Autonomous and Adaptive Systems

Monte Carlo Methods

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Monte Carlo Methods

- Monte Carlo methods are ways of solving the reinforcement learning problems based on averaging the sample returns.
- ▶ We focus on episodic tasks.
 - Only on the completion of episodes are values estimates and policies changed.
- Monte Carlo methods sample and average returns for each state-action pair much like the multi-armed bandits methods and average rewards for each action.
 - However, we now consider *multiple states*; and
 - The return after taking an action in one state depends on the actions taken in later states in the same episodes.
 - ▶ In other words: this is the *full* reinforcement learning problem.

Monte Carlo Methods

- If you have the full knowledge of the MDP you can compute the value functions (see Bellman equation, dynamic programming).
- We assume that we do not have full knowledge of the underlying MDP.
 - This is the case in general because the underlying dynamics and characteristics of the system are unknown (e.g., robot exploration) or because the system is too complex (e.g., games).
- Since we do not have full knowledge, we need to *learn* the values functions.

Monte Carlo Methods

- ► We consider three problems:
 - The prediction problem: the estimation of v_{π} and q_{π} for a fixed policy π .
 - The *policy improvement problem*: the estimation of v_{π} and q_{π} while trying at the same time to improve the policy π .
 - The *control problem*: the estimation of an optimal policy π_* .

Monte Carlo Prediction

▶ Goal: learning the state-value function for a given policy.

- Recall: the value of a state is the expected return (expected cumulative future discounted reward) from that state.
- Obvious/simple solution: average the returns observed after visiting that state. As more returns are observed, the average should converge to the expected value.

Monte Carlo Prediction

- More formally: we want to estimate $v_{\pi}(s)$, the value of a state *s* under policy π , given a set of episodes obtained by following π and passing through *s*.
- Each occurrence of a state s in an episode is called a *visit* to s.
- A state s can be visited multiple times.
- The first time a state *s* is visited in an episode is called the first visit to *s*.
- The *first-visit Monte Carlo method* estimates $v_{\pi}(s)$ as the average of the returns following first visits to *s*, whereas the *every-visit Monte Carlo method* averages the returns following all the visits to *s*.

First-visit Monte Carlo Prediction

Input: a policy π to be evaluated

Initialise:

 $V(s) \in \mathbb{R}$ arbitrarily for all $s \in \mathcal{S}$

 $Returns(s) \leftarrow$ empty list for all $s \in \mathcal{S}$

Loop forever (for each episode):

Generate an episode following $\pi: S_0, A_0, R_1, S_1, A_1, R_2, ..., S_{T-1}, A_{T-1}, R_T$

$G \leftarrow 0$

Loop for each step of episode t = T - 1, T - 2, ..., 0:

 $G \leftarrow \gamma G + R_{t+1}$

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If S_t does not appear in S_0, S_1, \ldots, S_{t-1}:
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Append G to Returns(S_t)
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V(S_t) \leftarrow average(Returns(S_t))
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Multi-visit Monte Carlo Prediction

- Every-visit Monte Carlo would be the same except without the check for S_t having occurred earlier in the episode.
- Both first-visit Monte Carlo and every-visit Monte Carlo converge to $v_{\pi}(s)$ as the number of visits (or first visits) to *s* goes to infinity.

Monte Carlo Estimation of Action Values

- The estimation of a state value makes sense when you have a model of the system.
 - ▶ With a model, state values alone are sufficient to determine a policy.
 - Without a model, it is necessary to estimate the value of each action in order for the value to be useful in suggesting a policy.
- The policy evaluation problem for action values is to estimate $q_{\pi}(s, a)$, the expected return when starting in state *s*, taking action *a* and then following policy π .

• Remember we assume that the policy π is fixed.

Monte Carlo Estimation of Action Values

- The methods for the Monte Carlo estimation of action values are essentially the same as those presented for state value, but now we talk about visits to the state-action pair instead of to a state.
- A state-action pair s, a is said to be visited in an episode if the state s is visited and the action a is taken in it.
- The first-visit method Monte Carlo method averages the returns following the first time in each episode that the state was visited and the action was selected.

Maintaining Exploration

- ▶ But there is a problem: many state-action pairs might never be visited.
- If π is a deterministic policy, then in following π, we will observe returns only for one of the actions of each state.
 - ▶ No return to average -> no improvement with experience.
 - ▶ We cannot compare alternatives, because no alternatives are explored.
- To compare alternatives, we need to estimate the value of all the actions from each state, not only the one we currently prefer (according to our policy).
- ▶ How can we address this problem?

Maintaining Exploration

▶ This is the general problem of *maintaining exploration*.

- One way to do this is to have the episodes starting in a state-action pair and that every pair has non-zero probability of being selected as start.
- This ensures that asymptotically (infinite number of episodes), all the state-action pairs will be visited an infinite number of times.

On-policy and Off-policy Exploration

- The method described above is useful, but it cannot applied in general.
 - Think about the case for example where you have exploration with the environment. You cannot start by "jumping" to a certain state-action pair at the beginning.
- The most common alternative is to ensure that all the state-action pairs are explored anyway.
- We need to explore these states, not following the current policy (for example with a stochastic policy). In other words, the exploration is not performed on-policy as done until now in this lecture, but off-policy.
 - Various methods are possible, such as off-policy prediction via Importance Sampling (not covered in this module - see Sutton and Barto Chapter 5.5).

Policy Improvement

- We will focus now back on on-policy exploration, i.e., the policy is used to make decisions and to explore the various states.
- Until now we assumed that the policy was fixed. However, the policy itself can be *improved* while learning the value functions and, potentially, we might have to aim at obtaining an optimal policy.
- Methods used for improving a policy in order to reach the optimal policy are usually referred to as *Monte Carlo control*.

Policy Improvement

Remember until now we assumed that the policy was fixed.

▶ Given that policy, we estimate the value functions.

- Now we consider how to improve the policy starting from an onpolicy method.
 - The policy that we use to make decisions is that we are trying to improve. We do not use a separate policy to explore the stateaction pairs (that would be an off-policy method).
- Policy improvement is done by making the policy greedy with respect to the current value functions.

Policy Improvement

▶ In this case we have an action-value function.

▶ No model is needed to construct the greedy policy.

For any action-value function q, the corresponding greedy policy is the one that, for each $s \in S$, deterministically chooses an action with maximal action-value:

$$\pi(s) \leftarrow \arg\max_{a} Q(s, a)$$

We have formal results that ensures that this process of policy improvement leads to optimal policy (usually referred to Monte Carlo control).

Monte Carlo Control without Exploring Starts

- An alternative method for Monte Carlo control (i.e., to achieve the optimal policy) is to use an *ε*-greedy approach.
- Most of the time the action that has the maximal estimated action value is chosen but with a probability *e* an action is chosen at random.
 - This means that all non-greedy actions are given the minimal probability of selection $\frac{\epsilon}{|\mathscr{A}(s)|}$ is given to all the non-greedy actions.
- Following these mechanism, it can be guaranteed that we will have an improvement of the policy (policy improvement theorem).

Monte Carlo Control with Exploring Starts

Initialise:

 $\pi(s) \in \mathscr{A}(s)$ arbitrarily for all $s \in \mathscr{S}$

 $Q(s, a) \in \mathbb{R}$ arbitrarily for all $s \in \mathcal{S}, a \in \mathcal{A}(s)$

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Returns(s, a) \leftarrow empty \text{ list for all } s \in \mathcal{S}, a \in \mathcal{A}(s)
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Loop forever (for each episode):

Choose $S_0 \in \mathcal{S}, A_0 \in \mathscr{A}(S_0)$ randomly

Generate an episode from S_0 , A_0 following π : S_0 , A_0 , R_1 , S_1 , A_1 , R_2 , ..., S_{T-1} , A_{T-1} , R_T

$$G \leftarrow 0$$

```
Loop for each step of episode t = T - 1, T - 2, ..., 0:
```

```
G \leftarrow \gamma G + R_{t+1}
```

```
If S_t does not appear S_0, A_0, R_1, A_1, R_2, \dots, S_{T-1}, A_{T-1}:
Append G to Returns(S_t, A_t)
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Q(S_t, A_t) \leftarrow average(Returns(S_t, A_t))
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$$\pi(S_t) \leftarrow \arg\max_a Q(S_t, a)$$

References

Chapter 5 of Sutton and Barto. Introduction to Reinforcement Learning. Second Edition. MIT Press. 2018.