reaction force $\vec{\mathbf{f}}_T$ that can be generated by T lies along the right ray of the friction cone. This cannot be confused with the forces generated by D and E (zero force), or C and G ($\vec{\mathbf{f}}_T$ lies

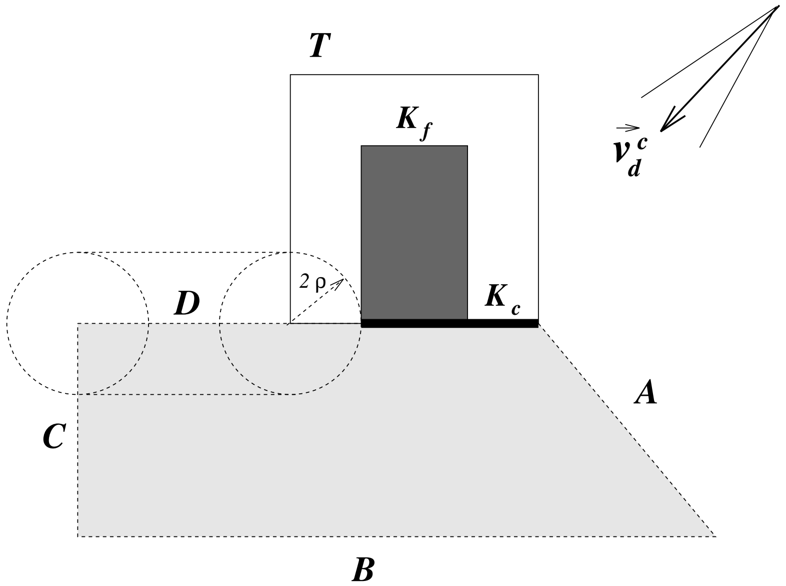

Figure 3.10: The kernel of a target lying both in free and in contact space

outside the friction cone of C or G). Edges F and B are parallel to T so they generate the same force $\vec{\mathbf{f}}_T$. They are confusable with T, and their cylspheres of radius 2ρ are subtracted from T. While B is too far away to have any effect, the cylsphere of F cuts a small part of T's left end. Edge A can also generate a single reaction force $\vec{\mathbf{f}}_A$. Since the angle α between $\vec{\mathbf{f}}_T$ and $\vec{\mathbf{f}}_A$ is smaller than the force uncertainty ε, edges T and A are confusable relative to force sensing. The cylsphere of A cuts a piece of length 2ρ from the right part of T. The remaining part of T, depicted as a solid black line in the figure, is the desired kernel.

As a further example consider Figure 3.10. Here, the target region $\mathcal{T}$ lies both in free and contact space. Its kernel consists of the free space set $\mathcal{K}_f$ and of the contact edge $\mathcal{K}_c$. Note that neither $\mathcal{K}_c$ is included in $\mathcal{K}_f$, nor $\mathcal{K}_f$ is included in $\mathcal{K}_c$. This means that by considering both position and force sensing simultaneously we create larger kernels (and consequently preimages) than either of them would produce on its own.

3.6 Computation of backprojections

After the kernel has been computed as a collection of free space regions and contact edges and vertices, we wish to compute its backprojection in order to find out whether it includes the initial region or not.

An algorithm for this task computes the backprojection of each kernel component independently and then merges them in a second pass [32, 52]. This algorithm has been generalized by Canny and Donald [23] to a one-pass algorithm that generates the backprojection of any region $\mathcal{K}$ described as the union of contact edges and polygonal regions in $\mathcal{C}_{feasible}$. The algorithm sweeps a line L across the plane perpendicularly to the commanded velocity $\vec{\mathbf{v}}_d^c$. The sweep starts at a position of L where it is tangent to $\mathcal{K}$, with $\mathcal{K}$ lying entirely on the side of L pointed to by the vector $-\vec{\mathbf{v}}_d^c$. The sweep proceeds in the direction of the vector $-\vec{\mathbf{v}}_d^c$. During the sweep, the algorithm maintains the status of L, i.e., the description of its intersection with $\mathcal{C}_{contact}$, $\mathcal{K}$, and rays parallel to the edges of the inverted velocity cone that are erected from contact vertices and goal vertices. This status changes at events occurring when L passes through a vertex of $\mathcal{C}_{contact}$, a vertex of $\mathcal{K}$, the intersection of two rays, the intersection of a ray and $\mathcal{C}_{contact}$, or the intersection of a ray and $\mathcal{K}$. At each of these events, the algorithm updates the status of the line and the list of future events. During the sweep, the algorithm traces out the backprojection. The sweep stops when the last event is encountered.

The key idea of this algorithm is that the backprojection's boundary consists of straight segments, and that at every event there exists a local criterion (which does not depend on what will happen later during the sweep) for deciding how to trace the backprojection's boundary. This criterion leads to erecting rays from both contact and goal vertices, in order to prevent the backprojection from containing configurations from where a motion may reach a vertex or an edge, where it may stick or slide away from the goal with no possibility of returning to it. The algorithm does not require the backprojection to be bounded, but it does require that the actual velocity of the robot never project negatively into the $\vec{\mathbf{v}}_d^c$ direction. This condition is achieved if the friction cone is larger than the control uncertainty cone, i.e., $\phi \geq \theta$.

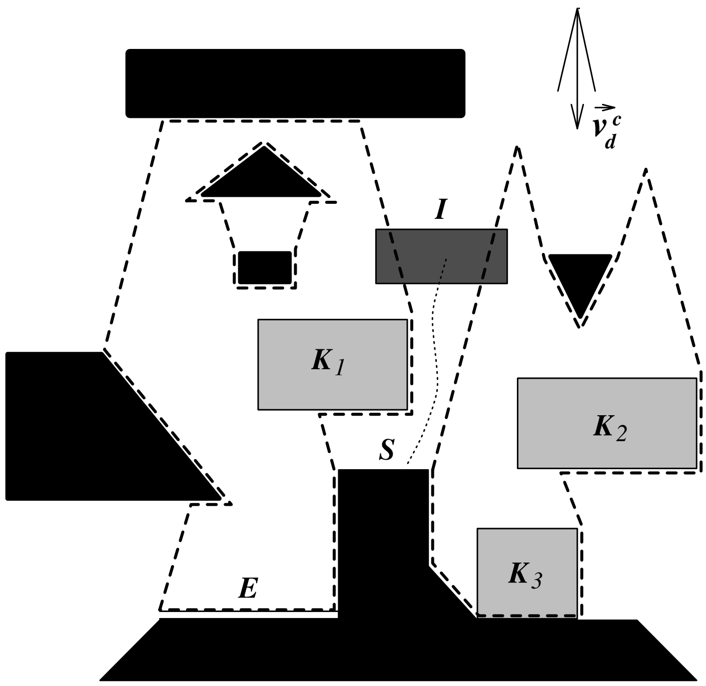

Figure 3.11: Computation of Backprojection

Let n be the number of vertices in $\mathcal{C}_{contact}$. Assume that the combined complexity of $\mathcal{C}_{contact}$ and $\mathcal{K}$ (the number of vertices in $\mathcal{C}_{contact}$ plus the number of vertices of $\mathcal{K}$ plus the number of intersections between $\mathcal{K}$ and $\mathcal{C}_{contact}$) is $O(n)$. The sweep-line algorithm erects $O(n)$ rays. The contour of the backprojection consists of goal edges (edges bounding $\mathcal{K}$), rays in free space and contact edges. Each erected ray can contribute at most one edge to the backprojection. Each goal edge contributes at most one backprojection edge. Each non-goal edge in contact space can contribute one or several backprojection edges, but whenever it contributes k backprojection edges, it also "consumes" $O(k)$ edges in free space. Hence, the boundary of the backprojection consists of $O(n)$ edges. Since each ray is interrupted at its first intersection with a goal edge, a contact edge, or another ray, the backprojection can be computed in $O(n \log n)$ time.

Figure 3.11 illustrates the backprojection of a kernel that consists of three regions K_1, K_2, K_3 and an edge E. The black polygons are C-obstacles. The backprojection (outlined with a dashed line) consists of two disconnected components, one of which contains a hole. Since the initial position (the grey rectangular region $\mathcal{I}$) is not completely contained in the backprojection, the robot is not guaranteed to reach the kernel if it starts from $\mathcal{I}$ and follows $\vec{\mathbf{v}}_d^c$. Indeed, it is possible for the robot to stick on vertex S and never reach the kernel.

3.7 Implementation and experimentation

We have implemented a motion planner based on the preimage backchaining approach (for details see [52, 54, 55]). This planner computes preimages using either the sticking termination condition, or the combination of position and force sensing, or all three together. The user inputs the description of the configuration space, the goal region $\mathcal{G}$, and the initial region $\mathcal{I}$. If successful, the planner returns a motion plan in the form of a sequence of unit commanded velocities and the associated sequence of computed preimages. The method used for computing the preimages determines the termination condition of every motion command in the plan. The planner is implemented in Allegro Common Lisp and runs on an Apple Macintosh II computer.

The planner constructs a graph of preimages by considering commanded velocities with the orientations $\{k\pi/N\}_{k=0,\ldots,2N-1}$, where N is input by the user. In our experiments, we used $N = 2$ or 4. The program searches the graph in a breadth-first fashion, but other (and perhaps more efficient) search techniques could have been used as well.

The algorithms implemented in the planner are essentially those described in the previous sections, with minor variations. In particular, the backprojection of a region is computed using an algorithm similar to that described in [52], rather than the sweep-line algorithm (which might be faster). The planner also approximates conservatively the generalized polygonal kernels by polygonal regions. These changes have no major impact on the visible operations of the planner.

In all the examples shown, the initial region is a single point; in the figures, it is the center of a disc that depicts position uncertainty. The control uncertainty cone and the force uncertainty cone are not depicted.

Figure 3.12(a) illustrates a plan generated with the sticking termination condition.

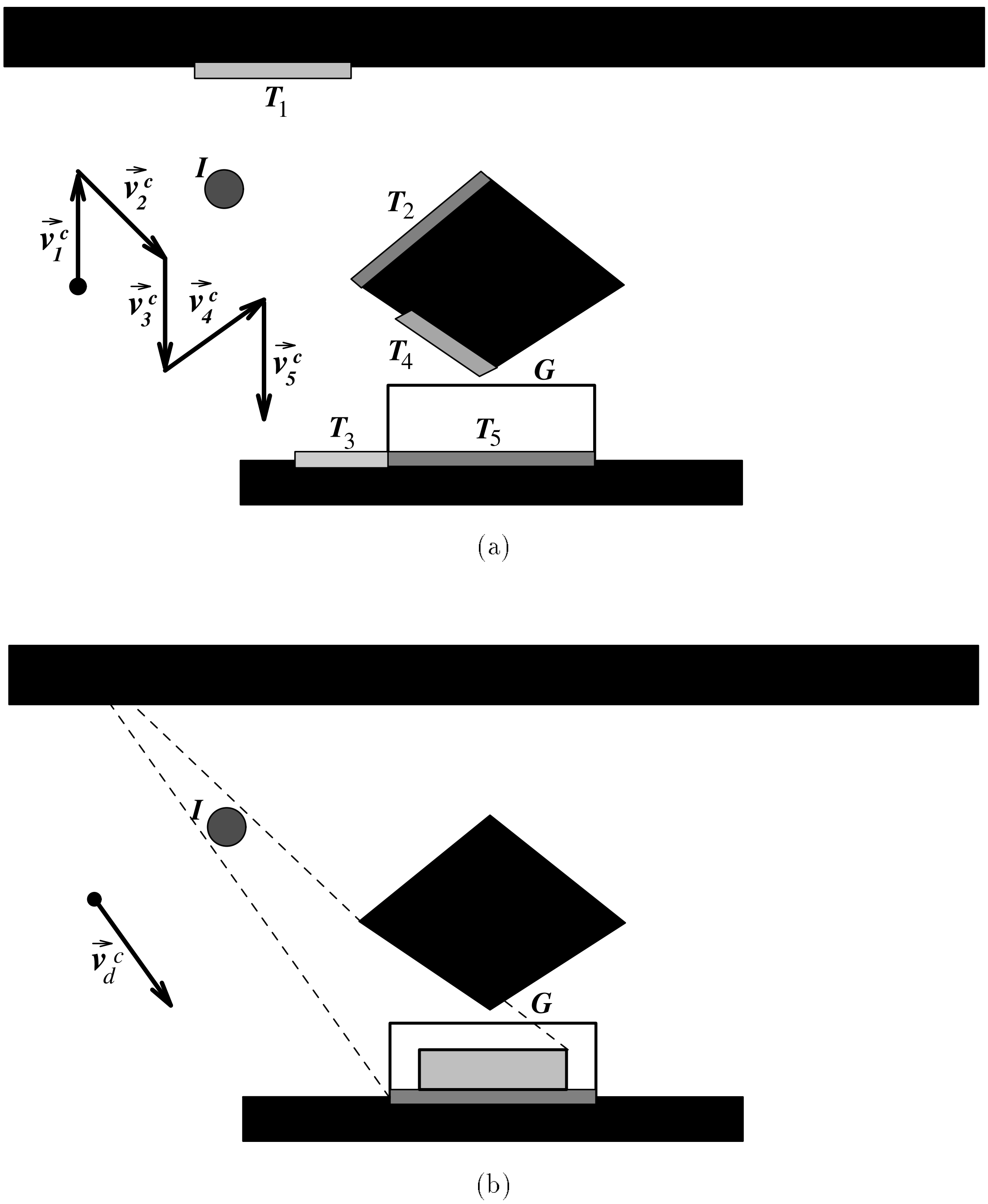

Figure 3.12: Example 1

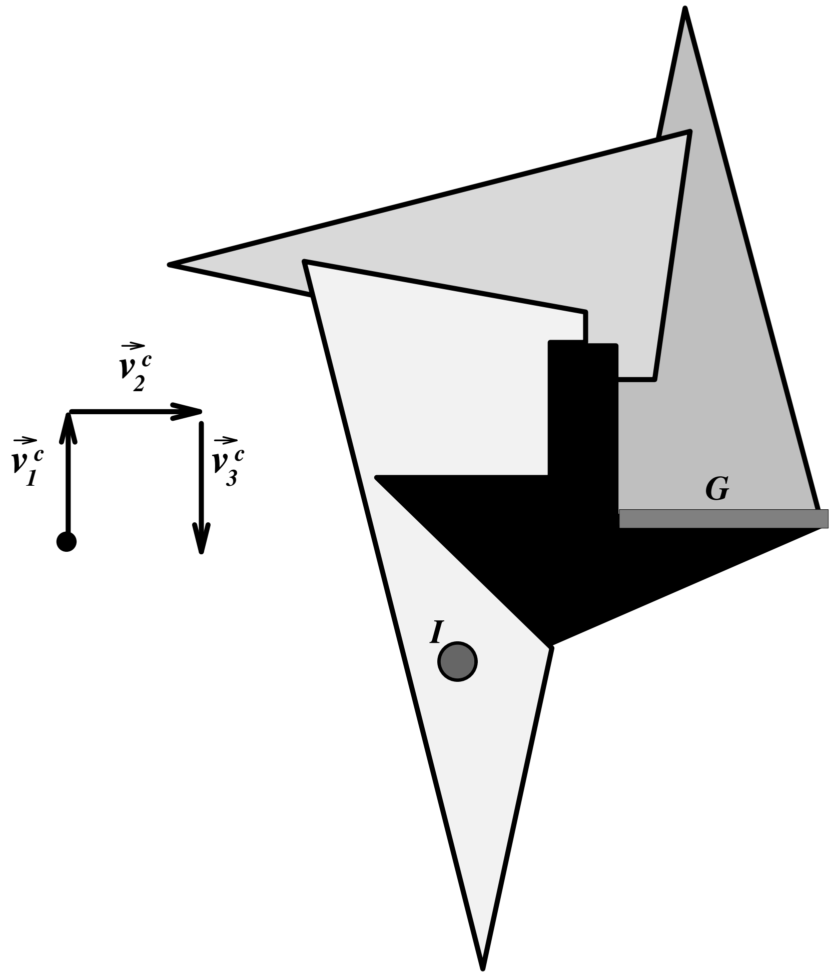

Figure 3.13: Example 2

The computed preimages are not shown in the figure. The goal is the rectangular region $\mathcal{G}$. The plan constructed consists of 5 steps, defined by the commanded velocities $\vec{\mathbf{v}}_1^c$ through $\vec{\mathbf{v}}_5^c$. The corresponding sticking edges are $\mathcal{T}_1$ through $\mathcal{T}_5$. ($\mathcal{T}_i$ is the intersection of the preimage of $\mathcal{T}_{i+1}$ and the sticking subset of $\mathcal{C}_{contact}$ for $\vec{\mathbf{v}}_i^c$. A motion commanded along $\vec{\mathbf{v}}_i^c$ and issued from within $\mathcal{T}_{i-1}$ is guaranteed to terminate in $\mathcal{T}_i$, but some configurations in $\mathcal{T}_i$ may not be reachable.) This plan is typical of the "bouncing on obstacles" that is caused by the sticking termination condition. On the other hand, if we use position and force sensing, the resulting plan consists of a single motion command along $\vec{\mathbf{v}}_d^c$ (Figure 3.12(b)).

Figure 3.13 shows an example which cannot be solved by using sticking on obstacle edges. The goal is just one obstacle edge marked with G in the figure. In this case, there are no edges facing each other so that the robot can bounce between them. The combination of position and force sensing creates the three-step plan shown in the figure.

3.8 On the termination condition's knowledge

We now return to the computation of the kernel in the case where only position sensing is available, and examine how the method described in this chapter relates to the generic method described in Chapter 2. Consider the simple case where the target set $\mathcal{T}$ is a rectangular region of the free space. According to the method described in this chapter its kernel $\mathcal{K}$ is found by shrinking $\mathcal{T}$ by 2ρ. Then $\mathcal{T}$'s preimage is computed as the backprojection of $\mathcal{K}$ (Figure 3.14(a)). The termination condition (i.e., the set of all sensory values for which the robot will stop) is found by isotropically growing the kernel by ρ. (In Chapter 2 we called this set $\hat{S}_{\mathbf{tc}}$.) The set of all configurations that are compatible with the termination condition ($\mathbf{INT}(\hat{S}_{\mathbf{tc}})$, according to the terminology of Chapter 2) is found by isotropically growing the kernel by 2ρ. This set represents configurations where the robot may stop, so all the points in this set that are not in the target set must be completely avoided. The method we proposed in Chapter 2 was to compute the backprojection of the kernel considering $\mathbf{INT}(\hat{S}_{\mathbf{tc}}) \setminus \mathcal{T}$ as an artificial obstacle that must be completely avoided (no sliding on it is allowed).

The kernel computation presented in this chapter always creates sets $\mathbf{INT}(\hat{S}_{\mathbf{tc}})$ which are subsets of the target; therefore, no artificial obstacles are induced. However, this method does not necessarily yield the biggest possible backprojection. Indeed, only obstacles that lie between the backprojection and the kernel cause the backprojection to shrink. It is possible to grow the kernel without inducing any obstacles between the backprojection and the kernel. In this case, a bigger kernel produces a bigger backprojection.

Consider again our simple example. If we use a kernel where only the top edge of

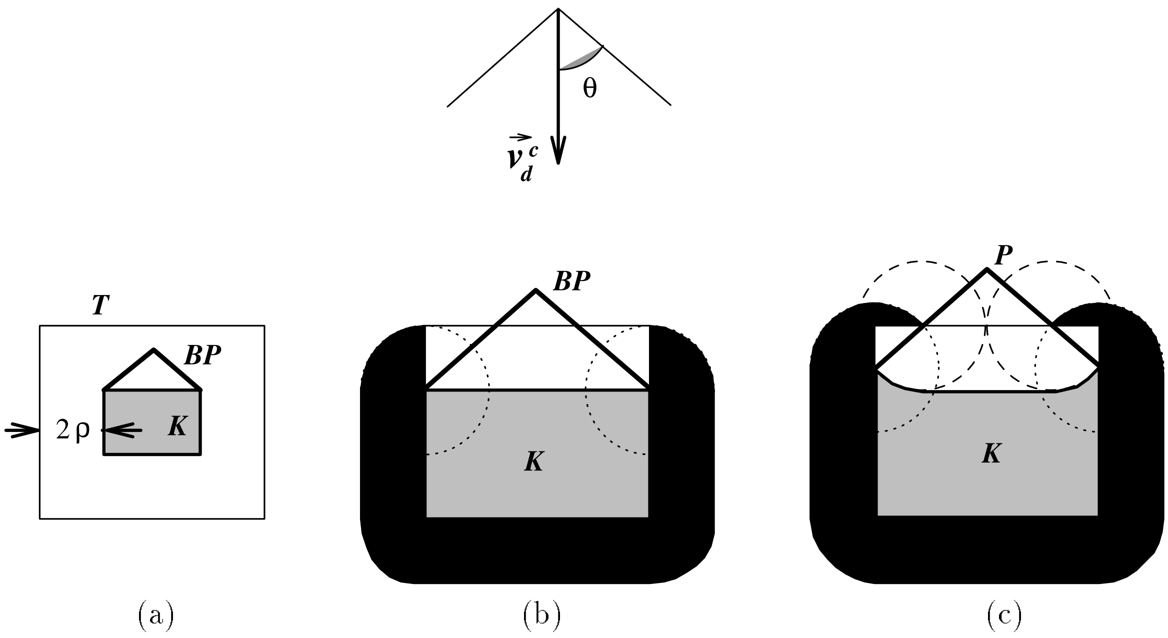

Figure 3.14: Rectangular goal in free space

the rectangle is pushed inwards by 2ρ, the induced artificial obstacle (depicted in black in Figure 3.14(b)) leaves the top edge of the kernel free. Thus, for the commanded velocity shown in the figure, we get a bigger backprojection.

Yet, we can do better. Figure 3.14(c) shows that we can keep pushing the rays of the backprojection outwards, as long as the artificial obstacle does not interfere with the achievement of the kernel. The artificial obstacle starts creeping on the upper edge of the target set, and at some point it reaches the rays of the backprojection. At this point, the length of the segments of the two rays of the backprojection that lie in the target set is equal to 2ρ. Each ray can be moved outwards, up to the point where the vertex of the kernel (where this ray is incident) starts creating an artificial obstacle along the direction of this ray in front of the target set. Clearly, it is not possible to push the rays any further. Therefore, this is the maximal backprojection for the selected commanded velocity, i.e., the *preimage* of $\mathcal{T}$.

3.8.1 Maximal preimage of a convex set

The above method is applicable to arbitrary convex sets. Given a commanded velocity $\vec{\mathbf{v}}_d^c$ and a convex target set $\mathcal{T}$, the preimage of $\mathcal{T}$ is found by erecting two rays (a left and a right one) parallel to the edges of the inverted control uncertainty cone, such that they both cut segments of length 2ρ in the target set. The points where these rays exit $\mathcal{T}$ are called the *extreme points* of the kernel. The computation of the extreme points of the kernel can be done with two line-sweeps, one for each ray. During a sweep, we track the length of the segment that is the intersection of the sweep line and the target set. The orientation of the sweep line remains constant. We call the mapping of the sweep line location onto the length of its intersection with the target set the *chord function* $\kappa(\delta)$. The location of the line is represented by its distance δ from the point where the line is tangent to $\mathcal{T}$, with $\mathcal{T}$ being ahead of the line (hence, $\kappa(\delta) = 0$ when $\delta < 0$). When $\mathcal{T}$ is a convex set, $\kappa(\delta)$ consists of an increasing and a decreasing part and has a single global maximum. It is continuous except in the case where the sweep line is parallel to an edge of $\mathcal{T}$. The algorithm stops when the chord function becomes equal to 2ρ while increasing, or at $\delta = 0$ if $\kappa(0) \geq 2\rho$. The worst-case performance of this algorithm is $O(n \log n)$, where n is the complexity of $\mathcal{T}$.

In the absence of obstacles and other target sets, the two rays found with the above method constitute the upper boundary of the preimage $\mathcal{P}$. It remains to construct the kernel $\mathcal{K}$ that corresponds to $\mathcal{P}$. The kernel, which is part of the preimage, will define the lower boundary of $\mathcal{P}$. As mentioned before, no kernel point must be closer than 2ρ to any point in $\mathcal{P} \setminus \mathcal{T}$. Therefore, we can find the kernel by growing $\mathcal{P} \setminus \mathcal{T}$ by 2ρ and subtracting it from $\mathcal{T}$. In Subsection 3.5.3 we defined the notion of the cylsphere of an edge to be the set of points whose distance from the edge is less or equal to some number. Similarly, the cylsphere of any set of points is the set which contains all the points whose distance from the first set is smaller than a given distance:

$$\mathcal{K} = \mathcal{T} \setminus cylsphere(\mathcal{P} \setminus \mathcal{T}, 2\rho)$$

In most cases this computation can be performed by connecting the extreme points of the kernel with two lines, the *lower* and the *upper*. The lower line follows the

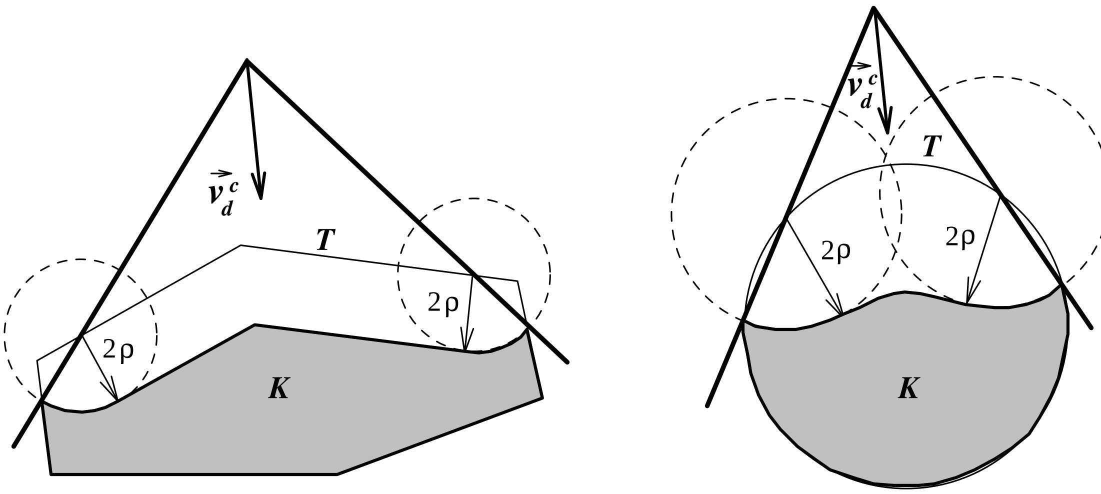

Figure 3.15: Maximal backprojections

lower boundary of the target set. The upper line is parallel to the upper boundary of the target set, at a distance 2ρ from it inwards. The upper line is connected to the extreme points of the kernel with circular arcs of radius 2ρ centered at the points where the rays enter the target set (see Figure 3.15). The termination condition is then found by growing the kernel isotropically by ρ.

Sometimes, though, the upper boundary may not lie entirely in the target set (see Figure 3.16). In this case, the above algorithm does not work. Then, we have to push the preimage rays inwards, until the part of it outside $\mathcal{T}$ and expanded by 2ρ is tangent to the lower boundary of $\mathcal{T}$. Thus, we obtain the widest possible preimage. The sweep-line algorithm used to find the original points can be adapted to handle this case as well.

3.8.2 About the existence of a unique maximal preimage

Another complication arises when the target consists of more than one convex sets. Consider the example of Figure 3.17(a). The target set consists of two parallelograms $\mathcal{T}_1$ and $\mathcal{T}_2$. If we use the above algorithm independently for each set, then the computed preimage may be incorrect. Assume, for example, that there exists a point A in the preimage of $\mathcal{T}_2$ and outside both $\mathcal{T}_1$ and $\mathcal{T}_2$, whose distance from the kernel of

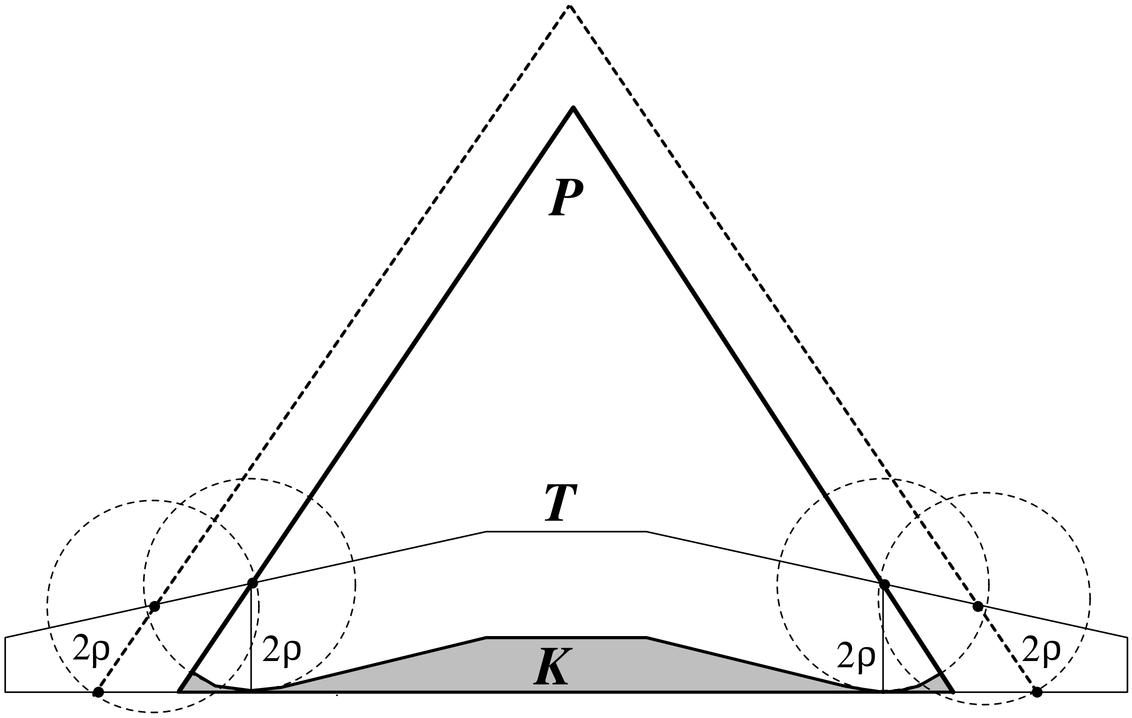

Figure 3.16: Preimage of a thin target region

$\mathcal{T}_1$ is less than 2ρ. A robot starting from A may actually think that it is at point B in the kernel of $\mathcal{T}_1$ and terminate its motion outside both target sets.

The problem is that the artificial obstacle created by the kernel of $\mathcal{T}_1$ intersects with the preimage of $\mathcal{T}_2$. When something like this occurs, there are two ways to proceed:

(a) We can leave the preimage of $\mathcal{T}_2$ intact and shrink the kernel of $\mathcal{T}_1$, so that their minimum distance is 2ρ. In practice, we can find the 2ρ-cylsphere of the preimage of $\mathcal{T}_2$, subtract it from $\mathcal{T}_1$, and then find the kernel and preimage of their difference. In the example of Figure 3.17(b), the kernel of $\mathcal{T}_1$ is split into two non-connected components, resulting in a much smaller preimage.[5]

(b) We can shrink the preimage of $\mathcal{T}_2$ up to the point that the upper boundary of the kernel of $\mathcal{T}_1$ remains intact (see Figure 3.17(c)). In this case, we find

[5] See the next subsection on how to find the preimage of a concave set.

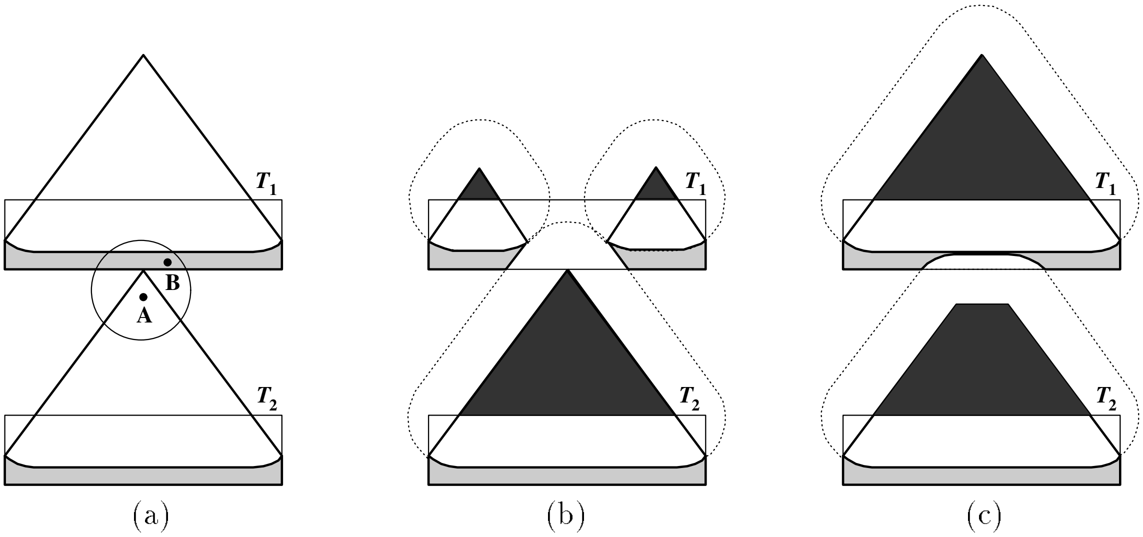

Figure 3.17: Preimage of non-connected convex sets

the 2ρ-cylsphere of the upper boundary of the kernel of $\mathcal{T}_1$ and subtract it from the preimage of $\mathcal{T}_2$.

We have thus computed two possible preimages of the target set $\mathcal{T}_1 \cup \mathcal{T}_2$. The union of these two preimages is not a preimage of the target set, because it contains points like A, where the robot may stop, although it is not in the target set. This means that there does not exist a unique two-dimensional set that contains all possible preimages of $\mathcal{T}_1 \cup \mathcal{T}_2$. In other words, we have proven that *there are cases when a unique maximal preimage*[6] *does not exist*. This result was first reported by Erdmann [32], who used non-compact sets to build a counterexample.

A complete algorithm must consider all possible maximal preimages, but it is very hard to do so because their number may be infinite (like in the example of Figure 3.17). One possibility is to compute the preimage independently for each set, and check for inclusion of all initial-region points. If this is true, there is hope that a guaranteed plan may exist. Then, we can find all points that are actually attainable by the robot during its motion (the forward projection), and check whether there is any chance for the robot to stop outside the goal. We can do this by computing the 2ρ-cylsphere of the part of the forward projection that lies outside the target set,

[6]A maximal preimage is a preimage that is not a proper subset of any other preimage.

and subtracting from the kernel any points that lie in this cylsphere. If the preimage of the reduced kernel still contains all initial-region points, then a guaranteed plan exists.

3.8.3 Maximal preimage of a concave set

The computation of the preimage of a concave set is significantly more complicated. The intersection of the sweep-line with the concave set may consist of more than one segments, so the chord function is a list of length functions $\kappa_i(\delta)$. These functions do not have the smooth behavior of the chord function of a convex set. They may have local maxima and minima, they may split into two new functions, or two of them may merge into one. Nevertheless, the sweep-line algorithm uses similar criteria to detect kernel extremities. It returns all points δ_0 such that $\kappa_i(\delta_0) = 2\rho$ while κ_i is increasing, $\delta_0 = 0$ if $\kappa_i(0) \geq 2\rho$, and all points δ_0 where two functions κ_i and κ_j merge, with $\lim_{\delta \to \delta_0^-} \kappa_i(\delta) \leq 2\rho$, $\lim_{\delta \to \delta_0^-} \kappa_j(\delta) \leq 2\rho$, and $\lim_{\delta \to \delta_0^-} (\kappa_i(\delta) + \kappa_j(\delta)) \geq 2\rho$.

The above criteria may be satisfied more than once, so we may end up with multiple kernel extreme points and multiple preimage components. Furthermore, these components may interact with each other, in which case a unique maximal preimage does not exist. Figure 3.18 presents a complicated example with a concave target set. The kernel and the preimage consist of four disjoint components each, but no unique maximal preimage exists.

3.8.4 Recognition power

The termination conditions created with the above method, may ask the robot to stop, despite a sensory reading that is compatible with configurations outside the target set. They can do that because of the additional knowledge that is embedded in them, knowledge about the *initial state* and knowledge about the *final state*. By knowing the initial state (i.e., the preimage), the robot can rule out interpretations that are not achievable. By knowing the final state (i.e., the kernel) the robot can rule out interpretations that are beyond the kernel, since it is guaranteed that the robot will stop in the kernel.

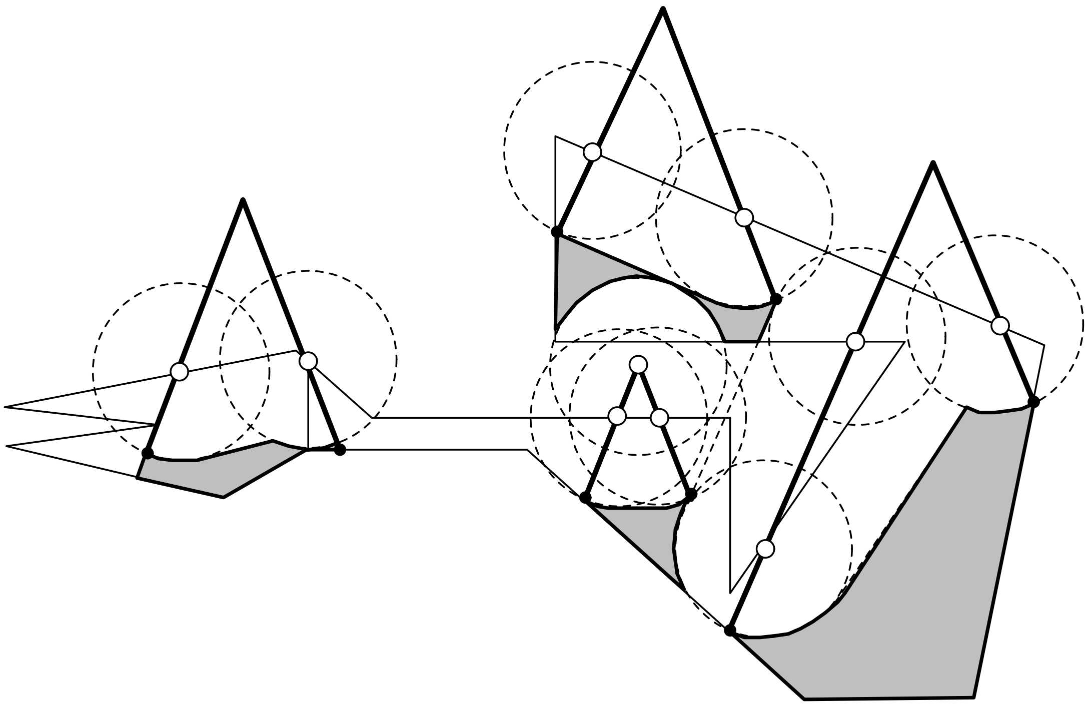

Figure 3.18: Maximal backprojection of a concave set

3.9 Omnidirectional backprojections

In the implemented planner of Section 3.7 the major bottleneck is the selection of the commanded velocity direction at each backchaining step. Not only do we lose completeness by discretizing the considered directions, but the resulting search takes exponential time in the number of backchaining steps.

The discretization problem can be solved if we find a way to efficiently compute for each point of the configuration space all commanded velocity directions that are guaranteed to bring a robot starting from that point into the target set $\mathcal{T}$.

Definition 3.1 (Nondirectional backprojection) *We call* nondirectional backprojection *of a target set $\mathcal{T}$ the two-dimensional set $\mathcal{NBP}(\mathcal{T}) \in \mathcal{C}$, consisting of all configurations, for which the above mapping produces a non-empty set of directions. In other words, $q \in \mathcal{NBP}(\mathcal{T}) \Leftrightarrow \exists d \in S^1 : \ q \in \mathcal{BP}(d;\mathcal{T})$.*

Definition 3.2 (Omnidirectional backprojection)
We call omnidirectional backprojection *of a target set $\mathcal{T}$ the three-dimensional set $\mathcal{OBP}(\mathcal{T}) \in \mathcal{C} \times S^1$, such that $(q, d) \in \mathcal{OBP}(\mathcal{T}) \Leftrightarrow q \in \mathcal{BP}(d; \mathcal{T})$.*

The knowledge of the nondirectional backprojection is sufficient to allow us to decide whether a one-step plan that connects a given initial region with a target set exists, but it does not facilitate the generation of the plan. In order to be able to derive the plan, we need to know the omnidirectional backprojection of the target set.[7] In the future, whenever we wish to refer specifically to the regular backprojection of a set, we will call it the *directional backprojection.*

The omnidirectional backprojection can be thought of as a stack of directional backprojections (*slices*) computed for various values of the commanded direction of motion. In order to create a one-step plan, we need to find out which of these slices completely include the initial region. This is the only reason we need the omnidirectional backprojection. Essentially, what we need is a one-dimensional boolean function that maps the various choices of the commanded direction of motion onto $[0, 1]$ (0 if the initial region is not included in the directional backprojection for this direction, 1 otherwise). It turns out that this function changes value only at a polynomial number of *critical orientations* of the commanded velocity vector. These values divide S^1 into regular sets (open intervals and points). If we compute the value of the boolean function for one point of such a set, we know it for all the other set points.

This observation was first made by Donald [25], though the notion of omnidirectional backprojections was introduced earlier in the original preimage backchaining paper [69]. The value of the boolean function can be computed by a sweep-line algorithm similar to the one that generates directional backprojections (Section 3.6). Moreover, this algorithm creates data structures that can be used to help update the value of the boolean function at each point considered, without having to call the sweep-line algorithm again. Donald proposed an $O(n^4 \log n)$ algorithm, with n being the number of obstacle edges, to compute nondirectional backprojections and

[7] There has been a fair amount of confusion between the notions of nondirectional and omnidirectional backprojection. Usually, the term nondirectional backprojection is used to describe both. We believe that the adjective "omnidirectional" is a better description of the three-dimensional set, since directional information is a vital part of it.

embedded this algorithm into a polynomial one-step planner. Briggs [8] reduced the time complexity of computing a nondirectional backprojection to $O(n^2 \log n)$.

Unfortunately, this algorithm cannot be generalized to produce multi-step plans. The reason is that as the backprojection varies with the choice of the commanded direction of motion, its kernel (the target set of the next backchaining step) varies too. Ultimately, of course, we are only interested in finding out whether a sequence of successive preimages includes the initial region or not. The answer to this question changes a large number of times (the number of changes is an exponential function of the number of backchaining steps), so it is not efficient to track it.

This result is typical of algorithms that use the preimage backchaining method. The core of our work, starting at the next chapter, shows how to avoid exponentiality by using carefully selected models of the robot, the workspace, and the representation of uncertainty.

Chapter 4

Landmark-Based Navigation

In this chapter we introduce the notion of *landmarks* as *islands of perfection* within an otherwise imperfect world. Using specific assumptions, we present a sound, complete and polynomial motion planning algorithm with uncertainty. The algorithm is based on the preimage backchaining method [69], and makes use of the idea of the omnidirectional preimage [25, 8]. We present an efficient way to compute and store a sufficient representation of an omnidirectional preimage, based on the idea of critical slices. Finally, we show several planning examples, and the simulated execution of the generated plans.

4.1 A new approach to motion planning with uncertainty

The problem of general motion planning with uncertainty is intrinsically hard. Several exponential lower bounds have been established, even for restricted subclasses of the problem. In particular, the three-dimensional problem attacked by the preimage backchaining method has been shown to be PSPACE-hard [77], as well as NEXPTIME-hard [15]. Canny gave a very general algorithm that takes double-exponential time in the number of plan steps (i.e., the number of motion commands in the plan) [14]. This algorithm assumes that the envelope of trajectories generated by the control system can be described algebraically, and that the robot has imperfect position and velocity sensing. The motion planning problem is reduced to the satisfiability of a semi-algebraic formula with alternating universal and existential quantifiers, like the one we introduced in Section 2.5. As Canny notes, "the existential quantifiers represent the actions taken by the plan executor, like choosing the direction to move or the time to stop moving, and the universals represent the actions of nature, namely the sensor readings." Its double-exponential bound stems precisely from the alternation of the quantifiers.

One exponential can be trimmed by eliminating sensing, so that the execution module does not need to make choices based on sensor readings. In this case, the number of alternations between universal and existential quantifiers is equal to the number of steps of the plan. Sensorless plans were first investigated in [37]. Donald [25] created a planner for a sensorless robot moving in a two-dimensional configuration space. The robot stops only by sticking on obstacle edges. Donald showed that this planner creates one-step motion plans in $O(n^4 \log n)$ time. Briggs [8] improved this bound to $O(n^2 \log n)$ using amortization techniques. Nevertheless, multi-step plans still require time exponential in the number of steps.

To make things worse, the number of steps for which a plan may exist grows with the complexity of the environment. Even in a three-dimensional polyhedral world, this number may grow exponentially with the environment complexity [15]. In order to avoid exponentiality, we need to find a way to limit the maximal number of steps of successful plans. Friedman [41] was able to do it by considering motions in the interior of a simple polygon, and assuming that sensing is possible only when the robot is in contact with one of the polygon edges. Under these assumptions, the maximum number of steps of a successful plan is bounded by the number of edges. Essentially, the polygon edges in Friedman's model are what in this work we call *landmarks*. Only by limiting sensing in a given number of regions in the workspace does it become possible to bound the number of planning steps, and thus to derive polynomial planning algorithms. This is exactly the key idea behind the work presented in this chapter. The framework, though, is significantly more general than Friedman's simple model.

Notions similar to landmarks have been introduced in the literature with a variety of names, e.g., *atomic regions* [12], *landmarks* [64], *signature neighborhoods* [70], *perceptual equivalence classes* [27], *sensory uncertainty field* [91], and *visual constraints* [49, 40]. Nevertheless, the relation between the use of landmarks and the development of complete polynomial planners has not been explored up to now. In this work we show that the use of landmarks considerably simplifies the selection of the set of states that the robot may traverse during plan execution, as well as the synthesis of state-recognition functions. It mainly reduces planning to selecting motion commands to navigate from landmark to landmark until the goal is achieved.

For a given problem, our planner generates a plan as a collection of motion commands, each attached to a specific region of the configuration space. When the robot enters one of these regions, it executes the associated motion command. Such a plan is reminiscent of the *reaction plans* proposed in [16, 28, 86] as a way to deal with uncertainty at planning time. In particular, Schoppers [86] developed the notion of *universal plans* whose rules cover all possible situations that may occur at execution time, much like our plans do. The interesting point is that our planner takes polynomial time, while the generation of universal plans is often believed to be exponential. However, unlike the planners mentioned above, our planner addresses a restricted family of planning problems in which uncertainty is well-bounded.

In this work we propose a new approach to robot motion planning with uncertainty. We focus on a restricted class of problems, for which we present a sound, complete, and polynomial planning algorithm. Then, we engineer a robot and its workspace so that these assumptions are met, and we perform experiments. Because the planner returns guaranteed plans, experimental failures can be tracked directly to violations of one or more of the assumptions. Thus, it is possible to isolate those assumptions that are the hardest (or the costliest) to satisfy, and then try to relax or eliminate them. The effect of the relaxation or elimination of a certain assumption on the complexity of the planning algorithm is an important part of our study. Since we do not wish to leave the realm of polynomial algorithms, we call this approach *the search for the limits of polynomiality in motion planning with uncertainty.*

4.2 Statement of the problem

We consider a circular omnidirectional robot moving in a planar workspace amongst circular forbidden regions that represent the obstacles. The definition of the obstacles as forbidden regions implies that the robot is not allowed to be (or move) in contact with them. Such a workspace maps onto a planar configuration space, where a point robot navigates amongst C-obstacles represented by forbidden circular regions (the *obstacle disks*). The configuration space also contains a set of *landmark disks* representing the regions of the configuration space, from where certain identifiable features of the workspace (the *landmarks*) are perceivable by the sensors of the robot.

The number of landmark disks is finite and equal to ℓ. The number of obstacle disks is also finite and in $O(\ell)$. (Throughout this chapter the number ℓ is used to measure the size of the input problem.) Landmark disks may intersect with each other; obstacle disks may intersect with each other, too. However, intersections between landmark and obstacle disks are not allowed. Maximal connected sets of landmark disks are called *landmark areas*, and their number is equal to s $(\leq \ell)$. We assume that the robot has accurate knowledge of the configuration space model, i.e., it knows the exact location and size of the obstacle and landmark disks at planning time. This model does not change over time, i.e., landmarks and obstacles are stationary.

When the robot lies within a landmark disk, it can sense its configuration with perfect accuracy. Using its perfect sensing, the robot is able to navigate between any two points in the same landmark area without control error. This mode of navigation is called the *perfect-control mode*. When the robot lies outside all landmark disks, it has no sensing at all and it navigates using the *imperfect-control mode*.

A command in the perfect-control mode (called a *P-command*) is a sequence of regions $(R_1, R_2, \ldots, R_n)$ of the configuration space. The P-command simply prescribes that the robot should visit all these regions in order. We are not interested in the details of the execution of a P-command, as long as we know that it is guaranteed to succeed.

A motion command in the imperfect-control mode (an *I-command*) is described by a pair $(d, \mathcal{L}_T)$, where $d \in S^1$ is a direction in the plane, called the *commanded* *direction of motion*, and $\mathcal{L}_T$ is a set of landmark disks, called the *termination set* of the command. This command can be executed from anywhere in the configuration space outside the obstacle disks. The robot follows a path whose tangent at any point makes an angle with the direction d that is no greater than some prespecified angle θ called the *directional uncertainty*. The cone of angle 2θ whose axis points along d is the *directional uncertainty cone*. The robot stops as soon as it enters a landmark disk in $\mathcal{L}_T$. The robot has no sense of time, which means that the modulus of its velocity is irrelevant to the planning problem.

The initial position of the robot is known to be anywhere in a specified region $\mathcal{I}$ (the *initial region*) that consists of one or several disks called the *initial-region disks*. The number of initial-region disks is assumed constant. At planning time, we only know that the robot will be in the initial region when the execution of the plan starts; but the robot may not be there yet. Furthermore, we do not want to make any assumption about how it will move into the initial region; perhaps it will be transported there, or it will use another control mode not considered here. Thus, each initial-region disk may be disjoint from the landmark areas, or it may overlap with some of them, or it may be entirely contained in one of them. The robot must move into a given region $\mathcal{G}$, called the *goal region*, which is any subset, connected or not, of the workspace whose intersection with the landmark disks is easily computable.

The problem is to generate a *motion plan*, i.e., an algorithm made up of motion commands in the perfect- and imperfect-control mode, which guarantees that the robot will be in $\mathcal{G}$ when the execution of the plan terminates.

The problem we have described above is a *planning problem*, and the model we have adopted is suitable for the purposes and capabilities of a planner. The model has to be correct (i.e., produce results that are compatible with the operation of a real-world system), but it does not have to describe every single capability of the system. This abstraction is exactly the focus of our research, which attempts to develop models simple enough for efficient planning, but in enough detail so that they are useful. The execution module need not be constrained to the use of our planning model; it may use its own (usually more precise) model to optimize execution efficiency. Nevertheless, both models must be able to recognize the language used to formulate plans.

4.3 Outline of a planning algorithm

4.3.1 An example

Consider the example of Figure 4.1(a). The black disks represent obstacle disks and the grey disks represent landmark disks. There are four obstacle and seven landmark disks, but only five landmark areas. The initial region consists of a single disk marked with $\mathcal{I}$. The goal region consists of a single disk marked with $\mathcal{G}$.

The points of $\mathcal{G}$ that belong to some landmark disk are distinguishable from points outside the goal by means of the perfect position sensing provided by the landmark. Since $\mathcal{G}$ and landmark disk C have a common region, the robot can get from any point of C into the goal with the P-command $(\mathcal{G} \cap C)$. Similarly, the robot can get from D to C and from there to the goal with the P-command $(D \cap C, \mathcal{G} \cap C)$. But from no other landmark disk can the robot get into the goal using only the perfect-control mode. We say that C and D form the *extension set* $\mathcal{E}(\mathcal{G})$ of the goal. The remaining landmark disks are called the *intermediate goals*.

Since we cannot use the perfect-control mode to bring the robot into the extension set, we must use the imperfect-control mode. We are interested in guaranteed plans, so we need to find a commanded direction of motion for which, if the robot starts from any point in $\mathcal{I}$, it is guaranteed to get (and stop) in either C or D despite control error. This is equivalent to finding the preimage of the extension set for some commanded direction d (as defined in Chapter 3), and making sure that it contains $\mathcal{I}$. In this case, the computation of the preimage is not subject to the complications that arise from the interaction of goal reachability and recognizability. The reason is that landmark recognition occurs immediately after the robot enters a landmark region, and is independent of the way the landmark was achieved. Therefore, the preimage of a set of disks is equal to the backprojection of the set.

In Figure 4.2(a) the extension disks are white. Several preimages, computed for various values of the commanded direction d, are shown in the figure, but none of them completely includes the initial region $\mathcal{I}$. In fact, in this example there is no commanded direction d for which the preimage of C and D contains $\mathcal{I}$. Nevertheless,