



Reinforcement Learning (3)

Maxime Berar

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Monte-Carlo Methods

MC learning through complete episodes, model-free

Episode : a state-reward sequence



Complete return of a state in an episode

$$G_t(e) \stackrel{.}{=} = R_{t+1} + \gamma R_{t+2} + \dots + \gamma^{T-t-1} R_T$$

is used to estimate v_{π}

$$v_{\pi}(s) = rac{1}{n_{ ext{episodes}}^{s}} \sum_{e=1}^{n_{ ext{episodes}}^{s}} G_{t}(e)$$

▶ 2 versions : first-visit *or* every-visits

Algorithm 1 First-visit MC prediction, for estimating $V pprox v_\pi$

- 1: Initialize:
- 2: $\pi \leftarrow$ policy to be evaluated
- 3: $V \leftarrow$ an arbitrary state-value function
- 4: $Returns(s) \leftarrow$ an empty list, $\forall s \in \mathcal{S}$
- 5: while forever do
- 6: Generate an episode using π
- 7: **for** each state *s* appearing in the episode **do**
- 8: $G \leftarrow$ the return that follows the first occurrence of s
- 9: Append G to Returns(s)
- 10: $V(s) \leftarrow \text{average}(Returns(s))$
- 11: end for
- 12: end while
 - The first-visit MC method estimates $v_{\pi}(s)$ as the average of the returns following first visits to s. It dates back to the 1940s.

Visits

- The first-visit MC ...
- ► The every-visit MC method averages the returns following all visits to s. Every-visit MC extends more naturally to function approximation and eligibility traces (as presented later).

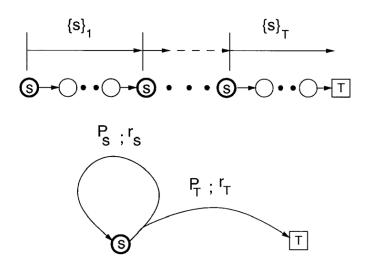
Both algorithms converge to $v_{\pi}(s)$ as the number of (first) visits to s goes to infinity. In the case of first-visit MC,

$$\hat{v}_{\pi}(s) = rac{1}{n_{ ext{ep}}^s} \sum_{e=1}^{n_{ ext{ep}}^s} G_t(e)$$

each return $G_t(e)$ is an iid estimate of $v_\pi(s)$ with finite variance. The sequence of averages of these estimates converges to their expected value. Each average is itself an unbiased estimate, and the standard deviation of its error falls as $1/n_{\rm ep}^s$, where $n_{\rm ep}^s$ is the number of returns averaged. The samples obtained by the every-visit strategy are dependent so they're not i.i.d. samples compared to first-visit strategy.

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Reduction to a Two-State Abstracted Markov Chain



Singh and Sutton, Reinforcement Learning with Replacing Eligibility Traces, Machine Learning 22 (1996)

First Visit MC

Let $\{x\}$ stand for the paired random sequence $(\{s\}, \{r\})$. The first-visit MC estimate for V(s) after one trial, $\{x\}$, is

$$V_1^F(s) = f(\lbrace x \rbrace) = r_{s_1} + r_{s_2} + \cdots + r_{s_k} + r_T,$$

where k is the random number of revisits to state s, r_{s_i} is the sum of the individual rewards in the sequence $\{r\}_i$, and r_T is the random total reward received after the last visit to state s. For all i, $\mathbb{E}[r_s] = R_s$. The first-visit MC estimate of V(s) after n trials, $\{x\}^1, \{x\}^2, \dots, \{x\}^n$, is

$$V_n^F(s) = f(\{x\}^1, \dots, \{x\}^n) = \frac{\sum_{i=1}^n f((\{x\}^i))}{n}$$

Every visit MC

$$V_1^{E}(s) = t(\lbrace x \rbrace) = \frac{t_{num}(\lbrace x \rbrace)}{k+1} = \frac{r_{s_1} + 2r_{s_2} + \cdots + kr_{s_k} + (k+1)r_T}{k+1},$$

where k is the random number of revisits to state s in the sequence $\{x\}$. Every visit to state s effectively starts another trial. The every-visit MC estimate after n trials, $\{x\}^1, \{x\}^2, \cdots, \{x\}^n$, is

$$V_n^E(s) = t(\{x\}^1, \cdots, \{x\}^n) = \sum_{i=1}^n \frac{t_{num,i}(\{x\})}{k_i + 1}$$

Some results on First Visit

From Bellman's equation

$$V(s) = \frac{P_s}{P_T} R_S + R_T$$

Expectation of $V_1^F(s)$

$$\mathbb{E}[f(\{x\})] = \sum_{\{x\}} P(\{x\})f(x) = \sum_{k} P(k)\mathbb{E}_{r}[f(\{x\})|k],$$

$$= \sum_{k} P_{T}P_{s}^{k}(kR_{S} + R_{T})$$

$$= P_{T}\left(\frac{P_{s}}{1 - P^{2}}R_{s} + \frac{1}{1 - P_{s}}R_{T}\right) = V(s)$$

Every-Visit MC is Biased

Expectation of $V_1^E(s)$

$$\mathbb{E}[t(\{x\})] = \sum_{\{x\}} P(\{x\})t(x)$$

$$= \sum_{k} P(k)\mathbb{E}_{r}[t(\{x\})|k],$$

$$= \sum_{k} P_{T}P_{s}^{k}(\frac{R_{s} + 2R_{s} + \dots + kR_{s} + (k+1)R_{T}}{k+1})$$

$$= \sum_{k} P_{T}P_{s}^{k}(\frac{k}{2}R_{s} + R_{T})$$

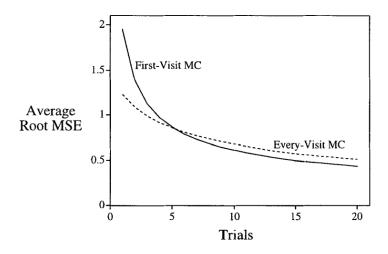
$$= \frac{P_{s}}{2P_{T}}R_{s} + R_{T} \neq V(s)$$

After *n* trials

$$\mathbb{E}\left[V_n^E(s)\right] = \frac{nP_s}{(n+1)P_T}R_S + R_T$$

every-visit MC algorithm is unbiased in the limit, $n \to \infty$

Comparison



Monte Carlo Estimation of Action Values

- without a model, state values alone are not sufficient.
- useful to estimate action values (the values of state-action pairs)

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 thereafter following policy π .

- ▶ The MC methods for this are essentially the same as just presented for state values, visits to a state-action pair rather than to a state
- first-visit MC vs every-visit MC

Pb! complication is that many state-action pairs may never be visited

- for policy evaluation to work for action values, we must assure continual exploration.
- One way to do this is by specifying that the episodes start in a state-action pair, and that every pair has a nonzero probability of being selected as the start.

the assumption of exploring starts.

alternative : consider only policies that are stochastic with a nonzero probability of selecting all actions in each state.

Monte Carlo Control

How Monte Carlo estimation can be used to approximate optimal policies?

- Generalized Policy Iteration : one maintains both an approximate policy and an approximate value function.
- The value function is repeatedly altered to more closely approximate the value function for the current policy, and the policy is repeatedly improved with respect to the current value function.
- These 2 improvement kinds of changes work against each other to some extent, but together they cause both policy and value function to approach optimality.

2 theorical assumptions

- the episodes have exploring starts,
- policy evaluation could be done with an infinite number of episodes

Policy evaluation is done exactly as described before.

Policy improvement is done by making the policy greedy with respect to the current value function.

Algorithm 2 Monte Carlo Exploring Starts, for estimating $\pi \approx \pi^*$

- 1: Initialize, $\forall s \in \mathcal{S}, a \in \mathcal{A}(s)$: 2: $Q(s, a) \leftarrow \text{arbitrary}$ 3: $\pi(s) \leftarrow \text{arbitrary}$ 4: $Returns(s, a) \leftarrow empty list$ 5: while forever do Choose $S_0 \in \mathcal{S}$ and $A_0 \in \mathcal{A}(S_0)$ (Pr(s, a) > 0) 6: Generate an episode starting from S_0 , A_0 , following π 7: for each pair s, a appearing in the episode do 8: $G \leftarrow$ the return that follows the first occurrence of s, a 9. Append G to Returns(s, a)10: $Q(s, a) \leftarrow average(Returns(s, a))$ 11: end for 12: 13: **for** each *s* in the episode **do** $\pi(s) \leftarrow \arg\max_a Q(s, a)$ 14: 15: end for

16: end while

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Monte Carlo Control without Exploring Start

How can we avoid the unlikely assumption of exploring starts? 2 approaches : on-policy methods and off-policy methods

- On-policy methods attempt to evaluate or improve the policy that is used to make decisions
- Off-policy methods evaluate or improve a policy different from that used to generate the data

In on-policy control methods the policy is generally soft, meaning that $\pi(a|s) > 0$ for all $s \in \mathcal{S}$ and all $a \in \mathcal{A}(s)$, but gradually shifted closer and closer to a deterministic optimal policy.

Example : ε -greedy policies (cf bandits)

Algorithm 3 On-policy first-visit MC control (for ε -soft policies) 1: **for** for all $s \in \mathcal{S}, a \in \mathcal{A}(s)$ **do**

- $Q(s,a) \leftarrow \text{arbitrary}$
 - $Returns(s, a) \leftarrow empty list$
- 4: end for
- 5: $\pi(a|s) \leftarrow$ an arbitrary ε —soft policy
- 6: while forever do
- Generate an episode using π
- - for each pair s, a appearing in the episode do
- 8: $G \leftarrow$ the return that follows the first occurrence of s, a 9:
- Append G to Returns(s, a) 10: $Q(s, a) \leftarrow average(Returns(s, a))$ 11:
- end for 12:
- 13: **for** each *s* in the episode **do**
- 14:
- $A^* \leftarrow \arg\max_a Q(s, a)$ (with ties broken arbitrarily) for all $a \in \mathcal{A}(s)$ do 15:
- $\pi(a|s) \leftarrow \left\{ \begin{array}{ll} 1 \varepsilon + \varepsilon/|\mathcal{A}(s)| & \text{if} \quad a = A^* \\ \varepsilon/|\mathcal{A}(s)| & \text{if} \quad a \neq A^* \end{array} \right.$ 16:
- end for 17:
- end for 18: 15/31

Off-policy MC

In off-policy methods these two functions are separated :

- ► The policy used to generate behavior, called the behavior policy, may in fact be unrelated to the policy that is evaluated and improved, called the target policy.
- An advantage of this separation is that the target policy may be deterministic (e.g., greedy), while the behavior policy can continue to sample all possible actions.
- require that the behavior policy has a nonzero probability of selecting all actions that might be selected by the target policy (coverage).

A potential problem is that this method learns only from the tails of episodes, when all of the remaining actions in the episode are greedy. If nongreedy actions are common, then learning will be slow, particularly for states appearing in the early portions of long episodes.

Algorithm 4 Off-policy MC prediction, for estimating $Q \approx q_{\pi}$

INPUT: an arbitrary target policy π for all $s \in \mathcal{S}$, $a \in \mathcal{A}(s)$ do

for all
$$s \in \mathcal{S}, a \in \mathcal{A}(s)$$
 d
 $Q(s, a) \leftarrow$ arbitrary
 $C(s, a) \leftarrow 0$

end for

while forever do

 $b \leftarrow$ any policy with coverage of π Generate an episode using b $G \leftarrow 0$

$$W \leftarrow 1$$

for
$$t = T - 1, T - 2, \cdots$$
 down to 0 do

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

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

 $C(S_t, A_t) \leftarrow C(S_t, A_t) + W$

$$Q(S_t, A_t) \leftarrow Q(S_t, A_t) + \frac{W}{C(S_t, A_t)} [G - Q(S_t, A_t)]$$

$$W \leftarrow W \frac{\pi(A_t|S_t)}{b(A_t|S_t)}$$

If W = 0 then exit For loop

Algorithm 5 Off-policy MC control, for estimating $\pi \approx \pi^*$ for all $s \in \mathcal{S}$, $a \in \mathcal{A}(s)$ do $Q(s, a) \leftarrow \text{arbitrary}$ $C(s,a) \leftarrow 0$ $\pi(s) \leftarrow \arg\max_a Q(s, a)$ (with ties broken consistently) end for while forever do $b \leftarrow \text{any soft policy}$ Generate an episode using b $G \leftarrow 0$ $W \leftarrow 1$ **for** $t = T - 1, T - 2, \cdots$ down to 0 **do**

for $t = T - 1, T - 2, \cdots$ down to 0 do $G \leftarrow \gamma G + R_{t+1}$ $G(S, A) \leftarrow G(S, A) + W(S, A) + W(S,$

$$C(S_t, A_t) \leftarrow C(S_t, A_t) + W$$
 $Q(S_t, A_t) \leftarrow Q(S_t, A_t) + \frac{W}{C(S_t, A_t)}[G - Q(S_t, A_t)]$
 $\pi(S_t) \leftarrow \arg\max_a Q(S_t, a)$ (with ties broken consistently)
If $A_t \neq \pi(S_t)$ then exit For loop

 $W \leftarrow W \frac{1}{b(A_t|S_t)}$ end for end while

Let's Play BlackJack

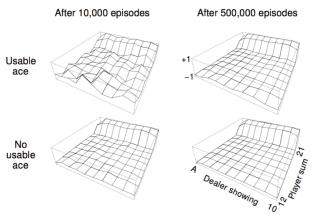
- ▶ 1 player (the agent) vs. 1 dealer
- Card values :
 - ▶ 10 for a face card (jack,queen,king),
 - ▶ 1 or 11 for an Ace (11 for the first encounter then 1),
 - Value on the card for any card between 2 and 9.
- Initialisation : 2 cards each, one is face down for the dealer
- ► Game : the player can request as many cards, one by one, as he wishes, until he either stop and exceed 21
- ▶ If the player has 21, he wins except if the dealer has 21 too.
- ightharpoonup Player goal : score \geq dealer, after both stops .
- Hypothesis : infinite deck

MDP for BlackJack

- ▶ 200 states :
 - ▶ player current sum : (12-21) [before hit]
 - ▶ Value of the face up card of the dealer : (1-10)
 - Ace already dealed (yes-no)
- 2 actions hit ou stick
- Reward:
 - ▶ 1 if win
 - -1 si lost (less than dealer or > 21)
 - 0 if null or continue
- $ightharpoonup \gamma = 1$
- ▶ Dealer policy : *hit* if total sum ≤ 17

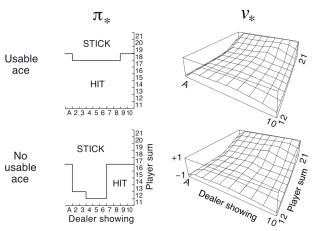
MC methods for BlackJack

Policy to be evaluated : stick if score \geq 20.



MC methods for BlackJack





Temporal Difference

TD learning is a combination of Monte Carlo ideas and dynamic programming (DP) ideas.

- learning with incomplete episodes
- model-free
- updating estimations with estimations

update from every-visit Monte Carlo method

$$V(S_t) \leftarrow V(S_t) + \alpha \left[G_t - V(S_t) \right]$$

Monte Carlo methods must wait until the end of the episode to determine the increment to $V(S_t)$. TD(0) will try to do the same with only one step of time

$$V(S_t) \leftarrow V(S_t) + \alpha \left[\underbrace{R_{t+1} + \gamma V(S_{t+1})}_{\hat{G}_t} - V(S_t)\right]$$

Algorithm 6 Algorithm TD(0) for estimating v_{π}

- 1: $v(s) \leftarrow 0 \forall s \in \mathcal{S}$
- 2: repeat
- 3: initialize *S*
- 4: repeat
- 5: $A \leftarrow \text{action given by } \pi \text{ for } s$
- 6: Take action A, observe R et s'

7:
$$v(s) \leftarrow v(s) + \alpha \left[R + \gamma v(s') - v(s) \right]$$

- 8: $s \leftarrow s'$
- 9: **until** s is terminal
- 10:
- 11: until all episodes are done

Advantages of TD Prediction Methods

- they do not require a model of the environment, of its reward and next-state probability distributions.
- they are naturally implemented in an on-line, fully incremental fashion.
- some Monte Carlo methods must ignore or discount episodes on which experimental actions are taken, which can greatly slow learning. TD methods are much less susceptible to these problems because they learn from each transition regardless of what subsequent actions are taken.

For any fixed policy π , TD(0) has been proved to converge to v_{π} , in the mean for a constant step-size parameter if it is sufficiently small, and with probability 1 if the step-size parameter decreases according to the usual stochastic approximation conditions.

Algorithm 7 Algorithm SARSA, on-policy TD control

- 1: initialize Q(s,a), for all $s \in \mathcal{S}$, $a \in \mathcal{A}(s)$, arbitrarily, and $Q(\text{terminal-state},\cdot) = 0$
- 2: repeat
- 3: Initialize *S*
- 4: Choose A from S using policy derived from Q (e.g., ε -greedy)
- 5: repeat
- 6: Take action A, observe R, S'
- 7: Choose A' from S' using policy derived from Q
- 8: $Q(S,A) \leftarrow Q(S,A) + \alpha (R + \gamma Q(S',A') Q(S,A))$
- 9: $S \leftarrow S'$, $A \leftarrow A'$
- 10: **until** S is terminal
- 11: until

Algorithm 8 Algorithm Q-Learning, off-policy TD control

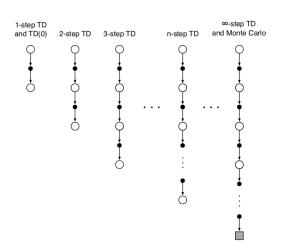
- 1: initialize Q(s,a), for all $s \in \mathcal{S}$, $a \in \mathcal{A}(s)$, arbitrarily, and $Q(\text{terminal-state}, \cdot) = 0$
- 2: while 1 do
- 3: Initialize *S*
- 4: $t \leftarrow 0$
- 5: repeat
- 6: Choose A from S using policy derived from Q
- 7: Take action A, observe R,S'
- 8: $Q(S,A) \leftarrow Q(S,A) + \alpha [R + \gamma \max_{a' \in A} Q(S',A') Q(S,A)]$
- 9: **until** *S* is terminal
- 10: end while

n-step Bootstrapping

unify the Monte Carlo (MC) methods and the one-step temporal-difference (TD) methods.

- free from the tyranny of the time step.
- a length of time in which a significant and recognizable state change has occurred.

The idea of n-step methods is usually used as an introduction to the algorithmic idea of eligibility traces, which enable bootstrapping over multiple time intervals simultaneously.



$$G_{t:t+n} = R_{t+1} + \gamma R_{t+2} + \dots + \gamma^{n-1} R_{t+n} + \gamma^n V_{t+n-1}(S_{t+n})$$

n-step TD for estimating $V \approx v_{\pi}$

```
Initialize V(s) arbitrarily, s \in S
Parameters: step size \alpha \in (0,1], a positive integer n
All store and access operations (for S_t and R_t) can take their index mod n
Repeat (for each episode):
   Initialize and store S_0 \neq \text{terminal}
   T \leftarrow \infty
   For t = 0, 1, 2, \dots:
       If t < T, then:
            Take an action according to \pi(\cdot|S_t)
            Observe and store the next reward as R_{t+1} and the next state as S_{t+1}
           If S_{t+1} is terminal, then T \leftarrow t+1
       \tau \leftarrow t - n + 1 (\tau is the time whose state's estimate is being updated)
       If \tau > 0:
           G \leftarrow \sum_{i=\tau+1}^{\min(\tau+n,T)} \gamma^{i-\tau-1} R_i
If \tau+n < T, then: G \leftarrow G + \gamma^n V(S_{\tau+n})
                                                                                                       (G_{\tau:\tau+n})
           V(S_{\tau}) \leftarrow V(S_{\tau}) + \alpha \left[ G - V(S_{\tau}) \right]
   Until \tau = T - 1
```

n-step Sarsa for estimating $Q \approx q_*$, or $Q \approx q_\pi$ for a given π

```
Initialize Q(s, a) arbitrarily, for all s \in S, a \in A
Initialize \pi to be \varepsilon-greedy with respect to Q, or to a fixed given policy
Parameters: step size \alpha \in (0,1], small \varepsilon > 0, a positive integer n
All store and access operations (for S_t, A_t, and R_t) can take their index mod n
Repeat (for each episode):
   Initialize and store S_0 \neq \text{terminal}
   Select and store an action A_0 \sim \pi(\cdot|S_0)
   T \leftarrow \infty
   For t = 0, 1, 2, \dots:
       If t < T, then:
           Take action A_t
           Observe and store the next reward as R_{t+1} and the next state as S_{t+1}
           If S_{t+1} is terminal, then:
               T \leftarrow t + 1
           else:
               Select and store an action A_{t+1} \sim \pi(\cdot|S_{t+1})
       \tau \leftarrow t - n + 1 (\tau is the time whose estimate is being updated)
       If \tau > 0:
           G \leftarrow \sum_{i=\tau+1}^{\min(\tau+n,T)} \gamma^{i-\tau-1} R_i
           If \tau + n < T, then G \leftarrow G + \gamma^n Q(S_{\tau+n}, A_{\tau+n})
                                                                                                       (G_{\tau \cdot \tau + n})
           Q(S_{\tau}, A_{\tau}) \leftarrow Q(S_{\tau}, A_{\tau}) + \alpha \left[ G - Q(S_{\tau}, A_{\tau}) \right]
           If \pi is being learned, then ensure that \pi(\cdot|S_{\tau}) is \varepsilon-greedy wrt Q
   Until \tau = T - 1
```

Off-policy n-step Sarsa for estimating $Q \approx q_*$, or $Q \approx q_\pi$ for a given π

```
Input: an arbitrary behavior policy b such that b(a|s) > 0, for all s \in S, a \in A
Initialize Q(s, a) arbitrarily, for all s \in S, a \in A
Initialize \pi to be \varepsilon-greedy with respect to Q, or as a fixed given policy
Parameters: step size \alpha \in (0,1], small \varepsilon > 0, a positive integer n
All store and access operations (for S_t, A_t, and R_t) can take their index mod n
Repeat (for each episode):
   Initialize and store S_0 \neq \text{terminal}
   Select and store an action A_0 \sim b(\cdot|S_0)
   T \leftarrow \infty
   For t = 0, 1, 2, \dots:
        If t < T, then:
            Take action A_t
            Observe and store the next reward as R_{t+1} and the next state as S_{t+1}
            If S_{t+1} is terminal, then:
                T \leftarrow t + 1
            else:
                 Select and store an action A_{t+1} \sim b(\cdot | S_{t+1})
        \tau \leftarrow t - n + 1 (\tau is the time whose estimate is being updated)
        If \tau > 0:
             \rho \leftarrow \prod_{i=\tau+1}^{\min(\tau+n-1,T-1)} \frac{\pi(A_i|S_i)}{b(A_i|S_i)} 
 G \leftarrow \sum_{i=\tau+1}^{\min(\tau+n,T)} \gamma^{i-\tau-1} R_i 
                                                                                                         (\rho_{\tau+1:t+n-1})
            If \tau + n < T, then: G \leftarrow G + \gamma^n Q(S_{\tau+n}, A_{\tau+n})
                                                                                                         (G_{\tau \cdot \tau \perp n})
            Q(S_{\tau}, A_{\tau}) \leftarrow Q(S_{\tau}, A_{\tau}) + \alpha \rho \left[ G - Q(S_{\tau}, A_{\tau}) \right]
            If \pi is being learned, then ensure that \pi(\cdot|S_{\tau}) is \varepsilon-greedy wrt Q
    Until \tau = T - 1
```