

RL Lecture 4 - Model free Learning

what we did earlier using DP, required the agent to know the entire model dynamics i.e. the ^{state probabilities} transition reward functions, etc. But since in most real life problems, the agent wouldn't know what the entire dynamics of the environment are, what can we do? That is discussed now! \rightarrow Model-free Learning

Just like earlier, we break ^{it} into 2 components:

- ① Policy evaluation \rightarrow model free prediction
 - ② Finding the optimal policy \rightarrow model free control (Lecture 5).
- estimate the value function of an unknown MDP.

Model free prediction

\rightarrow Monte Carlo Learning:

- learns directly from episodes of experience
- Look only at complete episodes
- Idea \rightarrow value = mean of the sample returns across many episodes.

Problem: Applicable to only episodic MDPs, i.e. those ^{where} episodes ~~which~~ terminate.

Monte Carlo Policy Evaluation:

- Policy already given $\Rightarrow \pi$ following
- Learn V_π from episodes of experience $\sim \pi$.
- $\# S_1, A_1, R_2, \dots S_k \sim \pi$.

Earlier: we had expected return:

$$V_\pi(s) = E_\pi(G_T | S_t = s).$$

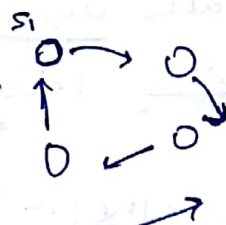
But here, we'll have empirical mean instead of expected.

Challenge: How to reset to a particular ~~step~~ ^{state} everytime and start our evaluation? And how to do this for every state?

2 methods:

① First visit: Imagine: circular MDP i.e. loop in an MDP.

Now, we consider only the time when a particular state, say s_1 , is 1st visited in an episode.



After that, update the count of # visits in that state: $N(s) \pm 1$
 $S(s) = S(s) + (G_t)$

← add reward from that time step onwards to the total reward for the state

estimated value for the state:

$$V(s) = \frac{S(s)}{N(s)}$$

According to ^{the} law of large numbers, the mean of a large # samples actually converges to the true mean

$$P\left(\left|\frac{X_1 + X_2 + \dots + X_n}{n} - \mu\right| < \epsilon\right) \rightarrow 1$$

$X_1, X_2, \dots, X_n \rightarrow n$ samples of the dists

$\mu \rightarrow$ actual mean of the distribution

as $n \rightarrow \infty, \forall \epsilon > 0$.
 $\epsilon \rightarrow$ any arbitrary small +ve value.

Clarifications :

Q) Does # states ~~depend~~ decide the # episodes required?

No!

From Central Limit Theorem, the variance decrease as a function of $1/n$. Hence, it doesn't depend on the # states, just how many times a particular state is visited

The trajectory in each episode should be such that it covers most of the states that we care about.

~~cover~~

Q) How to make sure that - we reach all the states we care about?

→ Problem of control. (next lecture)

→ Here, we care about states under policy π and we ^{make sure to} visit those states by following policy π .

② Every visit → instead of just the 1st visit, add on for every visit → i.e. multiple visits possible for each episode.

(Choice depends on the domain).

Incremental mean:

$$\mu_k = \frac{1}{k} \left(\sum_{i=1}^k x_i \right) = \frac{1}{k} \left[x_k + \sum_{i=1}^{k-1} x_i \right]$$

$$= \frac{1}{k} (x_k + (k-1) \mu_{k-1}) = \mu_{k-1} + \frac{1}{k} (x_k - \mu_{k-1})$$

Thus, we take a small step from our previous expectation following a error function.

↓ error term

↪ as difference betⁿ what we expected and what we actually got

Incremental Monte Carlo updates:

$$N(s) = N(s) + 1.$$

$$V(s) = V(s) + \frac{1}{N(s)} (G_T - V(s)).$$

(Idea \rightarrow want to do online learning instead of computing all values beforehand & then taking the mean).

Better \rightarrow add a weightage to discard old terms/values.

$$\begin{aligned} V(s) &= V(s) + \alpha (G_T - V(s)) \\ &= (1 - \alpha) V(s) + \alpha G_T \end{aligned}$$

We can weigh older terms accordingly.

Temporal Difference Learning: (TD)

* Learns from incomplete episodes \rightarrow bootstrapping
 \rightarrow don't need to get to the end and figure out the reward, rather, go to an intermediate step and estimate the reward from there to the end.

So, we have to make an initial guess and after making, taking some steps, we make another guess of the final reward from that point. Then we try to bring our initial guess closer to that guess.

Goal: Learn V_π online from experience under π .

Here, we update $V(s_t)$ towards the ~~expected~~ estimated return:

Simple TD learning: $V(s_t) = V(s_t) + \underbrace{(R_{t+1} + \gamma V(s_{t+1}) - V(s_t))}_{\text{TD target}}$

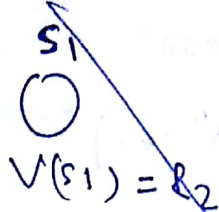
TD(0)

Advantages/Disadvantages

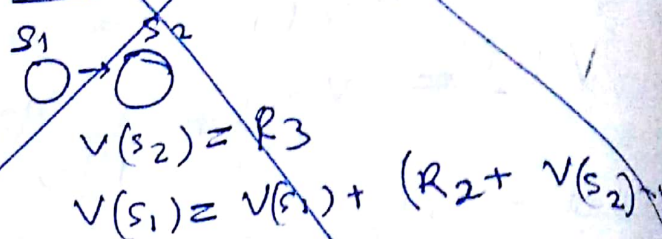
- \rightarrow TD can learn before knowing the final outcome. i.e. online after every step.
- \rightarrow TD can learn without the final outcome. i.e. it can be applied to processes that do not end.
Hence, learning from (episodes) incomplete sequences.

TD update:

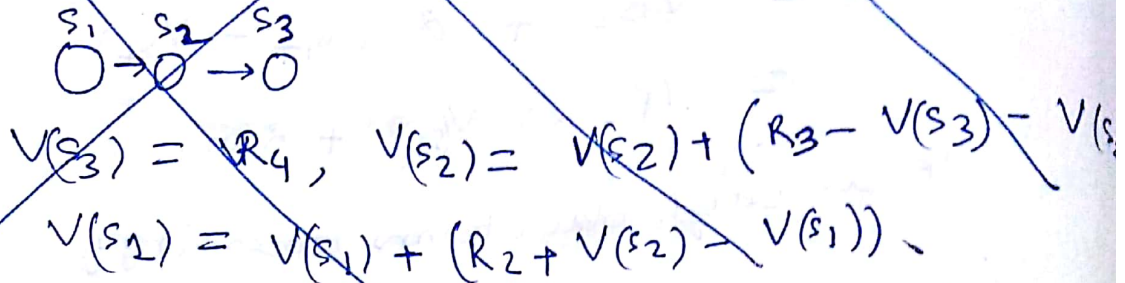
Step 1 →



Step 2:



Step 3:



i.e. update after each step.

TD → less noise

$$G_T = R_{t+1} + \gamma R_{t+2} + \dots$$

$\gamma^{T-1} R_T \rightarrow$ unbiased estimate of $V_{\pi}(s_t)$.

(bias of an estimate is given by: $\hat{\theta}_m$)

$$\text{Bias}(\hat{\theta}_m) = E(\hat{\theta}_m) - \theta, \quad \hat{\theta} \rightarrow \text{unbiased estimator when } E(\hat{\theta}_m) = \theta \text{ i.e.}$$

↳ expected

$$E(\hat{\theta}_m) = \theta \text{ i.e.}$$

$$\text{Bias}(\hat{\theta}_m) = 0$$

$$R_{t+1} + \gamma V_{\pi}(s_{t+1}) \rightarrow \text{unbiased estimate of } V_{\pi}(s_t)$$

But, we use:

$$R_{t+1} + \gamma V(s_{t+1}) \rightarrow \text{biased estimation.}$$

we don't know the true future $V_{\pi}(s_{t+1})$.

$$\text{But } G_T = R_{t+1} + \gamma R_{t+2} + \dots + \gamma^{T-1} R_T$$

↳ adds noise at each stage. (term)

$$R_{t+1} + \gamma V(s_{t+1}) \rightarrow \text{noise just in one term.}$$

Hence, TD → less noisy.

MC \rightarrow high variance, 0 bias

- \rightarrow Good convergence
- \rightarrow independent of initializato.

TD \rightarrow some bias, low variance.

- \rightarrow more efficient
- \rightarrow TD(0) \rightarrow $V_{\pi}(s)$
- \rightarrow sensitive to init value.

Batch MC / TD

Both converge as experience $\rightarrow \infty$

In batch case,

\rightarrow repeatedly sample

$$k \in [1, K]$$

Do they converge?

(eg. run 3 episodes & repeatedly iterate & learn from those 3).

eg. There are observed

1 A 0 B 0

2 B 1

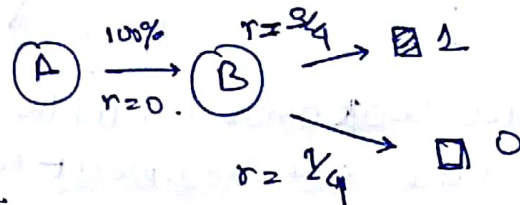
7 B 1

8 B 0

Monte carls for B = $\frac{6}{8}$

Monte carls for A = $\frac{0}{1}$

TD estimate \rightarrow tries to build an MDP that best explains the data.



MC \rightarrow reduces the mean squared error

• Fit fit to actual returns received $\rightarrow \sum_{k=1}^K \sum_{t=1}^{T_k} (r_t^k - V(s_t^k))$

TD \rightarrow converges to the MDP that best fits the data

\rightarrow set to max

likelihood Markov model

\rightarrow just by counting normal

$$P_{ss'}^a = \frac{1}{N(s,a)} \sum_{k=1}^K \sum_{t=1}^{T_k} \mathbb{I}(s_t^k, a_t^k, s_{t+1}^k = s', a, s)$$

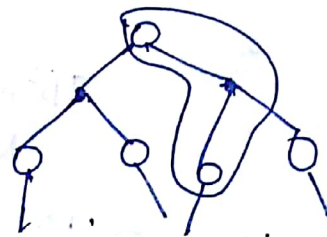
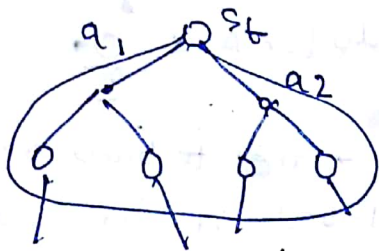
$$R_s^a = \frac{1}{N(s,a)} \sum_{k=1}^K \sum_{t=1}^{T_k} \mathbb{I}(s_t^k, a_t^k = s, a) r_t^k$$

TD makes use of markov property \rightarrow much more efficient in these environments.
situations may be non-markov.

ie. partially observed \overline{MDP} \rightarrow under the hood, there is an MDP, but what is observed is partial \rightarrow so TD(0) doesn't work well on such environments & MC still finds a reasonable solution.

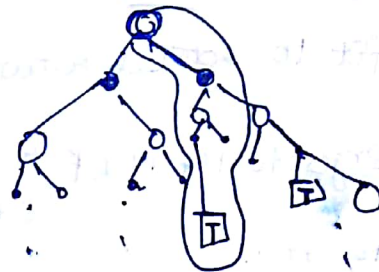
In DP \rightarrow we did a one-step lookahead over all possible next states.

Whereas, for eg. in TD, we just look over one step, one sample



In MC, we look over until the entire terminate but only take one sample at each stage.

We could take all the way to the bottom for DP too, which would be called exhaustive tree search



\rightarrow Bootstrap \rightarrow don't use the real value but the estimated value of the next step to update the current step.

- Not by MC
- By DP & TD

\rightarrow sampling \rightarrow don't do full width backup & only take samples.

- MC ^{TD} samples
- DP doesn't

Exhaustive search \rightarrow Bad!

Algorithms that are in the middle full width $\&$ shallow or deep backups $\&$ sample backups

TD(λ)

n-step prediction \rightarrow look n steps into the future [instead of one in TD(0)]

$n=1 \rightarrow TD(0): G_T^{(1)} = R_{t+1} + \gamma V(S_{t+1})$

$n=2: G_T^{(2)} = R_{t+1} + \gamma R_{t+2} + \gamma^2 V(S_{t+2})$

$n=\infty$ MC $\rightarrow G_T^{(\infty)} = R_{t+1} + \gamma R_{t+2} + \dots + \gamma^{T-t} R_T + \gamma^{T-t+1} V(S_T)$

n-step return:

$G_T^{(n)} = R_{t+1} + \gamma R_{t+2} + \dots + \gamma^{n-1} R_{t+n} + \gamma^n V(S_{t+n})$

n-step TD $\rightarrow V(S_t) = V(S_t) + \alpha (G_T^{(n)} - V(S_t))$

What value of n to be used? \rightarrow Hard to choose & depends on problem

online / offline update \rightarrow update the value immediately \rightarrow update the value only after reaching the end \rightarrow Use average n-step returns over several n.

how to do this efficiently for all n.

$G_T^{(2)} + G_T^{(4)} \rightarrow$ get the best of all n $\frac{2}{2}$

TD(λ)

$\lambda \rightarrow 0 < \lambda < 1$ \rightarrow return \rightarrow geometrically weighted average of all n decaying for each specific n.

(1-λ) \rightarrow normalizing factor:

$(1-\lambda) + (1-\lambda)\lambda + \dots$

Sum $= (1-\lambda) \frac{1}{(1-\lambda)} = 1$

~~$G_T^{(2)} = (1-\lambda) \lambda + (1-\lambda) \lambda^2 + \dots$~~

$$G_t^\lambda = (1-\lambda) \sum_{n=1}^{\infty} G_t^{(n)} \lambda^{n-1}$$

$$TD(\lambda) \Rightarrow V(s_t) = V(s_t) + \alpha (G_T^\lambda - V(s_t))$$

Why geometric mean? → computational complexity
 as $(1-\lambda), (1-\lambda)\lambda, (1-\lambda)\lambda^2 \dots$ → memoryless → doesn't require storing separate values for each timestep!
 (just store one value and keep multiplying λ).

This is the forward view $TD(\lambda)$ → needs to wait until completing the ~~episode~~ episode → suffers from same disadvantages as MC. ($\lambda = 1 \rightarrow MC$)

Backward view → $TD(\lambda)$: compute the above stuff more efficiently without

having to wait till the completion of the episode.

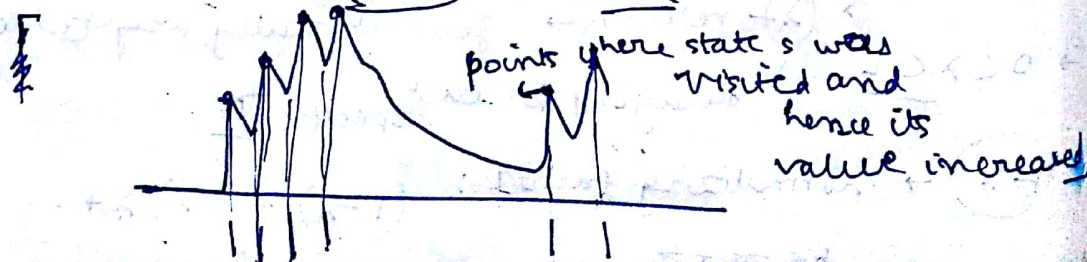
Eligibility traces

- Frequency heuristic → assign importance to most frequent states
 - Recency heuristic → importance to most recent states
- eligibility trace → combines both

$$E_0(s) = 0$$

$$E_t(s) = \gamma \lambda E_{t-1}(s) + I(s_t = s)$$

→ whenever state is visited we increase the trace & if not visited, we reduce its value



$$I(S_t = s) \rightarrow \text{Identity function} = \begin{cases} 1 & \text{if } S_t = s \\ 0 & \text{else} \end{cases}$$

For backward view:

→ Keep ET for each state

→ update $V(s)$ in proportion to $E_t(s)$ & TD-error δ_t

$$\delta_t = R_{t+1} + \gamma V(S_{t+1}) - V(S_t)$$

for each state.

$$V(s) = V(s) + \alpha E_t(s) \delta_t$$

→ $\lambda \rightarrow$ decaying factor.

→ $\gamma = 0 \rightarrow$ we squash the values to 0 of $S_t \neq s$

$$\therefore E_t(s) = I(S_t = s)$$

TP(0) Thus, value function updated ^{only} if we observe that state in that time step. visit

$$V(s) = V(s) + \alpha \delta_t E_t(s)$$

$\lambda = 1 \rightarrow$ same total update as MC.

Theorem \rightarrow total updates same for backward as well as Forward views.

$$\sum_{t=1}^T \alpha \delta_t E_t(s) = \sum_{t=1}^T \alpha (G_t - V(S_t)) I(S_t = s)$$

TD(λ) \rightarrow Spectrum. betw TD(0) & MC.

$E_t(s) \rightarrow$ can be computed efficiently too.

