# Feature-discovering Approximate Value Iteration Methods

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**Abstract.** Sets of features in Markov decision processes can play a critical role in approximately representing value and in abstracting the state space. Selection of features is crucial to the success of a system and is most often conducted by a human. We study the problem of automatically selecting problem features, and propose and evaluate a simple approach reducing the problem of selecting a new feature to standard classification learning. We learn a classifier that predicts the sign of the Bellman error over a training set of states. By iteratively adding new classifiers as features with this method, training between iterations with approximate value iteration, we find a Tetris feature set that outperforms randomly constructed features significantly, and obtains a score of about three-tenths of the highest score obtained by using a carefully hand-constructed feature set. We also show that features learned with this method outperform those learned with the previous method of Patrascu et al. [4] on the same SysAdmin domain used for evaluation there.

## 1 Introduction

Decision-theoretic planning and reinforcement-learning methods facing astronomically large state spaces typically rely on approximately represented value functions (see, e.g., [2, 7]). Many such approximate representations rely on an appropriate set of problem features; for example, by taking a weighted combination of the feature values as the value function [1]. Human engineering of the problem features used has repeatedly proven critical to the success of the resulting system. For example, in [8, 7], TD-gammon exploits human-constructed problem-specific features to achieve a playing strength that can compete with world-class players.

A Markov decision process (MDP) is a formal model of a single agent facing a sequence of action choices from a pre-defined action space, and living within a pre-defined state space. After each action choice is made, a state transition within the state space occurs according to a pre-defined, stochastic action-transition model. The agent receives reward after each action choice according to the state visited (and possibly the action chosen), and has the objective of accumulating as much reward as possible (possibly favoring reward received sooner, using discounting, or averaging over time, or requiring that the reward be received by a finite horizon).

MDP solution techniques often critically rely on finding a good value function; this is a mapping from states to real numbers with the intent that desirable states receive

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high values. Informally, a good value function should respect the action transitions in that good states will either have large immediate rewards or have actions available that lead to other good states; this property is formalized in *Bellman equations* that define the optimal value function (see below). The degree to which a given value function fails to respect action transitions in this way, formalized below, is referred to as the *Bellman error* of that value function.

Unfortunately, virtually every interesting MDP problem has an extremely large state space, preventing direct table-based representation of the value function. As a result, abstraction, approximation, and problem reformulation play a critical role in successfully representing and finding good value functions. A typical approach is to find useful *features*, which also map the state space to real numbers, and take the value function at each state to be a weighted combination of the features at that state. Here, good values for the weights can often be found using machine learning techniques involving search and gradient descent. In this paper, we address the problem of finding good features automatically: in most previous work the features are simply selected by the human designer. Learning features for this purpose can be regarded as learning an abstraction of the MDP state space: differences between states with the same feature values have been abstracted away.

Here, we study the problem of automatically selecting problem features for use in approximately representing value in Markov decision processes. We focus our initial work on this problem on binary features, i.e., mappings from state to Boolean values. We view each such feature as a set of states, those states where the feature is true.

We propose and evaluate a simple, greedy approach to finding new binary features for a linear-combination value estimate. Our heuristic approach assumes an initial "base" value estimate described by a linear approximation where the weights have already been tuned to minimize Bellman error. We attempt to reduce the Bellman error magnitude of this value estimate further by learning a new feature that is true in statespace regions of positive statewise Bellman error and correspondingly false in regions of negative statewise Bellman error, or vice versa. The learning problem generated is a standard supervised classification problem, and for this work we address this problem using the decision-tree learner C4.5 [5].

One view of this approach is that we are conducting approximate value iteration with an added mechanism for extending the available feature set. Given an initial feature set, imagine a sufficient period of approximate value iteration (or any similar weight adjustment method) to achieve convergence of the approximation to a value function  $\tilde{V}$ . We can think of the approximate value iteration process as "stuck", in that it can represent  $\tilde{V}$  but not the Bellman update of  $\tilde{V}$ . (Of course, this assumes that the inductive updates being performed would find  $\tilde{V}$  if they could represent it, which is only heuristically true.) We are then trying to induce features to enable representation of the Bellman update of  $\tilde{V}$ , so that the approximate value iteration process can continue to reduce Bellman error, with the larger feature space.

If the learner succeeds in capturing features that describe the statespace regions of positive and negative Bellman error, we can guarantee that adding these features makes available weight assignments closer to the Bellman update of the base value estimate.

Our practical method retrains the weights including the new feature(s), using approximate value iteration (AVI), and then repeats the process of selecting a new feature.

We have found surprisingly little previous work on selecting features automatically in MDPs. Patrascu et al. [4] give a linear programming technique for selecting new features. In their work, the primary technique for selecting weights is approximate linear programming (ALP). It is observed in [4] that ALP "only minimizes  $L_1$  error"; perhaps for this reason, that work proposes to construct features aimed to minimize the  $L_1$  error of the resulting approximation. It is stated there that there is a "hope that this leads to a reduction in Bellman error as a side effect."

Our technique works instead directly to reduce the Bellman error magnitude of the resulting approximation by trying to identify regions that contribute to large statewise Bellman error magnitude. Because the Bellman error at any given state is easily computable in many domains of interest, unlike the  $L_1$  error, we are able to convert the problem of minimizing Bellman error magnitude to a supervised classification problem.

Patrascu et al. [4] do not mention any reduction to supervised-classification learning, nor is it clear how to construct such a reduction from the approach they describe, because the computation of the  $L_1$  error for particular states requires knowledge of the optimal value of those states. Thus, although decision trees are suggested as a possible representation for candidate features in that work, there is no suggestion that these trees be acquired by a supervised classification method like C4.5 and no discussion of how to overcome the need to know the optimal value if the  $L_1$  error is to be used for supervision.

For another view of the contrast between our method and that previous method, consider that we seek a feature, via supervised classification, that corresponds to the region of states that have significantly positive (or alternatively, negative) Bellman error. There is no such declarative characterization of the desired feature in the Patrascu et al. work; rather, the region sought is that region that results in the best improvement in  $L_1$  error after retraining via ALP (or less expensive-to-compute approximations of this "error after retraining"). This does not represent a reduction to classification and is a substantially different approach, using  $L_1$  error after retraining (or a less expensive stand-in) as a scoring function.

Both approaches are reasonable, of course. We show below that our technique empirically outperforms this previous work on the planning domain used in their evaluation. Specifically, we require fewer new features to achieve the same Bellman error, and can achieve a lower overall Bellman error given enough features.

There is related previous work on function approximation in which new features are automatically added during supervised learning of real-valued functions [10]. It would be reasonable to consider, in comparison to our work here, plugging in such a function approximator at the learning step in approximate value iteration—this would result in an overall method similar in spirit to what we design directly here. We have not done an empirical comparison along these lines at this time.

We also evaluate our technique in the computer-game domain of Tetris. Starting from a constant value function based on only the uniformly true feature, our technique can add features automatically to produce performance that is significantly better than a randomly constructed feature set, and is at about three-tenths of the performance of a carefully hand-constructed feature set.

In what follows, we first provide technical background on Markov decision processes and value-function approximation, then describe our technique for inducing new features to reduce approximation error, and finally present empirical results on two domains showing improvement over the state of the art, before concluding.

## 2 Technical Background

#### 2.1 Markov Decision Processes

We define here our terminology for Markov decision processes. For a more thorough discussion of Markov decision processes, see [2] and [7]. A Markov decision process (MDP) D is a tuple (S,A,R,T) where state space S is a finite set of states, action space A is a finite set of actions,  $R:S\times A\times S\to \mathbb{R}$  is the reward function, and  $T:S\times A\to \mathcal{P}(S)$  is the transition probability function that maps (state, action) pairs to probability distributions over S.  $R(s_1,a,s_2)$  represents how much immediate reward is obtained by taking action a from state  $s_1$  and ending up in state  $s_2$ .  $T(s_1,a,s_2)$  represents the probability of ending up in state  $s_2$  if the action a is taken from state  $s_1$ .

A policy  $\pi$  for an MDP is a mapping  $\pi:S\to A$ . Given policy  $\pi$ , the value function  $V^\pi(s)$  gives the expected discounted reward obtained starting from state s and selecting action  $\pi(s)$  at each state encountered. Rewards after the first time step are discounted by a factor  $\gamma$  where  $0\le\gamma<1$ . A Bellman equation relates  $V^\pi$  at any state s and successor states s':

$$V^{\pi}(s) = \sum_{s' \in S} T(s, \pi(s), s') [R(s, \pi(s), s') + \gamma V^{\pi}(s')].$$

There is at least one optimal policy  $\pi^*$  for which  $V^{\pi^*}(s)$ , abbreviated  $V^*(s)$ , is no less than  $V^{\pi}(s)$  at every state s, for any other policy  $\pi$ . Another Bellman equation governs  $V^*$ :

$$V^*(s) = \max_{a \in A} \sum_{s' \in S} T(s, a, s') [R(s, a, s') + \gamma V^*(s')].$$

From any value function V, we can compute a policy  $\operatorname{Greedy}(V)$  that selects, at any state s, the greedy look-ahead action  $\operatorname{arg} \max_{a \in A} \sum_{s' \in S} T(s, a, s')[R(s, a, s') + \gamma V(s')]$ . The policy  $\operatorname{Greedy}(V^*)$  is an optimal policy. Value iteration iterates the operation  $V'(s) = \max_{a \in A} \sum_{s' \in S} T(s, a, s')[R(s, a, s') + \gamma V(s')]$ , computing V' from V, producing a sequence of value functions converging to  $V^*$ , regardless of the initial V used.

We define the statewise Bellman error B(V,s) for a value function V at a state s to be V'(s)-V(s). We will be inducing new features based on the sign of the statewise Bellman error. The sup-norm distance of a value function V from the optimal value function  $V^*$  can be bounded using the Bellman error magnitude, which is defined as  $\max_{s\in S}|B(V,s)|$  (e.g., see [12]).

Linear Approximation of Value Functions. We assume that the states of the MDP have structure. In particular, we assume a state is a vector of basic properties with Boolean, integer, or real values, and that the state space is the set of all such vectors. We call these basic properties *state attributes*. This factored form for states is essential to enable compact representation of approximate value functions.

A common solution to the problem of representing value functions (e.g., value iteration) in very large, structured state spaces is to approximate the value V(s) with a linear combination of features extracted from s, i.e., as  $\tilde{V}(s) = \sum_{i=0}^p w_i f_i(s)$ , where  $w_i$  is a real-valued weight for the ith feature  $f_i(s)$ . Our goal is to find features  $f_i$  (each mapping states to boolean values) and weights  $w_i$  so that  $\tilde{V}$  closely approximates  $V^*$ .

Many methods have been proposed to select weights  $w_i$  for linear approximations [6, 11]. Here, we use a trajectory-based approximate value iteration (AVI) approach. Other training methods can be substituted and this choice is orthogonal to our main purpose.

The AVI method we deploy constructs a fixed-length sequence of value functions  $V^1, V^2, \ldots, V^T$ , and returns the last one. Each value function  $V^\beta$  is defined by weight values  $w_0^\beta, w_1^\beta, \ldots, w_p^\beta$  as  $V^\beta(s) = \sum_{i=0}^p w_i^\beta f_i(s)$ . Value function  $V^{\beta+1}$  is constructed from  $V^\beta$  by drawing a training set of trajectories under the policy Greedy( $V^\beta$ ) and updating the weights according to this training set as follows.

Let  $s_1, s_2, \ldots, s_n$  be a sequence of training states, which we generate by drawing a set of trajectories under the current greedy policy. A training set for weight adjustment is defined as  $\{(s_j, V'(s_j)) \mid 1 \leq j \leq n\}$ . We adjust weights in iterations of batch training. In the l'th batch training iteration, the weight update for the i'th weight in the training set is  $w_{i,l+1} = w_{i,l} + \frac{1}{n} \sum_j \alpha f_i(s_j)(V'(s_j) - \sum_{i=0}^p w_{i,l} f_i(s_j))$ . Here,  $\alpha$  is the learning rate parameter. We take the initial weights  $w_{i,1}$  to be  $w_i^\beta$ . After  $\kappa$  iterations we set  $w_i^{\beta+1} = w_{i,\kappa+1}$ .

#### 2.2 Decision Tree Classification

A detailed discussion of classification using decision trees can be found in [3]. A decision tree is a binary tree with internal nodes labelled by state attributes (and, in our case, learned features), and leaves labelled with classes (in our case, either zero or one). A path through the tree from the root to a leaf with label l identifies a partial assignment to the state attributes—each state consistent with that partial assignment is viewed as labelled l by the tree. We learn decision trees from training sets of labelled states using the well known C4.5 algorithm [5]. This algorithm induces a tree greedily matching the training data from the root down. We use C4.5 to induce new features—the key to our algorithm is how we construct suitable training sets for C4.5 so that the induced features are useful in reducing Bellman error.

<sup>&</sup>lt;sup>1</sup> The source of this set is a parameter of the algorithm, and it could for example be drawn by sampling initial states from some state distribution and then simulating  $\pi$  to some horizon from each initial state.

## **3 Feature Construction for MDPs**

We propose a simple method for constructing new features given a current set of features and an MDP for which we desire an approximation of  $V^*$ . We first use AVI, as described above, to select heuristically best weights to approximate  $V^*$  with  $\tilde{V}$  based on the current feature set. We then use the sign of the statewise Bellman error at each state as an indication of whether the state is undervalued or overvalued by the current approximation. If we can identify a collection of undervalued states (ideally, all such states) as a new feature, then assigning an appropriate positive weight to that feature should reduce the Bellman error magnitude. The same effect should be achieved by identifying overvalued states with a new feature and assigning a negative weight. We note that the domains of interest are generally too large for statespace enumeration, so we will need classification learning to generalize the notions of overvalued and undervalued across the statespace from training sets of sample states. Also, to avoid blurring the concepts of overvalued and undervalued with each other, we discard states with statewise Bellman error near zero from either training set.

More formally, we draw a training set of states  $\Sigma$  from which we will select training subsets  $\Sigma_+$  and  $\Sigma_-$  for learning new features. The training set  $\Sigma$  can either be drawn uniformly at random from the state space, or drawn by collecting all states in sample trajectories starting at uniformly random start states under a policy of interest (typically Greedy( $\tilde{V}$ )). If using trajectories, each trajectory must be terminated at some horizon. The horizon and the size of  $\Sigma$  are parameters of our algorithm.

For each state s in  $\Sigma$ , we compute the statewise Bellman error  $B(\tilde{V},s)$ . We then discard from  $\Sigma$  those states s with statewise Bellman error near zero, i.e., those states for which  $|B(\tilde{V},s)|<\delta$  for a non-negative real-valued parameter  $\delta$ , and then divide the remaining states into sets  $\Sigma_+$  and  $\Sigma_-$  according to the sign of  $B(\tilde{V},s)$ . So,  $\Sigma_+$  is the set  $\{s|B(\tilde{V},s)\geq\delta\}$  and  $\Sigma_-$  is the set  $\{s|B(\tilde{V},s)\leq-\delta\}$ .

We note that computing statewise Bellman error exactly can involve a summation over the entire state space, whereas our fundamental motivations require avoiding such summations. In many MDP problems of interest, the transition matrix T is sparse in a way that set of states reachable in one step with non-zero probability is small, for any current state. In such problems, statewise Bellman error can be computed effectively using an appropriate representation of T. More generally, when T is not sparse in this manner, the expectation can be effectively approximately evaluated by sampling next states according to the distribution represented by T.

We then use  $\Sigma_+$  as the positive examples and  $\Sigma_-$  as the negative examples for a supervised classification algorithm; in our case, C4.5 is used. The hypothesis space for classification is built from the primitive attributes defining the state space; in our case, we use decision trees over these attributes. We can also interchange the roles of  $\Sigma_+$  and  $\Sigma_-$ , using the latter as positive examples. In our experiments, we do this interchanging for every other feature constructed.

The concept resulting from supervised learning is then treated as a new feature for our linear approximation architecture, with an initial weight of zero. The process can then be repeated, of course, resulting in larger and larger feature sets, and, hopefully, smaller and smaller Bellman error magnitude.

To conclude our description of our algorithm, we discuss setting the parameter  $\delta$  dynamically, once in each iteration of feature construction. Rather than directly specify  $\delta$ , we specify  $\delta$  in terms of the standard deviation  $\sigma$  of the statewise Bellman error over the same distribution used in selecting states for the training set  $\Sigma$ . The value of  $\sigma$  is easily estimated by sampling the training distribution and computing the Bellman error. We then set  $\delta$ , at each iteration, to be a fixed multiple  $\eta$  of  $\sigma$ . This approach removes  $\delta$  as a parameter of the algorithm, replacing it with the parameter  $\eta$ . This dynamic selection of  $\delta$  allows adaptation to the decrease in Bellman error magnitude over the run of the algorithm.

## 4 Experiments

In this section, we present some experimental results for our feature construction algorithm. We use two domains in the experiments. The first domain is an  $8 \times 8$  game of Tetris (Tetris). The second domain is a computer network optimization problem called SysAdmin, which we use primarily in order to compare to the closest previous related work; that work [4] used SysAdmin as a testing domain. Both the state attributes and the learned features in the experiments are binary features.

Tetris. For the Tetris domain, we start with 71 state attributes; 64 attributes which represent if the 64 squares are occupied or not, and 7 attributes which represent which of the 7 pieces is currently being dropped. We select training sets for feature construction by drawing trajectories from an initial state with an empty board and collecting 600,000 states on these trajectories as  $\Sigma$ . The training sets for AVI are selected by drawing 100 trajectories from an initial state with an empty board and allowing each trajectory to extend to the end of the game. We draw the trajectories using the Greedy( $\tilde{V}$ ) policy. The discount factor  $\gamma$  is 0.9 for this experiment, and the parameter  $\eta$  is set to 0.3. In addition,  $\kappa$  is fixed at 100 and  $\alpha$  at 0.01. AVI is stopped (appearing to have converged) after 1,200 training sets are drawn; at that point, a new feature is learned.

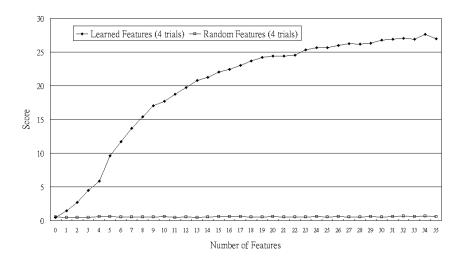
The results are shown in Figure 1. The score is determined by the average number of lines erased during a sequence of games. The performance of the learned features are evaluated by the 2,000-game average score for  $\operatorname{Greedy}(\tilde{V})$  using the weights learned by AVI. Figure 1 displays the average of such evaluations over 4 separate trials of feature learning. In addition, we also show in Figure 1 the result of using sets of randomly generated features; such features are generated following the same procedure described in the previous section, but label examples in the training set  $\Sigma$  randomly instead of deciding the labeling by statewise Bellman error. Value functions constructed from randomly generated features perform poorly, and do not show improvement as the number of features used increases. Thus, our use of statewise Bellman error to label the training examples plays an important role in the performance of our feature construction algorithm.

We also tested AVI on human-constructed features in this domain. The features we used in this case were provided by Bertsekas in [2]. These features are useful features as considered by a human, and according to [2] they were selected after some testing. We tested the performance of the weights learned after each AVI iteration by running 2,000

games and taking the average score. The maximum 2,000-game average performance was 92.9, which was achieved after 22,634 iterations of AVI. This performance was substantially better than the best performance our learned feature set exhibited, which was 27.6 (using 34 learned binary features).

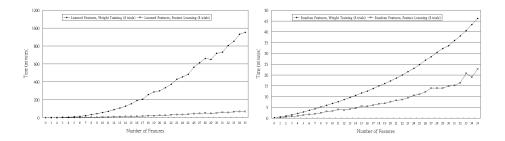
We note that the human-selected features are all integer-valued, apparently giving the human set a clear advantage over our binary features (especially per feature). Clearly one approach for further improvement in feature learning is to design a feature-learning approach that can produce integer-valued features.

We also study the the runtime behavior for our algorithm in this domain. We show in Figure 2 the execution time for generating the features and adjusting the weights. How long it takes a feature to be learned depends on how many training examples are collected and how many features exist. Since the number of training examples does not grow as we learn additional features, the time required to learn a new feature is eventually dwarfed by growing time required to train the weights for the growing feature set.



**Fig. 1.** Score (average number of lines erased in 2,000 games) plot for the learned features and randomly generated features in the  $8 \times 8$  Tetris domain. For reference, the maximum score for the human-selected feature set from [2] was 92.9.

SysAdmin. For the SysAdmin domain, two different kinds of topologies are tested: 3-legs and cycle. There are 10 nodes in each topology. We follow the settings used in [4] for testing this domain. The target of learning in this domain is to keep as many machines operational as possible, since the number of operating machines directly affects the reward for each step. Since there are only 10 nodes, the on/off status of each node is used as a basic feature, which means there are a total of 1024 states. We simply use



**Fig. 2.** Number of minutes required to generate the n'th feature, and to train the weights after the n'th feature is added to the feature set.

all states as the training set for feature construction. To enable direct comparison to the previous work in [4], we use Bellman error magnitude to measure the performance of the feature construction algorithm here.

For the experiments that use the whole state space as a training set, the plot of average Bellman error for 10 separate trials over the number of features learned is shown in Figure 3. We used  $\gamma$  equal to 0.95,  $\eta$  equal to 1,  $\alpha$  equal to 0.1, and  $\kappa$  equal to 100. In this experiment there are 50 trajectories drawn in each AVI training set, each drawn from a random initial state, and using trajectory length 2. AVI was stopped, appearing to have converged, after 1,000 iterations.

Also included in Figure 3 are the results from [4]. We select the best result they show (from various algorithmic approaches) from the 3-legs and cycle domains shown in their paper (their "d-o-s" setting for the cycle domain and their "d-x-n setting" for the 3-legs domain).

Compared to the results in [4], also shown here, our feature construction algorithm achieves a lower Bellman error magnitude in these domains for the same number of features, throughout, and a lower converged Bellman error magnitude when new features stop improving that measure. This is another encouraging result for this proposed feature construction algorithm.

# 5 Conclusions and Future Work

From the experiments, the results show that our feature construction algorithm can generate features that show significantly better performance in  $8\times 8$  Tetris than randomly constructed features, and can produce features that outperform the features produced by the algorithms in [4] for the SysAdmin domain. However, our algorithm cannot learn a feature set for  $8\times 8$  Tetris that competes well with the human-constructed feature set provided in [2].

Our technique depends critically on the generalization ability of the classification learner to cope with large state spaces. The features generated by the feature-construction algorithm are currently represented as decision trees. Although the experiments showed that these features are useful in some problems, they are still hard to interpret. One

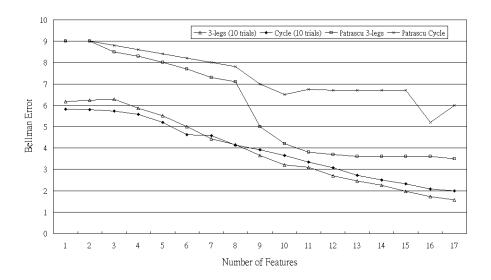


Fig. 3. Bellman error for SysAdmin domain (10 nodes).

goal for designing a good feature-construction algorithm is to be able to produce features that are understandable by humans. Furthermore, decision tree learning might not be adequate to find good generalizations in complex domains. We observed that the human-constructed features in [2] can be represented compactly using relational languages. Some of them, e.g. the number of "holes" in Tetris, are quite awkward to represent using the decision tree structure in this paper. One way we are considering for improving our algorithm is to use a relational classification or function-approximation algorithm combined with an expressive knowledge representation instead of using C4.5 with decision trees.

We note that the performance of machine-learned feature set appears to converge, with little added benefit per new feature, at a point where the policy corresponding to the learned value function is far short of optimal. This suggests a lack of state-space exploration during the feature learning stage. Another direction we are considering for improving our algorithm is to develop new exploration strategies for generating the training sets of states for feature learning.

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