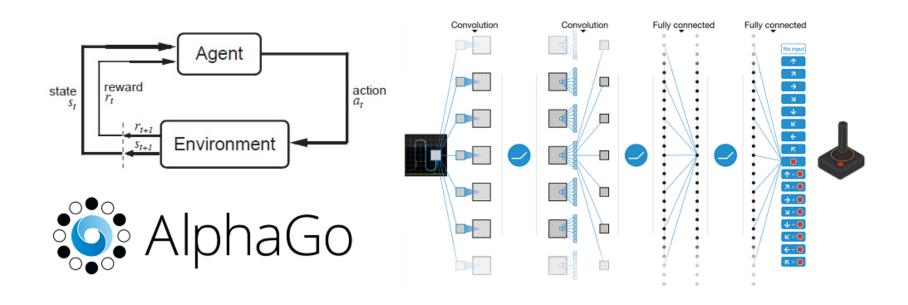
DSCI 552, Machine Learning for Data Science

University of Southern California

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Lesson 13 Reinforcement Learning



Overview

- Supervised Learning: Immediate feedback (labels provided for every input).
- Unsupervised Learning: No feedback (no labels provided).
- Reinforcement Learning: Delayed scalar feedback (a number called reward).

Overview

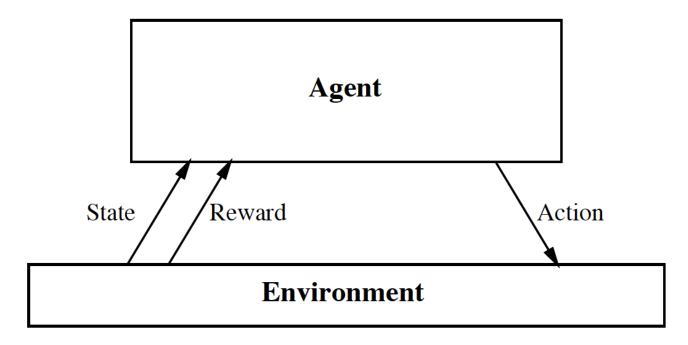
- RL deals with agents that must sense & act upon their environment.
- •This combines classical agent-based Al and machine learning techniques.

It is a very comprehensive problem setting.

Overview

- Examples:
 - A robot cleaning my room and recharging its battery
 - Robot-soccer
 - How to invest in shares
 - Modeling the economy through rational agents
 - Learning how to fly a helicopter
 - Scheduling planes to their destinations
 - and so on

The Big Picture



$$s_0 \stackrel{a_0}{\longrightarrow} s_1 \stackrel{a_1}{\longrightarrow} s_2 \stackrel{a_2}{\longrightarrow} \dots$$

Your action influences the state of the world which determines its reward

Complications

- The outcome of your actions may be uncertain
- You may not be able to perfectly sense the state of the world
- The reward may be stochastic.
- Reward is delayed (i.e. finding food in a maze)

Complications

- You may have no clue (model) of how rewards are being paid off.
- The world may change while you try to learn it
- How much time do you need to explore uncharted territory before you exploit what you have learned?

The Task

• To learn an optimal *policy* that maps states of the world to actions of the agent.

I.e., if this patch of room is dirty, I clean it. If my battery is empty, I recharge it.

$$\pi: \mathcal{S} \to \mathcal{A}$$

What is it that the agent tries to optimize?
 Answer: the total future discounted reward:

The Task

What is it that the agent tries to optimize?
 Answer: the total future discounted reward:

$$V^{\pi}(s_{t}) = r_{t} + \gamma r_{t+1} + \gamma^{2} r_{t+2} + \dots$$

$$= \sum_{i=0}^{\infty} \gamma^{i} r_{t+i} \qquad 0 \leq \gamma < 1$$

Note: immediate reward is worth more than future reward.

What would happen to mouse in a maze with gamma = 0 ?

Value Function

- Let's say we have access to the optimal value function that computes the total future discounted reward $V^*(s)$
- What would be the optimal policy $\pi^*(s)$?
- Answer: we choose the action that maximizes:

$$\pi^*(s) = \operatorname{argmax} \left[r(s,a) + \gamma V^*(\delta(s,a)) \right]$$

Value Function

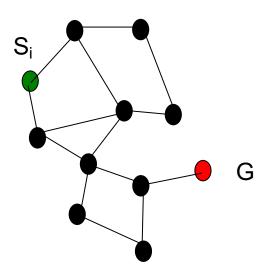
 We assume that we know what the reward will be if we perform action "a" in state "s":

 We also assume we know what the next state of the world will be if we perform action "a" in state "s":

$$S_{t+1} = \delta(S_t, a)$$

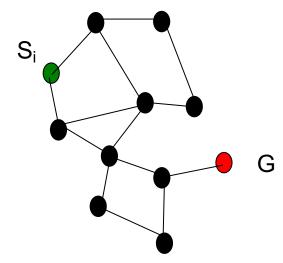
Example I

- Consider some complicated graph, and we would like to find the shortest path from a node S_i to a goal node G.
- Traversing an edge will cost you "length edge" dollars.



Example I

- The value function encodes the total remaining distance to the goal node from any node s, i.e. V(s) = "1 / distance" to goal from s.
- If you know V(s), the problem is trivial. You simply choose the node that has highest V(s).

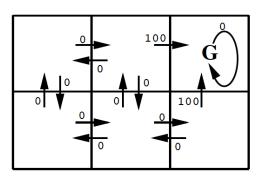


Example II

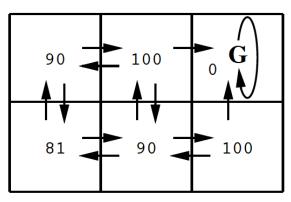
- •A simple deterministic world.
- Each grid square represents a distinct state, each arrow a distinct action.
- •The immediate reward function, r(s, a) gives reward 100 for actions entering the goal state G, and zero otherwise. Values of $V^*(s)$ follow from r(s, a), and the discount factor $\gamma = 0.9$.
- An optimal policy is also shown.

Example II

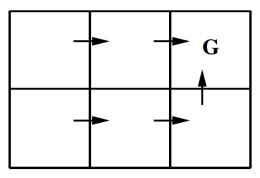
Find your way to the goal.



r(s, a) (immediate reward) values



$$V^*(s)$$
 values



One optimal policy

$$V^{\pi}(\mathbf{S}_{t}) = \mathbf{r}_{t} + \gamma \mathbf{r}_{t+1} + \gamma^{2} \mathbf{r}_{t+2} + \dots$$

$$= \sum_{i=0}^{\infty} \gamma^{i} \mathbf{r}_{t+i} \qquad 0 \le \gamma < 1$$

One approach to RL is then to try to estimate V*(s).
 Bellman Equation:

$$V^*(s) \leftarrow \max_{a} \left[r(s,a) + \gamma V^*(\delta(s,a)) \right]$$

- However, this approach requires you to know r(s,a) and $\delta(s,a)$.
- This is unrealistic in many real problems. What is the reward if a robot is exploring mars and decides to take a right turn?

•Fortunately we can circumvent this problem by exploring and experiencing how the world reacts to our actions. We need to *learn* $r \& \delta$.

• We want a function that directly learns good state-action pairs, i.e. what action should I take in this state. We call this Q(s,a).

- Let us define the evaluation function Q(s, a) so that its value is the maximum discounted cumulative reward that can be achieved starting from state s and applying action a as the first action.
- In other words, the value of Q is the reward received immediately upon executing action a from state s, plus the value (discounted by γ) of following the optimal policy thereafter.

$$Q(s,a) = r(s,a) + \gamma V^*(\delta(s,a))$$

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• Why is this rewrite important? Because it shows that if the agent learns the Q function instead of the V^* function, it will be able to select optimal actions even when it has no knowledge of the functions r and δ .

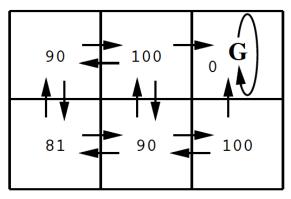
• It need only consider each available action a in its current state s and choose the action that maximizes Q(s, a).

$$\pi^*(s) = \underset{a}{\operatorname{argmax}} Q(s,a)$$
 $V^*(s) = \underset{a}{\operatorname{max}} Q(s,a)$

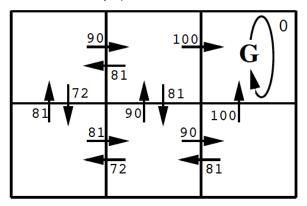
Example

- •To illustrate, the figure in the next slide shows the Q values for every state and action in the simple grid world.
- •The Q value for each state-action transition equals the *r* value for this transition plus the *V** value for the resulting state discounted by *γ*.
- The optimal policy shown in the figure corresponds to selecting actions with maximal Q values.

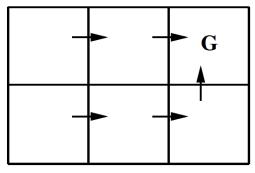
Example II



 $V^*(s)$ values



Q(s,a) values



One optimal policy

Check that
$$\pi^{*}(s) = \operatorname*{argmax}_{a} Q(s,a)$$

$$V^{*}(s) = \operatorname*{max}_{a} Q(s,a)$$

- Learning the Q function corresponds to learning the optimal policy. How can Q be learned?
- The key problem is finding a reliable way to estimate training values for Q, given only a sequence of immediate rewards r spread out over time.

- This can be accomplished through iterative approximation.
- To see how, notice the close relationship between Q and V*,

$$V^*(s) = \max_{a'} Q(s, a')$$

which allows rewriting the Q function as:

$$Q(s,a) = r(s,a) + \gamma V^*(\delta(s,a))$$

$$= r(s,a) + \gamma \max_{a'} Q(\delta(s,a),a')$$

This still depends on r(s,a) and $\delta(s,a)$; however,

$$Q(s,a) = r(s,a) + \gamma V^*(\delta(s,a))$$

$$= r(s,a) + \gamma \max_{a'} Q(\delta(s,a),a')$$

this recursive definition of Q provides the basis for algorithms that iteratively approximate Q.

Q refers to the learner's estimate, or hypothesis, of the actual Q function.

$$Q(s,a) = r(s,a) + \gamma V^*(\delta(s,a))$$

$$= r(s,a) + \gamma \max_{a'} Q(\delta(s,a),a')$$

- Imagine the robot is exploring its environment, trying new actions as it goes.
- At every step it receives some reward "r", and it observes the environment change into a new state s' for action a.
- •How can we use these observations, (s,a,s',r) to learn a model?

- •The learner represents its hypothesis \hat{Q} by a large table with a separate entry for each state-action pair.
- •The table entry for the pair (s, a) stores the value for $\hat{Q}(s,a)$, learner's current hypothesis about the actual but unknown value Q(s,a).
- •The table can be initially filled with random values (though it is easier to understand the algorithm if one assumes initial values of zero).

•The agent repeatedly observes its current state s, chooses some action a, executes this action, then observes the resulting reward r = r(s, a) and the new state $s' = \delta(s, a)$.

• It then updates the table entry for Q(s,a) following each such transition, according to the rule:

$$\hat{Q}(s,a) \leftarrow r + \gamma \max_{a'} \hat{Q}(s',a')$$
 $s' = s_{t+1}$

$$\hat{Q}(s,a) \leftarrow r + \gamma \max_{a'} \hat{Q}(s',a')$$
 $s' = s_{t+1}$

- This equation continually estimates Q at state s consistent with an estimate of Q at state s', one step in the future: temporal difference (TD) learning.
- Note that s' is closer to goal, and hence more "reliable", but still an estimate itself.

Q-Learning Summary

Q learning algorithm

For each s, a initialize the table entry $\hat{Q}(s, a)$ to zero.

Observe the current state s

Do forever:

- Select an action a and execute it
- Receive immediate reward r
- Observe the new state s'
- Update the table entry for $\hat{Q}(s, a)$ as follows:

$$\hat{Q}(s,a) \leftarrow r + \gamma \max_{a'} \hat{Q}(s',a')$$

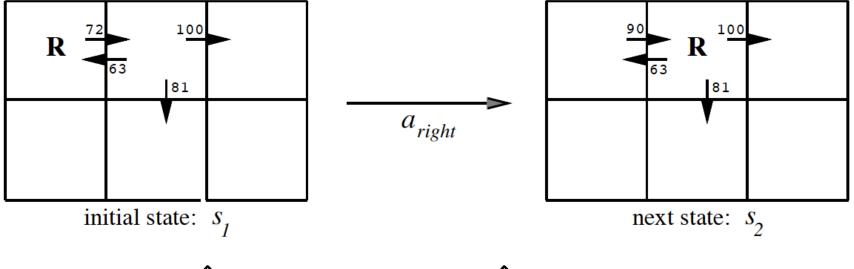
• $s \leftarrow s'$

$$\hat{Q}(s,a) \leftarrow r + \gamma \max_{a'} \hat{Q}(s',a')$$

$$s' = s_{t+1}$$

- We do an update after each state-action pair. I.e., we are learning online!
- We are learning useful things about explored state-action pairs. These are typically most useful because they are likely to be encountered again.
- Under suitable conditions, these updates can actually be proved to converge to the real answer.

Example: Q-Learning



$$\hat{Q}(s_1, a_{right}) \leftarrow r + \gamma_{\max_{a'}} \hat{Q}(s_2, a')$$

$$\leftarrow 0 + 0.9 \max\{66, 81, 100\}$$

$$\leftarrow 90$$

Q-learning propagates Q-estimates 1-step backwards

- •The algorithm does not specify how actions are chosen by the agent.
- •One strategy would be in state s to select the action a that maximizes $\hat{Q}(s,a)$, thereby exploiting its current approximation \hat{Q} .

- •Using this strategy the agent runs the risk that it will overcommit to actions that are found during early training to have high Q values, while failing to explore other actions that have even higher values.
- •It is common in Q learning to use a probabilistic approach to selecting actions.

•Actions with higher Q values are assigned higher probabilities, but every action is assigned a nonzero probability. One way to assign such probabilities is

 $P(a_i | s) = \frac{e^{\hat{Q}(s,a_i)/T}}{\sum_{j} e^{\hat{Q}(s,a_j)/T}}$

•where $P(a_i | s)$ is the probability of selecting action a_i , given that the agent is in state s, and where T > 0 is a constant that determines how strongly the selection favors actions with high Q values.

•Hence it is good to try new things so now and then, e.g. If *T* large lots of exploring, if *T* small, exploit current policy. One can decrease *T* over time to first explore, and then converge and exploit.

•For example T = c/k + d where k is iteration of the

algorithm

$$P(a_{i} | s) = \frac{e^{\hat{Q}(s, a_{i})/T}}{\sum_{j} e^{\hat{Q}(s, a_{j})/T}}$$

•Decreasing *T* over time is sometimes called simulated annealing, which is inspired by annealing process in metals. *T* is sometimes called the *Temperature*.

Improvements

- One can trade-off memory and computation by cashing (s,s',r) for observed transitions. After a while, as Q(s',a') has changed, you can "replay" the update: $\hat{Q}(s,a) \leftarrow r + \gamma \max_{a'} \hat{Q}(s',a')$
- One can actively search for state-action pairs for which Q(s,a) is expected to change a lot (prioritized sweeping).

- •Q learning algorithm learns by iteratively reducing the discrepancy between Q value estimates for adjacent state
- •Q learning is a special case of *temporal* difference algorithms that learn by reducing discrepancies between estimates made by the agent at different times.

- •The raining rule we studied reduces the difference between the estimated Q values of a state and its immediate successor
- •However, we could design an algorithm that reduces discrepancies between this state and more distant descendants or ancestors.

- •Recall that our Q learning training rule calculates a training value for $Q(s_t, a_t)$ in terms of the values for $Q(s_{t+1}, a_{t+1})$ where s_{t+1} is the result of applying action a_t to the state s_t .
- Let $Q^{(1)}(s_t, a_t)$ denote the training value calculated by this one-step lookahead:

$$Q^{(1)}(s_t, a_t) \equiv r_t + \gamma \max_{a} \hat{Q}(s_{t+1}, a)$$

•One alternative way to compute a training value for $Q(s_t, a_t)$ is to base it on the observed rewards for two steps

$$Q^{(2)}(s_t, a_t) \equiv r_t + \gamma r_{t+1} + \gamma^2 \max_{a} \hat{Q}(s_{t+2}, a)$$

• or, in general, for *n* steps

$$Q^{(n)}(s_t, a_t) \equiv r_t + \gamma r_{t+1} + \dots + \gamma^{(n-1)} r_{t+n-1} + \gamma^n \max_{a} \hat{Q}(s_{t+n}, a)$$

- •A general method for blending these alternative training estimates, called $TD(\lambda)$.
- •The idea is to use a constant $0 \le \lambda \le 1$ to combine the estimates obtained from various lookahead distances in the following fashion

$$Q^{\lambda}(s_t, a_t) \equiv (1 - \lambda) \left[Q^{(1)}(s_t, a_t) + \lambda Q^{(2)}(s_t, a_t) + \lambda^2 Q^{(3)}(s_t, a_t) + \cdots \right]$$

•An equivalent recursive definition for Q^{λ} is

$$Q^{\lambda}(s_t, a_t) = r_t + \gamma [(1 - \lambda) \max_{a} \hat{Q}(s_t, a_t) + \lambda Q^{\lambda}(s_{t+1}, a_{t+1})]$$

$$Q^{\lambda}(s_{t}, a_{t}) = r_{t} + \gamma [(1 - \lambda) \max_{a} \hat{Q}(s_{t}, a_{t}) + \lambda Q^{\lambda}(s_{t+1}, a_{t+1})]$$

- •If $\lambda = 0$ we have our original training estimate $Q^{(1)}$, which considers only one-step discrepancies in the Q estimates.
- •As λ is increased, the algorithm places increasing emphasis on discrepancies based on more distant lookaheads.

$$Q^{\lambda}(s_t, a_t) = r_t + \gamma [(1 - \lambda) \max_{a} \hat{Q}(s_t, a_t) + \lambda Q^{\lambda}(s_{t+1}, a_{t+1})]$$

- •At the extreme value $\lambda = 1$, only the observed r_{t+i} values are considered, with no contribution from the current Q estimate.
- •The motivation for the TD(λ) method is that in some settings training will be more efficient if more distant lookaheads are considered.

Extensions

• To deal with stochastic environments, we need to maximize *expected* future discounted reward:

$$Q(s,a) = E[r(s,a)] + \gamma \sum_{s'} P(s'|s,a) \max_{a'} Q(s',a')$$

 Often the state space is too large to deal with all states and adopt a table-lookup approach. In this case we need to learn a function:

$$Q(s,a) \approx f_{\theta}(s,a)$$

Extensions

$$Q(s,a) \approx f_{\theta}(s,a)$$

- Neural network with back-propagation have been quite successful.
- For instance, TD-Gammon is a back-gammon program that plays at expert level.
- •state-space very large, trained by playing against itself, uses NN to approximate value function, uses $TD(\lambda)$ for learning.

More on Function Approximation

For instance: linear function:

$$Q(s,a) \approx f_{\theta}(s,a) = \sum_{k=1}^{K} \theta_{k}^{a} \Phi_{k}(s)$$

The features Φ are fixed measurements of the state (e.g. # stones on the board).

We only learn the parameters theta.

 Update rule: (start in state s, take action a, observe reward r and end up in state s')

$$\theta_k^a \leftarrow \theta_k^a + \alpha \left(r + \gamma \max_{a'} \hat{Q}(s', a') - \hat{Q}(s, a) \right) \Phi_k(s)$$

$$change in Q$$

Conclusion

 Reinforcement learning addresses a very broad and relevant question:

How can we learn to survive in our environment?

 We have looked at Q-learning, which simply learns from experience.

No model of the world is needed.

Conclusion

•We made simplifying assumptions: e.g. state of the world only depends on last state and action. This is the *Markov* assumption. The model is called a *Markov Decision Process (MDP)*.

Conclusion

- We assumed deterministic dynamics, reward function, but the world really is stochastic.
- There are many extensions to speed up learning.
- There have been many successful real world applications.