

Lecture 9: Model-Free Prediction To

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A, f(a) + (1-7,) f(b)

Outline

- Introduction
- Monte-Carlo Learning
- Temporal-Difference Learning
- 4 n-step TD Methods
- TD (λ)
- 6 References

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- Introduction
- 2 Monte-Carlo Learning
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Model-Free Reinforcement Learning

- Last lecture:
 - Planning by dynamic programming
 - ▶ Solve a known MDP
- This lecture:
 - Model-free prediction
 - Estimate the value function of an unknown MDP
- Next lecture:
 - Model-free control
 - Optimize the value function of an unknown MDP

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Monte-Carlo Reinforcement Learning

- MC methods learn directly from episodes of experience
- MC is model-free: no knowledge of MDP transitions / rewards
- MC learns from complete episodes: no bootstrapping
- MC uses the simplest possible idea: value = mean return
- To learn values & policies, MC can be used in two ways:
 - model-free: no model necessary and still attains optimality
 - simulated: needs only a simulation, not a full model
- Caveat: can only apply MC to episodic MDPs
 - All episodes must terminate

Fluite Stage

Monte-Carlo Policy Evaluation

• Goal: learn v_{π} from episodes of experience under policy π

$$S_1, A_1, R_2, \ldots, S_k \sim \pi$$

Recall that the return is the total discounted reward:

$$\underline{G_t} = R_{t+1} + \gamma R_{t+2} + \ldots + \gamma^{T-1} R_T$$

• Recall that the value function is the expected return:

$$v_{\pi}(s) = \mathbb{E}_{\pi}[G_t|S_t = s]$$

 Monte-Carlo policy evaluation uses empirical mean return instead of expected return



Monte-Carlo Policy Evaluation

- Goal: learn $v_{\pi}(s)$
- Given: some number of episodes under π which contains s
- Idea: average returns observed after visits to s
- Every-Visit MC: average returns for every time s is visited in an episode
- First-visit MC: average returns only for first time s is visited in an episode
- Both converge asymptotically

First-Visit Monte-Carlo Policy Evaluation

- To evaluate state s
- The first time-step <u>t</u> that state <u>s</u> is <u>visited</u> in an episode
- Increment counter $N(s) \leftarrow N(s) + 1$
- Increment total return $S(s) \leftarrow S(s) + G_t$
- Value is estimated by mean return V(s) = S(s)/N(s)
- ullet By law of large numbers, $V(s)
 ightarrow
 u_\pi(s)$ as $N(s)
 ightarrow \infty$

First-Visit Monte-Carlo Policy Evaluation

```
First-visit MC prediction, for estimating V \approx v_{\pi}

Input: a policy \pi to be evaluated Initialize:

V(s) \in \mathbb{R}, \text{ arbitrarily, for all } s \in \mathcal{S}
Returns(s) \leftarrow \text{ an empty list, for all } s \in \mathcal{S}
Loop forever (for each episode):
Generate \text{ an episode following } \pi\colon S_0, A_0, R_1, S_1, A_1, R_2, \ldots, S_{T-1}, A_{T-1}, R_T
G \leftarrow 0
Loop for each step of episode, t = T-1, T-2, \ldots, 0:
G \leftarrow \gamma G + R_{t+1}
Unless S_t appears in S_0, S_1, \ldots, S_{t-1}:
Append G \text{ to } Returns(S_t)
V(S_t) \leftarrow \text{average}(Returns(S_t))
```

Every-Visit Monte-Carlo Policy Evaluation

- To evaluate state s
- Every time-step t that state s is visited in an episode
- Increment counter $N(s) \leftarrow N(s) + 1$
- Increment total return $S(s) \leftarrow S(s) + G_t$
- ullet Value is estimated by mean return V(s) = S(s)/N(s)
- ullet Again, $V(s)
 ightarrow v_\pi(s)$ as $N(s)
 ightarrow \infty$

Incremental Mean

The mean μ_1, μ_2, \ldots of a sequence x_1, x_2, \ldots can be computed incrementally,

$$\underline{\mu_k} = \frac{1}{k} \sum_{j=1}^k x_j$$

$$= \frac{1}{k} \left(x_k + \sum_{j=1}^{k-1} x_j \right)$$

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

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

Incremental Monte-Carlo Updates

- Update V(s) incrementally after episode $S_1, A_1, R_2, \dots, S_T$
- For each state S_t with return G_t

$$N(S_t) \leftarrow N(S_t) + 1 \ V(S_t) \leftarrow V(S_t) + rac{1}{N(S_t)}(G_t - V(S_t))$$

 In non-stationary problems, it can be useful to track a running mean, i.e. forget old episodes.

$$V(S_t) \leftarrow V(S_t) + QG_t - V(S_t)$$

$$Step-size$$



Monte-Carlo Estimation of Action Values

- Monte Carlo is most useful when a model is not available: we want to learn q_*
- $q_{\pi}(s, a)$: average return starting from state s and action a following policy π
- Converges asymptotically if every state-action pair is visited
- Exploring Starts: every state-action pair has a non-zero probability of being the starting pair

Backup Diagram for Monte-Carlo

- Entire rest of episode included
- Only one choice considered at each state (unlike DP)
- thus, there will be an explore/exploit dilemma
- Does not bootstrap from successor states's values (unlike DP)
- Time required to estimate one state does not depend on the total number of states



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Temporal-Difference Learning

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- TD methods learn directly from episodes of experience
- TD is model-free: no knowledge of MDP transitions / rewards
- TD learns from incomplete episodes, by bootstrapping
- TD updates a guess towards a guess

Existing samples: 2, 4,5,6,6

Vesamply with replacement

2.5, 5.6,6

4.5.6,6.6

2,2,4,5,7

bootstrap
samples.

MC and TD

- Goal: learn v_{π} online from experience under policy π
- Incremental every-visit Monte-Carlo

▶ Update value $V(S_t)$ toward actual return G_t

$$V(S_t) \leftarrow V(S_t) + \alpha(\underline{G_t} - V(S_t)) \checkmark$$

- Simplest temporal-difference learning algorithm: TD(0)
 - ▶ Update value $V(S_t)$ toward estimated return $R_{t+1} + \gamma V(S_{t+1})$

$$V(S_t) \leftarrow V(S_t) + \alpha(R_{t+1} + \gamma V(S_{t+1}) - V(S_t))$$

- $R_{t+1} + \gamma V(S_{t+1})$ is called the *TD target*
- $\delta_t = R_{t+1} + \gamma V(S_{t+1}) V(S_t)$ is called the *TD error*

Advantages and Disadvantages of MC vs. TD

- TD can learn before knowing the final outcome
 - ► TD can learn online after every step (less memory & peak computation)
 - MC must wait until end of episode before return is known
- TD can learn without the final outcome
 - TD can learn from incomplete sequences
 - MC can only learn from complete sequences
 - ► TD works in continuing (non-terminating) environments
 - MC only works for episodic (terminating) environments

Bias/Variance Trade-Off

- Return $G_t = R_{t+1} + \gamma R_{t+2} + \ldots + \gamma^{T-1} R_T$ is unbiased estimate of $v_{\pi}(S_t)$
- True TD target $R_{t+1} + \gamma (S_{t+1})$ is unbiased estimate of $v_{\pi}(S_t)$
- TD target $R_{t+1} + \gamma V(S_{t+1})$ is biased estimate of $v_{\pi}(S_t)$
- TD target is much lower variance than the return:
 - Return depends on many random actions, transitions, rewards
 - ▶ TD target depends on one random action, transition, reward

Advantages and Disadvantages of MC vs. TD (2)

- MC has high variance, zero bias
 - Good convergence properties
 - (even with function approximation)
 - Not very sensitive to initial value
 - Very simple to understand and use
- TD has low variance, some bias
 - Usually more efficient than MC
 - ▶ TD(0) converges to $v_{\pi}(s)$
 - (but not always with function approximation)
 - More sensitive to initial value

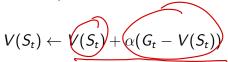
Advantages and Disadvantages of MC vs. TD (3)

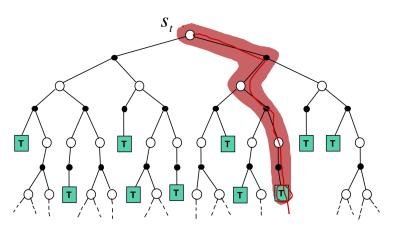
- TD exploits Markov property
 - Usually more efficient in Markov environments
- MC does not exploit Markov property
 - Usually more effective in non-Markov environments
- MC has lower error on past data, but higher error on future data

Bellman Backup

- The term "Bellman backup" comes up quite frequently in the RL literature.
- The Bellman backup for a state (or a state-action pair) is the right-hand side of the Bellman equation: the reward-plus-next-value.
- Under different algorithms, we obtain
 - Monte-Carlo Backup
 - Temporal-Difference Backup
 - Dynamic Programming Backup

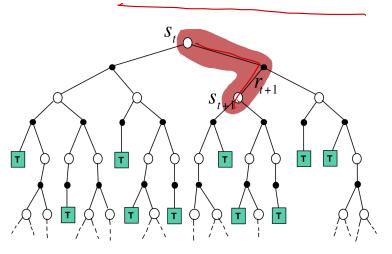
Monte-Carlo Backup



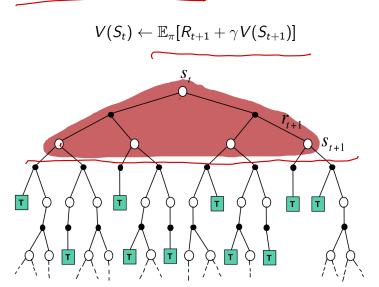


Temporal-Difference Backup

$$V(S_t) \leftarrow V(S_t) + \alpha(R_{t+1} + \gamma V(S_{t+1}) - V(S_t))$$



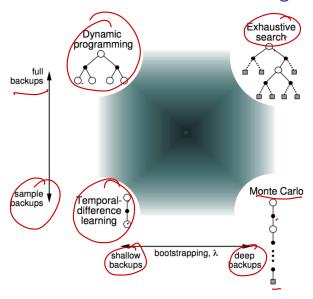
Dynamic Programming Backup



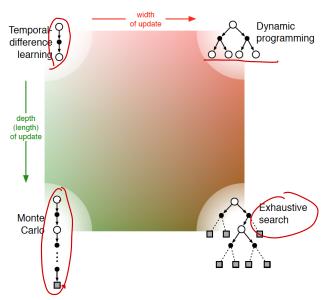
Bootstrapping and Sampling

- Bootstrapping: update involves an estimate
 - MC does not bootstrap
 - DP bootstraps
 - TD bootstraps
- Sampling: update samples an expectation
 - MC samples
 - DP does not sample
 - TD samples

Unified View of Reinforcement Learning



Unified View of Reinforcement Learning

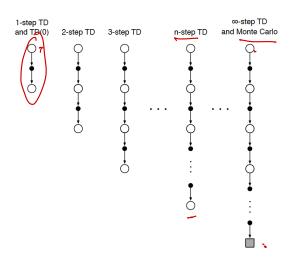


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n-Step Prediction

• Let TD target look *n* steps into the future



n-Step Return

• Consider the following *n*-step returns for $n = 1, 2, \infty$:

$$n = 1 \quad (TD) \quad G_{t}^{(1)} = R_{t+1} + \gamma V(S_{t+1})$$

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

$$\vdots \quad \vdots$$

$$n = \infty \quad (MC) \quad G_{t}^{(\infty)} = R_{t+1} + \gamma R_{t+2} + \dots \gamma^{T-1} R_{T}$$

Define the *n*-step return

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

• *n*-step temporal-difference learning

$$V(S_t) \leftarrow V(S_t) + \alpha \underbrace{(G_t^{(n)} - V(S_t))}$$

n-step TD

Recall the *n*-step return:

$$G_{t}^{(n)} \doteq R_{t+1} + \gamma R_{t+2} + \dots + \gamma^{n-1} R_{t+n} + \gamma^{n} V_{t+n-1} \left(S_{t+n} \right), \ n \geq 1, 0 \leq t < T-n$$

- Of course, this is <u>not available</u> until time t + n
- The natural algorithm is thus to wait until then

$$V_{t+n}(S_t) \doteq V_{t+n-1}(S_t) + \alpha \left[G_t^{(n)} - V_{t+n-1}(S_t) \right], \quad 0 \le t < T$$

• This is called *n*-step TD

n-step TD Algorithm

```
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
                                                                                           (G_{\tau}^{(n)})
           If \tau + n < T, then: G \leftarrow G + \gamma^n V(S_{\tau+n})
           V(S_{\tau}) \leftarrow V(S_{\tau}) + \alpha \left[G - V(S_{\tau})\right]
   Until \tau = T - 1
```

Error Reduction Property

Error reduction property of n-step returns

$$\max_{s} \left| \mathbb{E}_{\pi} \left[G_{t}^{(n)} \middle| S_{t} = s \right] - v_{\pi}(s) \right| \leq \gamma^{n} \max_{s} \left| V_{t}(s) - v_{\pi}(s) \right|$$

$$\text{Maximum error using } n\text{-step return} \qquad \text{Maximum error using V}$$

- Using this property, we can show that n-step TD methods converge
- n-step TD methods: a family of sound methods including one-step TD methods & MC methods as extreme members

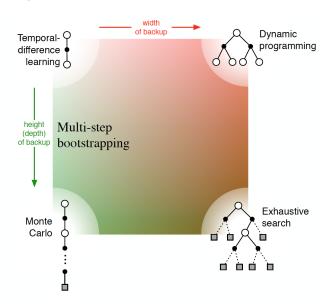
Summary of n-step TD Methods

- Generalize Temporal-Difference and Monte Carlo learning methods, sliding from one to the other as *n* increases
 - ▶ n = 1 is TD
 - ▶ $n = \infty$ is MC
 - ▶ an intermediate *n* is often much better than either extreme
 - applicable to both continuing and episodic problems
- There is some cost in computation
 - need to remember the last n states
 - learning is delayed by n steps
 - per-step computation is small and uniform, like TD
- Everything generalizes nicely: error-reduction theory

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- $5 \text{ TD}(\lambda)$
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Unified View

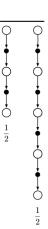


Averaging *n*-Step Returns

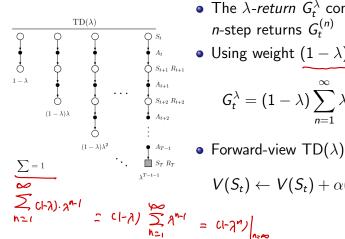
- We can average n-step returns over different n
- e.g. average the 2-step and 4-step returns

$$\underbrace{\frac{1}{2}G^{(2)} + \frac{1}{2}G^{(4)}}_{}$$

- Combines information from two different time-steps
- Can we efficiently combine information from all time-steps?



λ -return



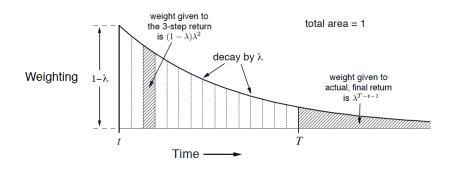
- The λ -return G_t^{λ} combines all *n*-step returns $G_t^{(n)}$
- Using weight $(1-\lambda)\lambda^{n-1}$

Using weight
$$(1-\lambda)\lambda^{n-1}$$
 S_{t+1}
 S_{t+1}
 S_{t+2}
 S_{t+2

$$V(S_t) \leftarrow V(S_t) + \alpha(G_t^{\lambda}) V(S_t)$$



$\mathsf{TD}(\lambda)$ Weighting Function



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



Relation to TD(0) & MC

• The λ -return can be rewritten as:

$$G_t^{\lambda} = \underbrace{(1-\lambda)\sum_{n=1}^{T-t-1}\lambda^{n-1}G_t^{(n)}}_{\text{Until termination}} + \underbrace{\lambda^{T-t-1}G_t}_{\text{After termination}}$$

• if $\lambda = 1$, you get the MC target:

$$G_t^{\lambda} = (1-1) \sum_{n=1}^{T-t-1} 1^{n-1} G_t^{(n)} + 1^{T-t-1} G_t = G_t$$

• If $\lambda = 0$, you get the TD(0) target:

$$G_t^{\lambda} = (1-0) \sum_{n=1}^{T-t-1} 0^{n-1} G_t^{(n)} + 0^{T-t-1} G_t = G_t^{(1)}$$

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Summary of $TD(\lambda)$ algorithms

- Another way of interpolating between MC and TD methods
- A way of implementing compound λ -return targets

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Main References

- Reinforcement Learning: An Introduction (second edition), R. Sutton & A. Barto, 2018.
- RL course slides from Richard Sutton, University of Alberta.
- RL course slides from David Silver, University College London.