Modular Reinforcement Learning: An Application to a Real Robot Task

Zsolt Kalmár[1], Csaba Szepesvári[2], and András Lőrincz[3]

[1] Dept. of Informatics JATE, Szeged, Aradi vrt. tere 1, Hungary, H-6720
{kalmar,szepes,lorincz}@iserv.iki.kfki.hu
[2] Research Group on Art. Int., JATE, Szeged, Aradi vrt. tere 1, Hungary, H-6720
[3] Dept. of Chemical Physics, Inst. of Isotopes, HAS, Budapest, P.O. Box 77, Hungary, H-1525

Abstract. The behaviour of reinforcement learning (RL) algorithms is best understood in completely observable, finite state- and action-space, discrete-time controlled Markov-chains. Robot-learning domains, on the other hand, are inherently infinite both in time and space, and moreover they are only partially observable. In this article we suggest a systematic design method whose motivation comes from the desire to transform the task-to-be-solved into a finite-state, discrete-time, "approximately" Markovian task, which is completely observable too. The key idea is to break up the problem into subtasks and design controllers for each of the subtasks. Then operating conditions are attached to the controllers (together the controllers and their operating conditions which are called modules) and possible additional features are designed to facilitate observability. A new discrete time-counter is introduced at the "module-level" that clicks only when a change in the value of one of the features is observed. The approach was tried out on a real-life robot. Several RL algorithms were compared and it was found that a model-based approach worked best. The learnt switching strategy performed equally well as a handcrafted version. Moreover, the learnt strategy seemed to exploit certain properties of the environment which could not have been seen in advance, which predicted the promising possibility that a learnt controller might overperform a handcrafted switching strategy in the future.

1 Introduction

Reinforcement learning (RL) is the process of learning the coordination of concurrent behaviours and their timing. A few years ago Markovian Decision Problems (MDPs) were proposed as the model for the analysis of RL [17] and since then a mathematically well-founded theory has been constructed for a large class of RL algorithms. These algorithms are based on modifications of the two basic dynamic-programming algorithms used to solve MDPs, namely the value- and policy-iteration algorithms [25, 5, 10, 23, 18]. The RL algorithms learn via experience, gradually building an estimate of the optimal value-function, which is known to encompass all the knowledge needed to behave in an optimal way according to a fixed criterion, usually the expected total discounted-cost criterion.

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The basic limitations of all of the early theoretical results of these algorithms was that they assumed finite state- and action-spaces, and discrete-time models in which the state information too was assumed to be available for measurement. In a real-life problem however, the state- and action-spaces are infinite, usually non-discrete, time is continuous and the system's state is not measurable (i.e. with the latter property the process is only partially observable as opposed to being completely observable). Recognizing the serious drawbacks of the simple theoretical case, researchers have begun looking at the more interesting yet theoretically more difficult cases (see e.g. [11, 16]). To date, however, no complete and theoretically sound solution has been found to deal with such involved problems. In fact the above-mentioned learning problem is indeed intractable owing to partial-observability. This result follows from a theorem of Littman's [9].

In this article an attempt is made to show that RL can be applied to learn real-life tasks when *a priori* knowledge is combined in some suitable way. The key to our proposed method lies in the use of high-level modules along with a specification of the operating conditions for the modules and other "features", to transform the task into a finite-state and action, completely-observable task. Of course, the design of the modules and features requires a fair amount of *a priori* knowledge, but this knowledge is usually readily available. In addition to this, there may be several possible ways of breaking up the task into smaller subtasks but it may be far from trivial to identify the best decomposition scheme. If all the possible decompositions are simultaneously available then RL can be used to find the best combination. In this paper we propose design principles and theoretical tools for the analysis of learning and demonstrate the success of this approach via *real-life* examples. A detailed comparison of several RL methods, such as Adaptive Dynamic Programming (ADP), Adaptive Real-Time Dynamic Programming (ARTDP) and Q-learning is provided, having been combined with different exploration strategies.

The article is organized in the following way. In the next section (Section 2) we introduce our proposed method and discuss the motivations behind it. In it the notion of "approximately" stationary MDPs is also introduced as a useful tool for the analysis of "module-level" learning. Then in Section 3 the outcome of certain experiments using a mobile robot are presented. The relationship of our work to that of others is contrasted in Section 4, then finally our conclusions and possible directions for further research are given in Section 5. Due to the lack of space some details were left out from this article, but these can be found in [8].

2 Module-based Reinforcement Learning

First of all we will briefly run through Markovian Decision Problems (MDPs), a value-function approximation-based RL algorithm to learn solutions for MDPs and their associated theory. Next the concept of recursive-features and time discretization based on these features are elaborated upon. This is then followed

by a sensible definition and principles of module-design together with a brief explanation of why the modular approach can prove successful in practice.

2.1 Markovian Decision Problems

RL is the process by which an agent improves its behaviour from observing its own interactions with the environment. One particularly well-studied RL scenario is that of a single agent minimizing the expected-discounted total cost in a discrete-time finite-state, finite-action environment, when the theory of MDPs can be used as the underlying mathematical model. A finite MDP is defined by the 4-tuple (S, A, p, c), where S is a finite set of states, A is a finite set of actions, p is a matrix of transition probabilities, and c is the so-called immediate cost-function. The ultimate target of learning is to identify an optimal policy. A policy is some function that tells the agent which set of actions should be chosen under which circumstances. A policy π is optimal under the *expected discounted total cost criterion* if, with respect to the space of all possible policies, π results in a minimum expected discounted total cost for all states. The optimal policy can be found by identifying the optimal value-function, defined recursively by

$$v^*(s) = \min_{a \in U(s)} \left(c(s,a) + \gamma \sum_{s'} p(s,a,s')\, v^*(s') \right)$$

for all states $s \in S$, where $c(s,a)$ is the immediate cost for taking action a from state s, γ is the discount factor, and $p(s,a,s')$ is the probability that state s' is reached from state s when action a is chosen. $U(s)$ is the set of admissible actions in state s. The policy which for each state selects the action that minimizes the right-hand-side of the above fixed-point equation constitutes an optimal policy. This yields the result that to identify an optimal policy it is sufficient just to find the optimal value-function v^*. The above simultaneous non-linear equations (non-linear because of the presence of the minimization operator), also known as the *Bellman equations* [3], can be solved by various dynamic programming methods such as the value- or policy-iteration methods [15].

RL algorithms are generalizations of the DP methods to the case when the transition probabilities and immediate costs are unknown. The class of RL algorithms of interest here can be viewed as variants of the value-iteration method: these algorithms gradually improve an estimate of the optimal value-function via learning from interactions with the environment. There are two possible ways to learn the optimal value-function. One is to estimate the model (i.e., the transition probabilities and immediate costs) while the other is to estimate the optimal action-values directly. The optimal action-value of an action a given a state s is defined as the total expected discounted cost of executing the action from the given state and proceeding in an optimal fashion afterwards:

$$Q^*(s,a) = c(s,a) + \gamma \sum_{s'} p(s,a,s') v^*(s'). \tag{1}$$

The general structure of value-function-approximation based RL algorithms is given in Table 1. In the RL algorithms various models are utilized along with

Initialization: Let $t = 0$, and initialize the utilized model (M_0) and the Q-function (Q_0)
Repeat Forever
1. Observe the next state s_{t+1} and reinforcement signal c_t .
2. Incorporate the new experience (s_t, a_t, s_{t+1}, c_t) into the model and into the estimate of the optimal Q-function: $(M_{t+1}, Q_{t+1}) = F_t(M_t, Q_t, (s_t, a_t, s_{t+1}, c_t))$.
3. Choose the next action to be executed based on (M_{t+1}, Q_{t+1}): $a_{t+1} = S_t(M_{t+1}, Q_{t+1}, s_{t+1})$ and execute the selected action.
4. $t := t + 1$.

Table 1. The structure of value-function-approximation based RL algorithms.

an update rule F_t and action-selection rule S_t.

In the case of the Adaptive Real-Time Dynamic Programming (ARTDP) algorithm the model consists (M_t) of the estimates of the transition probabilities and costs, the update-rule F_t being implemented e.g. as an averaging process. Instead of the optimal Q-function, the optimal value-function is estimated and stored to spare storage space, and the Q-values are then computed by replacing the true transition probabilities, costs and the optimal value-function in Equation 1 by their estimates. An update of the estimate for the optimal value-function is implemented by an asynchronous dynamic programming algorithm using an inner-loop in Step 2 of the algorithm.

Another popular RL algorithm is Q-learning, which does not employ a model but instead the Q-values are updated directly according to the iteration procedure [25]

$$Q_{t+1}(s_t, a_t) = (1 - \alpha_t(s_t, a_t))Q_t(s_t, a_t) +$$

$$\alpha_t(s_t, a_t)(c_t + \gamma \min_a Q_t(s_{t+1}, a)),$$

where $\alpha_t(s_t, a_t) \geq 0$, and satisfies the usual Robins-Monro type of conditions. For example, one might set $\alpha_t(s, a) = \frac{1}{n_t(s,a)}$ but often in practice $\alpha_t(s, a) =$ const is employed which while yielding increased adaptivities no longer ensures convergence.

Both algorithms mentioned previously are guaranteed to converge to the optimal value-/Q-function if each state-action pair is updated infinitely often. The action selection procedure S_t should be carefully chosen so that it fits the dynamics of the controlled process in a way that the condition is met. For example, the execution of random actions meets this "sufficient-exploration" condition when the MDP is ergodic. However, if on-line performance is important then

more sophisticated exploration is needed which, in addition to ensuring sufficient exploratory behaviour, exploits accumulated knowledge. For more details the interested reader is referred to [8].

2.2 Recursive Features and Feature-based Time-Discretization

In case of a real-life robot-learning task the dynamics cannot be formulated exactly as a *finite* MDP, nor is the state information available for measurement. This latter restriction is modelled by Partially-Observable MDPs (POMDPs) where (in the simplest case) one extends an MDP with an *observation function* h which maps the set of states S into a set X, called the observation set (which is usually non-countable, just like S). The defining assumption of a POMDP is that the full state s can be observed only through the observation function, i.e. only $h(s)$ is available as input and this information alone is usually insufficient for efficient control since h is usually a non-injection (i.e. h may map different states to the same observations). *Features* which mathematically are just well-designed observation functions, are known to be efficient in dealing with the problem of infinite state-spaces. Moreover, when their definitions are extended in a sensible way they become efficient in dealing with partial observability.

It is well known that POMDPs optimal policies can depend on their whole past histories. This leads us onto a mere generalization of features, such that the feature's values can depend on all past observations, i.e. mathematically a feature becomes an infinite sequence of mappings $(f^0, f^1, \ldots f^t, \ldots)$, with $f^t : (X \times A)^t \times X \to F$, where F and X are the feature- and observation-spaces. Since RL is supposed to work on the output of features, and RL requires finite spaces it means that F should be finite. Features that require infinite memory are clearly impossible to implement, so features used in practice should be restricted in such a way that they require a finite "memory". For example, besides stationary features which take the form $(f^0, f^0, \ldots, f^0, \ldots)$ (i.e. $f^t = f^0$ for all $t \geq 1$) and are called *sensor-based* features, *recursive* features (in control theory these are called *filters* or *state-estimators*) are those that can be implemented using a finite memory[1]. For example, in the case of a one-depth recursive feature the value at the t^{th} step is given by $f_t = R(x_t, a_{t-1}, f_{t-1})$, where $R : X \times A \times F \to F$ defines the recursion and $f_0 = f^0(x_0)$ for some function $f^0 : X \to F$.[2] Features whose values depend on the past-observations of a finite window form a special class of recursive filters. Later in the application-section some example features will be presented.

Instead of relying on a single feature, it is usually more convenient to define and employ a set of features, each of which indicates a certain event of interest. Note that a finite set of features can always be replaced by a single feature whose

[1] If the state space is infinite then not all sensor-based features can be realized in practice.

[2] If time is continuous then recursive features should be replaced by features that admit continuous-time dynamics. Details concerning continous time systems are given in [8].

output-space is the Cartesian product of the output spaces of the individual features and whose values can be computed componentwise by the individual single features. That is to say, the new feature's values are the 'concatenated' values of the individual features.

Since the feature-space is finite, a natural discretization of time can be obtained. The new time-counter clicks only when the feature value jumps in the feature-space. This makes it useful to think of such features as event-indicators which represent the actuality of certain conditions of interest. This interpretation gives us an idea of how to define features in such a way that the dynamics at the level of the new counter are simplified.

2.3 Modules

So far we have realized that the new "state-space" (the feature-space) and "time" can be made discrete. However, the action-space may still be infinite. Uniform discretization which lacks a priori knowledge would again be impractical in most of the cases (especially when the action space is unbounded), so we would rather consider an idea motivated by a technique which is often applied to solve large search-problems efficiently. The method in question divides the problem into smaller subproblems which are in turn divided into even smaller sub-problems, etc., then at the end routines are provided that deal with the resulting mini-problems. The solution of the entire problem is then obtained by working backwards: At every moment the mini-routine corresponding to the actual state of the search problem is applied. To put it in another way, the problem-solver defines a set of sub-goals, sub-sub-subgoals, etc. in such a way that if one of the sub-goals is satisfied then the resolution of the main-goal will be easy to achieve. In control-tasks the same decomposition can usually be done with respect to the main control objective without any difficulty. The routines that resolve the very low-level subgoals should be provided by closed-loop controllers which achieve the given subgoal under the set conditions (the conditions are usually given implicitly, e.g. the condition is fulfilled if the "predecessor" subgoals have already been achieved). The process of breaking up the problem into small subtasks can be repeated several times before the actual controllers are designed, so that the complexity of the individual controllers can be kept low. The controllers together with their operating conditions, which may serve as a basic set of features, will be called modules. In principle a consistent transfer of the AI-decomposition yields the result that the operating conditions of the situation are exclusive and cover every situation. However, such a solution would be very sensitive to perturbations and unmodelled dynamics.

A more robust solution can be obtained by extending the range of operating conditions and may mean that more than one controller can be applied at the same time. This calls for the introduction of a mechanism (the switching function) which determines which controller has to be activated if there are more than one available. More specifically, in our case a switching function $\mathcal{S}$ maps feature-vectors (which are composed of the concatenation of the operating conditions of the individual modules and some possible additional features) to the

indeces of modules and at any time the module with index $\mathcal{S}(f)$ where f id observed feature vector is activated. Of course only those modules can be activated whose operating conditions are satisfied. The operation of the whole mechanism is then the following. A controller remains active until the switching function switches to another controller. Since the switching function depends on the observed feature-values the controllers will certainly remain active till a change in the feature-vector is observed. We further allow the controllers to manipulate the own observation-process. In this the controllers may inhibit an observation from occurring and thus may hold up their activity for a while.

The goal of the design procedure is to set up the modules and additional features in such a way that there exists a proper switching controller $\mathcal{S} : F \to \{1, 2, \ldots, n\}$ which for any given history results in a closed-loop behaviour which fulfills the "goal" of control in time. It can be extremely hard to prove even the *existence* of such a valid switching controller. One approach is to use a so-called *accessibility decision problem* for this purpose which is a discrete graph with its node set being in our case the feature set F and the edges connect features which can be observed in succession. Then standard DP techniques can be used to decide the existence of a proper switching controller [8].

Of course, since the definitions of the modules and features depend on the designer, it is reasonable to assume that by clever design a satisfactory decomposition and controllers could be found even if only qualitative properties of the controlled object were known. RL could then be used for two purposes: either to find the best switching function assuming that at least two proper switching functions exist, or to decide empirically whether a valid switching controller exists at all. The first kind of application of RL arises as result of the desire to guarantee the existence of a proper switching function through the introduction of more modules and features than is minimally needed. But then good switching which exploits the capabilities of all the available modules could well become complicated to find manually.

If the accessibility decision problem were extendible with transition-probabilities to turn it to an MDP [3] then RL could be rightly applied to find the best switching function. For example if one uses a fixed (maybe stochastic) stationary switching policy and provided that the system dynamics can be formulated as an MDP then there is a theoretically well-founded way of introducing transition-probabilities (see [16]). Unfortunately, the resulting probabilities may well depend on the switching policy which can prevent the convergence of the RL algorithms. However, the following "stability" theorem shows that the difference of the cost of optimal policies corresponding to different transition probabilities is proportional to the extent the transition probabilities differ, so we may expect that a slight change in the transition probabilities does not result in completely different optimal switching policies and hence, as will be explained shortly after the theorem, we may expect RL to work properly, after all.

[3] Note that as the original control problem is deterministic it is not immediate when the introduction of probabilities can be justified. One idea is to refer to the ergodicity of the control problem.

Theorem 21 *Assume that two MDPs differ only in their transition-probability matrices, and let these two matrices be denoted by p_1 and p_2. Let the corresponding optimal cost-functions be v_1^* and v_2^*. Then*

$$||v_1^* - v_2^*|| \leq \gamma \frac{nC||p_1 - p_2||}{(1-\gamma)^2},$$

where $C = ||c||$ is the maximum of the immediate costs, $||\cdot||$ denotes the supremum-norm and n is the size of the state-space.

Proof. Let T_i be the optimal-cost operator corresponding to the transition-probability matrix p_i, i.e.

$$(T_i v)(s) = \min_{a \in U(x)} \left(c(s,a) + \gamma \sum_{s' \in X} p_i(s,a,s') v(s') \right),$$

$$v : S \to \Re,\ i = 1, 2.$$

Proceeding with standard fixpoint and contraction arguments (see e.g. [19]) we get that $\|v_1^* - v_2^*\| \leq \|T_1 v_1^* - T_1 v_2^*\| + \|T_1 v_2^* - T_2 v_2^*\|$ and since T_1 is a contraction with index γ, and the inequality $\|T_1 v - T_2 v\| \leq \gamma \|p_1 - p_2\| \sum_{y \in X} |v(y)|$ we obtain $\delta = \|v_1^* - v_2^*\| \leq \gamma\delta + \gamma \|p_1 - p_2\| |X| C/(1-\gamma)$, where $\|v_1^*\| \leq C/(1-\gamma)$ has been employed [15]. Rearranging the inequality in terms of δ then yields Theorem 21.

Motivated by the previous theorem we define ε-stationary MDPs as the quadruple (S, A, p, c), where S, A and c are as before but p, the transition probability matrix, may vary in time but with $||p_t - p^*|| \leq \varepsilon$ holding for all $t > 0$. Our expectations are that although the transitions cannot be modelled with a fixed transition probability matrix (i.e. stationary MDP), they can be modelled by an ε-stationary one even if the switching functions are arbitrarily varied and we conjecture that RL methods would then result in oscillating estimates of the optimal value-function, but with the oscillation being asymptotically proportional to ε. Note that ε-stationarity was clearly observed in our experiments which we will describe now.

3 Experiments

The validity of the proposed method was checked with actual experiments carried out using a Khepera-robot. The robot, the experimental setup, general specifications of the modules and the results are all presented in this section.

3.1 The Robot and its Environment

The mobile robot employed in the experiments is shown in Figure 1. It is a Khepera[4] robot equipped with eight IR-sensors, six in the front and two at the

[4] The Khepera was designed and built at Laboratory of Microcomputing, Swiss Federal Institute of Technology, Lausanne, Switzerland.

Fig. 1. The Khepera and the experimental environment. The task was to grasp a ball and hit the stick with it.

back, the IR-sensors measuring the proximity of objects in the range 0-5 cm. The robot has two wheels driven by two independent DC-motors and a gripper which has two degrees of freedom and is equipped with a resistivity sensor and an object-presence sensor. The vision turret is mounted on the top of the robot as shown. It is an image-sensor giving a linear-image of the horizontal view of the environment with a resolution of 64 pixels and 256 levels of grey. The horizontal viewing-angle is limited to about 36 degrees. This sensor is designed to detect objects in front of the robot situated at a distance spanning 5 to 50 cm. The image sensor has no tilt-angle, so the robot observes only those things whose height exceeds 5 cm.

The learning task was defined as follows: find a ball in an arena, bring it to one of the corners marked by a stick and hit the stick with the ball. The robot's environment is shown in Figure 1. The size of the arena was 50 cm x 50 cm with a black coloured floor and white coloured walls. The stick was black and 7 cm long, while three white-coloured balls with diameter 3.5 cm were scattered about in the arena. The task can be argued to have been biologically inspired because it can be considered as the abstraction of certain foraging tasks or a "basketball game". The environment is highly chaotic because the balls move in an unpredictable manner and so the outcome of certain actions is not completely predictable, e.g. a grasped ball may easily slip out from the gripper.

3.2 The Modules

Subtask decomposition Firstly, according to the principles laid down in Section 2, the task was decomposed into subtasks. The following subtasks were naturally: (T1) to find a ball, (T2) grasp it, (T3) bring it to the stick, and (T4) hit the stick with the grasped ball. Subtask (T3) was further broken into two subtasks, that of (T3.1) 'safe wandering' and (T3.2) 'go to the stick', since the robot cannot see the stick from every position and direction. Similarly, because of the robot's limited sensing capabilities, subtask (T1) was replaced by safe-wandering and subtask (T2) was refined to 'when an object nearby is sensed examine it and grasp it if it is a ball'. Notice that subtask 'safe wandering' is used for two purposes (to find a ball or the stick). The operating conditions of the corresponding controllers arose naturally as (T2) – an object should be nearby, (T3.2) – the stick should be detected, (T4) – the stick should be in front of the robot, and (T1,T3.1) – no condition. Since the behaviour of the robot must differ

before and after locating a ball, an additional feature indicating when a ball was held was supplied. As the robot's gripper is equipped with an 'object-presence' sensor the 'the ball is held" feature was easy to implement. If there had not been such a sensor then this feature still could have been implemented as a switching-feature: the value of the feature would be `'on'` if the robot used the grasping behaviour and hence not the hitting behaviour. An 'unstuck' subtask and corresponding controller were also included since the robot sometimes got stuck. Of course yet another feature is included for the detection of "goal-states". The corresponding feature indicates when the stick was hit by the ball. This feature's value is `'on'` iff the gripper is half-closed but the object presence sensor does not give a signal. Because of the implementation of the grasping module (the gripper was closed only after the grasping module was executed) this implementation of the "stick has been hit by the ball" feature was satisfactory for our purposes, although sometimes the ball slipped out from the gripper in which case the feature turned 'on' even though the robot did not actually reach the goal. Fortunately this situation did not happen too often and thus did not affect learning.

The resulting list of modules and features is shown in Table 2. The controllers work as intended, some fine details are discussed here (for more complete description see [8]). For example, the observation process was switched off until the controller of `Module 3`was working so as the complexity of the module-level decision problem is reduced. The dynamics of the controller associated with `Module 1` were based on the maximization of a function which depended on the proximity of objects and the speed of both motors[5]. If there were no obstacles near the robot this module made the robot go forward. This controller could thus serve as one for exploring the environment. `Module 2` was applicable only if the stick was in the viewing-angle of the robot, which could be detected in an unambiguous way because the only black thing that could get into the view of the robot was the stick. The range of allowed behaviour associated with this module was implemented as a proportional controller which drove the robot in such a way that the angle difference between the direction of motion and line of sight to the stick was reduced. The behaviour associated with `Module 3` was applicable only if there was an object next to the robot, which was defined as a function of the immediate values of IR-sensors. The associated behaviour was the following: the robot turned to a direction which brought it to point directly at the object, then the gripper was lowered. The "hit the stick" module (`Module 4`) lowers the gripper which under appropriate conditions result in that the ball jumps out of the gripper resulting in the goal state. `Module 5` was created to handle stuck situations. This module lets the robot go backward and is applicable if the robot has not been able to move the wheels into the desired position for a while. This condition is a typical time-window based feature.

Simple case-analysis shows that there is no switching controller that would reach the goal with complete certainty (in the worst-case, the robot could re-

[5] Modules are numbered by the identification number of their features.

FNo	'on'	Behaviour
1	always	explore while avoiding obstacles
2	if the stick is in the viewing angle	go to the stick
3	if an object is near	examine the object grasp it if it is a ball
4	if the stick is near	hit the stick
5	if the robot is stuck	go backward
6	if the ball is grasped	-
7	if the stick is hit with the ball	-

Table 2. Description of the features and the modules. 'FNo.' means 'Feature No.', in the column labelled by 'on' the conditions under which the respective feature's value is `'on'` are listed.

turn accidentally to state "10000000" from any state when the goal feature was `'off'`), so that an almost-sure switching strategy should always exist. On the other hand, it is clear that a switching strategy which eventually attains the target does indeed exist.

3.3 Details of learning

A dense cost-structure was applied: the cost of using each behaviour was one except when the goal was reached, whose cost was set to zero. Costs were discounted at a rate of $\gamma = 0.99$. Note that from time to time the robot by chance became stuck (the robot's 'stuck feature' was `'on'`), and the robot tried to execute a module which could not change the value of the feature-vector. This meant that the robot did not have a second option to try another module since by definition the robot could only make decisions if the feature-representation changed. As a result the robot could sometimes get stuck in a "perpetual" or so-called "jammed" state. To prevent this happening we built in an additional rule which was to stop and reinitialize the robot when it got stuck and could not unjam itself after 50 sensory measurements. A cost equivalent to the cost of never reaching the goal, i.e. a cost of $\frac{1}{1-\gamma}$ $(= 100)$ was then communicated to the robot, which mimicked in effect that such actions virtually last forever.

Experiments were fully automated and organized in trials. Each trial run lasted until the robot reached the goal or the number of decisions exceeded 150 (a number that was determined experimentally), or until the robot became jammed. The 'stick was hit' event was registered by checking the state of the gripper (see also the description of `Feature 7`).

During learning the Boltzmann-exploration strategy was employed where the temperature was reduced by $T_{t+1} = 0.999\, T_t$ uniformly for all states [2]. During the experiments the cumulative number of successful trials were measured and compared to the total number of trials done so far, together with the average number of decisions made in a trial.

3.4 Results

Two sets of experiments were conducted. The first set was performed to check the validity of the module based approach, while the second was carried out to compare different RL algorithms. In the first set the starting exploration parameter T_0 was set to 100 and the experiment lasted for 100 trials. These values were chosen in such a way that the robot could learn a good switching policy, the results of these experiments being shown in Figure 2. One might

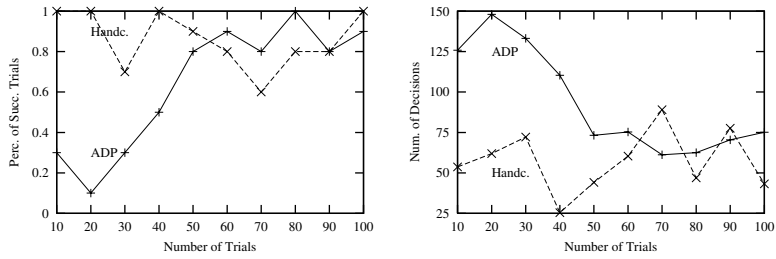

Fig. 2. Learning curves. In the first graph the percentage of successful trials out of ten are shown as a function of the number of trials. In the second graph the number of decisions taken by the robot and averaged over ten trials are both shown, as well as a function of the number of learning trials.

conclude from the left subgraph which shows the percentage of task completions in different stages of learning that the robot could solve the task after 50 trials fairly well. Late fluctuations were attributable to unsuccessful ball searches: as the robot could not see the balls if they were far from it, the robot had to explore to find one and the exploration sometimes took more than 150 decisions, yielding trials which were categorized as being failures. The evaluation of behaviour-coordination is also observed in the second subgraph, which shows the number of decisions per trial as a function of time. The reason for later fluctuations is again due to a ticklish ball search. The performance of a handcrafted switching policy is shown on the graphs as well. As can be seen the differences between the respective performances of the handcrafted and learnt switching functions are eventually negligible. In order to get a more precise evaluation of the differences the average number of steps to reach the goal were computed for both switchings over 300 trials, together with their standard deviations. The averages were 46.61 and 48.37 for the learnt and the handcrafted switching functions respectively, with nearly equal std-s of 34.78 and 34.82, respectively.

Theoretically, the total number of states is $2^7 = 128$, but as learning concentrates on feature-configurations that really occur this number transpires to be just 25 here. It was observed that the learnt policy was always consistent with a set of handcrafted rules, but in certain cases the learnt rules (which however can not be described here due to the lack of space) are more refined than their

handcrafted counterparts. For example, the robot learnt to exploit the fact that the arena was not completely level and as a result balls were biased towards the stick and as a result if the robot did not hold a ball but could see the stick it moved towards the stick.

In the rest of the experiments we compared two versions of ARTDP and three versions of *real-time Q-learning* (RTQL). The two variants of ARTDP which we call ADP, and "ARTDP", corresponding to the cases when in the inner loop of ARTDP the optimal value function associated with the actual estimated model (transition probabilities and imediate cost) is computed and when only the estimate of the value of the actual state is updated. Note that due to the small number of states and module-based time discretization even ADP could be run in real-time. But variants of RTQL differ in the choice of the learning-rate's time-dependence. RTQL-1 refers to the choice of the so-called *search-then-converge* method, where $\alpha_k(s,a) = \frac{50}{100+n_k(s,a)}$, $n_k(s,a)$ being the number of times the event $(s,a) = (s_t, a_t)$ happened before time k plus one (the parameters 50 and 100 were determined experimentally as being the best choices). In the other two cases (the corresponding algorithms were denoted by RTQL-2 and RTQL-3 respectively) constant learning rates (0.1 and 0.25, respectively) were utilized.

The online performances of the algorithms were measured as the cumulative number of unsuccessful trials, i.e. the regret. The regret R_t at time t is the difference between the performance of an optimal agent (robot) and that of the learning agent accumulated up to trial t, i.e. it is the price of learning up to time t. A comparison of the different algorithms with different exploration ratios is given in Table 3. All algorithms were examined with all the four different exploration parameters since the same exploration rate may well result in different regrets for different algorithms, as was also confirmed in the experiments.

Method	$T_0 = 100$	$T_0 = 50$	$T_0 = 25$	$T_0 = 0$
ADP	(61;19;6)	(52;20;4)	(45;12;5)	*(44;16;4)*
ARTDP	36	29	50	24
RTQL-1	53	69	47	66
RTQL-2	71	65	63	73
RTQL-3	(83;1;2)	(79;26:3)	(65;24;4)	*(61;14;2)*

Table 3. Regret. The table shows the number of unsuccessful trials among the first 100 trials. The entries with three number in them show cases when more than one experiment was run. In those entries the first number shows the average of the number of unsuccessful trials, the second is the standard deviation while the third is the number of experiments run.

First note that in order to evaluate statistically the differences observed for different exploration strategies much more experiments would be needed but running these experiments would require an enormous amount of time (approximately 40 days) and have not been performed yet. Thus we performed the

following procedure: Based on the first runs with every exploration-parameter and algorithm the algorithms that seemed to perform the best were selected (these were the ADP and the RTQL-3 algorithms) and some more experiments were carried out with these. The results of these experiments (15 more for the ADP and 7 more for the RTQL-3) indicated that the difference between the performances of the RTQL-3 and ADP is indeed significant at the level $p = 0.05$ (Student's t-test was applied for testing this).

We have also tested another exploration strategy which Thrun found the best among several undirected methods[6] [21]. These runs reinforced our previous findings that estimating a model (i.e. running ADP or ARTDP instead of Q-learning) could reduce the regret rate by as much as 40%.

4 Related Work

There are two main research-tracks that influenced our work. The first was the introduction of features in RL. Learning while using features were studied by Tsitsiklis and Van Roy to deal with large finite state spaces, and also to deal with infinite state spaces [22]. Issues of learning in partially observable environments have been discussed by Singh et al. [16].

The work of Connell and Mahadevan complements ours in that they set-up subtasks to be learned by RL and fixed the switching controller [13].

Asada et al. considered many aspects of mobile robot learning. They applied a vision-based state-estimation approach and defined "macro-actions" similar to our controllers [1]. In one of their papers they describe a goal-shooting problem in which a mobile robot shot a goal while avoiding another robot [24]. First, the robot learned two behaviours separately: the "shot" and "avoid" behaviours. Then the two behaviours were synthetised by a handcrafted rule and later this rule was refined via RL. The learnt action-values of the two behaviours were reused in the learning process while the combination of rules took place at the level of state variables.

Matarić considered a multi-robot learning task where each robot had the same set of behaviours and features [14]. Just as in our case, her goal was to learn a good switching function by RL. She considered the case when each of the robots learned separately and the ultimate goal was that learning should lead to a good collective behaviour, i.e. she concentrated mainly on the more involved multi-agent perspective of learning. In contrast to her work, we followed a more engineer-like approach when we suggested designing the modules based on well-articulated and simple principles and contrary to her findings it was discovered that RL can indeed work well at the modular level.

In the AI community there is an interesting approach to mobile robot control called Behaviour-Based Artificial Intelligence in which "competence" modules or behaviours have been proposed as the building blocks of "creatures" [12, 4]. The

[6] An exploration strategy is called undirected when the exploration does not depend on the number of visits to the state-action pairs.

decision-making procedure is, on the other hand, usually quite different from ours.

The technique proposed here was also motivated by our earlier experiences with a value-estimation based algorithm given in the form of "activation spreading" [20]. In this work activation spread out along the edges of a dynamically varying graph, where the nodes represented state transitions called triplets. Later the algorithm was extended so that useful subgoals could be found by learning [6, 7]. In the future we plan to extend the present module-based learning system with this kind of generalization capability. Such an extension may in turn allow the learning of a hierarchical organization of modules.

5 Summary and Conclusions

In this article module-based reinforcement learning was proposed to solve the coordination of multiple "behaviours" or controllers. The extended features served as the basis of time- and space discretization as well as the operating conditions of the modules. The construction principles of the modules were: decompose the problem into subtasks; for each subtask design controllers and the controllers' operating conditions; check if the problem could be solved by the controllers under the operating and observability conditions, add additional features or modules if necessary, set-up the reinforcement function and learn a switching function from experience.

The idea behind our approach was that a partially observable decision problem could be usually transformed into a completely observable one if appropriate features (filters) and controllers were employed. Of course some *a priori* knowledge of the task and robot is always required to find those features and controllers. It was argued that RL could work well even if the resulting problem was only an almost stationary Markovian. The design principles were applied to a real-life robot learning problem and several RL-algorithms were compared in practice. We found that estimating the model and solving the optimality equation at each step (which could be done owing to the economic, feature-based time-discretization) yielded the best results. The robot learned the task after 700 decisions, which usually took less than 15 minutes in real-time. We conjecture that using a rough initial model good initial solutions could be computed off-line which could further decrease the time required to learn the optimal solution for the task.

The main difference between earlier works and our approach here is that we have established principles for the design modules and found that our subsequent design and simple RL worked spendidly. Plans for future research include extending the method via module learning and also the theoretical investigation of almost stationary Markovian decision problems using the techniques developed in [10].

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