Reinforcement Learning Free model Algorithms

CS-UPC

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Recap

- Definition of RL
- Framework
- Concepts learned:
 - Model
 - ► Policy: deterministic and non-deterministic
 - Reward functions, immediate reward
 - Discounted and undiscounted Long-term reward
 - ... γ , π , markovian condition
- Value functions
- Bellman equation
- Policy evaluation: Value iteration and algebraic method
- Optimal policy, greedy policy and relation with Value functions
- Dynamic programming methods: Value iteration, Policy iteration, asynchronous methods

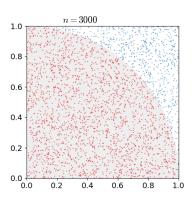
Goal of this lecture

- Problems with Dynamic Programming methods
 - Sweep of full steps or random steps
 - Need to know the model
- We'll see now methods that do not require a model but only experiences to build evaluations of policies and also to find optimal policies
- Methods we'll see:
 - ► Monte-Carlo
 - ► Q-learning
 - Temporal differences; n-steps and TD(λ)
 - Sarsa, Expected Sarsa
- Off-line vs. on-line learning
- Importance Sampling

Monte-Carlo methods

Monte-Carlo reinforcement learning

- Monte Carlo (MC): When interested in some complex value, instead
 of computing it, try to estimate it by sampling.
- Example: measure an irregular area.



• Refresher about expectations:

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For continuous variables

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Expectation computation by sampling

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Monte-Carlo Policy Evaluation

ullet 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:

$$R_t = r_{t+1} + \gamma r_{t+2} + \ldots + \gamma^{T-1} r_T$$

• Recall that the value function is the expected return:

$$V^{\pi}(s) = \mathbb{E}_{\pi}[R_t|S_t = s] pprox rac{1}{N}\sum_{i=1}^N R_i$$

where R_i is obtained from state s under π distribution (following π)

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

Monte-Carlo reinforcement learning

- MC uses the simplest possible idea: value = mean return. Instead of computing expectations, sample the long term return under the policy
- MC methods learn directly from episodes of experience
- MC is model-free: no explicit knowledge of environment mechanisms
- MC learns from complete episodes
 - Caveat: can only be applied to complete episodic environments (all episodes must terminate).

Monte-Carlo policy evaluation

Monte Carlo policy evaluation

```
Given \pi, the policy to be evaluated, initialize V randomly Returns(s) \longleftarrow empty list, \forall \in S repeat

Generate trial using \pi for each s in trial do

R \leftarrow return following the first occurrence of s

Append R to Returns(s)

V(s) \leftarrow average(Returns(s))

end for

until true
```

Monte-Carlo Policy Evaluation

- How to average results for V(s)? Every time-step t that state s is visited in an episode:
 - ▶ Increment counter $N(s) \leftarrow N(s) + 1$
 - ▶ Increment total return $S(s) \leftarrow S(s) + R_t$
 - ▶ Value is estimated by mean return V(s) = S(s)/N(s)
- ullet By law of large numbers, $V(s) o V^\pi(s)$ as $N(s) o \infty$ for all states
- ullet However, for each state you should store S and N.

• Update V(s) incrementally:

$$V_n(S_t) = \frac{1}{n} \sum_{i=1}^n R_i$$

$$V_n(S_t) = \frac{1}{n} \left(R_n + \sum_{i=1}^{n-1} R_i \right)$$

$$V_n(S_t) = \frac{1}{n} \left(R_n + (n-1)V_{n-1}(S_t) \right)$$

$$V_n(S_t) = \frac{1}{n} R_n + \frac{1}{n} \left((n-1)V_{n-1}(S_t) \right)$$

$$V_n(S_t) = \frac{1}{n} R_n + V_{n-1}(S_t) - \frac{1}{n} V_{n-1}(S_t)$$

$$V_n(S_t) = V_{n-1}(S_t) + \frac{1}{n} (R_n - V_{n-1}(S_t))$$

- Compute return R_t
- For each state S_t with return R_t

$$N(S_t) \leftarrow N(S_t) + 1$$

 $V(S_t) \leftarrow V(S_t) + \alpha(S_t)(R_t - V(S_t))$

where

$$\alpha(S_t) = \frac{1}{N(S_t)}$$

• Still we have to store the number of visits to each state: $N(S_t)$. Usually a **constant parameter** α in (0..1) is used:

$$V(S_t) \leftarrow V(S_t) + \alpha(R_t - V(S_t))$$

• Value of α

- ... with side effect of forgetting old episodes: as higher the value, higher the influence of recent experiences in the estimations
- Notice that:

$$V_n(S) = V_{n-1}(S) + \alpha(R_n - V_{n-1}(S)) = \alpha R_n + (1 - \alpha)V_{n-1}(S)$$

So,

$$V_{n}(S) = \alpha R_{n} + (1 - \alpha)V_{n-1}(S)$$

$$V_{n}(S) = \alpha R_{n} + (1 - \alpha)(\alpha R_{n-1} + (1 - \alpha)V_{n-2}(S))$$

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$$V_{n}(S) = \alpha R_{n} + \alpha (1 - \alpha)R_{n-1} + \alpha (1 - \alpha)^{2}R_{n-2} + \dots$$

$$V_{n}(S) = \alpha \sum_{i=1}^{n-1} \left[(1 - \alpha)^{i}R_{n-i} \right] + (1 - \alpha)^{n}R_{0}$$

- Trick used not only in Monte Carlo but on all methods
- ullet Choose lpha carefully. Remember

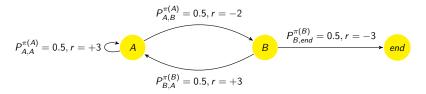
$$V_n(S) = \alpha \sum_{i=0}^{n-1} [(1-\alpha)^i R_{n-i}] + (1-\alpha)^n$$

• Usually α is low (0.01..0.2), but depends on the problem. Sometimes is good to forget old Long-term-Returns, for instance, when you change the policy!

Monte-Carlo policy evaluation: First visit vs. Every visit

- What to do when in a trial the state *s* appears several times? Two approaches: First-visit and Every-visit updates.
- Algorithm presented is called *First Visit MC* because it updates V(s) using return from the first visit.
- On the other hand, Every-visit MC updates states with all returns obtained

Monte-Carlo policy evaluation: First visit vs. Every visit



• Example: V for states A and B using both algorithms on these trials:

$$A \xrightarrow{+3} A \xrightarrow{-2} B \xrightarrow{+3} A \xrightarrow{-2} B \xrightarrow{-3} end$$

$$B \xrightarrow{+3} A \xrightarrow{-2} B \xrightarrow{-3} end$$

first-visit V(A) =
$$1/2(-1 - 5) = -3$$

V(B) = $1/2(-2 + -2) = -2$
every-visit V(A) = $1/4(-1 + -4 - 5 - 5) = -3.75$
V(B) = $1/4(-2 + -3 + -2 + -3) = -2.5$

ullet Both MC methods can be proved to converge to V^π

• Can we use the MC policy evaluation to learn a policy (like with PI)?

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- If we want to take the *greedy action*, like in PI, to improve the policy then we need the model!

$$\pi(s) = \operatorname*{arg\,max}_{s \in A} \sum_{s'} P_{ss'}^{a} \left[R(s') + \gamma V^{\pi}(s') \right]$$

- Can we use the MC policy evaluation to learn a policy (like with PI)?
- If we want to take the *greedy action*, like in PI, to improve the policy then we need the model!

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- **Solution**: estimate Q^{π} function instead of V^{π}
- Now we can *greedify* the policy without the model:

$$\pi(s) = \underset{a \in A}{\operatorname{arg max}} Q^{\pi}(s, a)$$

• Apply the *improvement-of-the-policy* idea to learn the optimal policy.

Caution! Wrong Monte Carlo policy control

```
Initialize \pi and Q randomly:
repeat
  Generate trial using \pi
  for each s, a in trial do
     R \leftarrow return following the first occurrence of s
     Q(s,a) \leftarrow Q(s,a) + \alpha(R - Q(s,a))
  end for
  for each s, a in trial do
     \pi(s) = \operatorname{arg\,max}_{a \in A} Q(s, a)
  end for
until true
```

- What's wrong?
 - ► Algorithm tries to implement asynchronous version of policy iteration... but remember... there states are selected for updating **randomly**.
 - Now states to be updated depend on the current policy, so we cannot guarantee convergence.
- New important concept: Exploration vs. Exploitation
 - ▶ All pairs (s,a) should have probability non-zero to be updated.
 - ► At same time, we want to evaluate the current policy
- Several ways to balance two concepts.

ϵ -greedy exploration

- Simplest idea for ensuring continual exploration
- All m actions are tried with non-zero probability
- ullet With probability $1-\epsilon$ choose the greedy action
- ullet With probability ϵ choose an action at random

$$\pi(a|s) = egin{cases} \epsilon/m + 1 - \epsilon, & ext{if } a = rg \max_{a' \in \mathcal{A}} Q(s, a') \ \epsilon/m, & ext{otherwise} \end{cases}$$

where $m = |\mathcal{A}(s)|$

Apply the *improvement-of-the-policy* idea to learn the optimal policy.
 Caution!

Monte Carlo policy control

```
Initialize \pi and Q randomly:
repeat
  Generate trial using \epsilon-greedy strategy on \pi
  for each s, a in trial do
     R \leftarrow return following the first occurrence of s
     Q(s,a) \leftarrow Q(s,a) + \alpha(R - Q(s,a))
  end for
  for each s, a in trial do
     \pi(s) = \arg \max_{a \in A} Q(s, a) // ties randomly broken
  end for
until true
```

- Ideally, exploration should not be constant during training.
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- It should be larger at the beginning and lower after a lot of experience is accumulated (why?)... but never disappear (why?)
- This requirement is asked in convergence proofs of most RL algorithms, f.i. in MC.
- In ϵ -greedy, this is implemented with variable ϵ starting from 1 and decreasing with number of experiences until a minimum ϵ value from which does not decrease further, f.i:

$$\epsilon = \max(1/(\alpha T), 0.1)$$

where T is the number of Trials done and α is a constant that controls decrease of exploration

- Another popular way to explore is using Softmax exploration or Gibb's exploration or Boltzman exploration.
- Idea is that probability depends on the value of actions, with bias of exploration towards more promising actions
- Softmax action selection methods grade action probabilities by estimated values

$$P(s,a) = \frac{e^{Q(s,a)/\tau}}{\sum_{a' \in A} e^{Q(s,a')/\tau}}$$

where parameter τ is called *temperature* and decreases with experience

• When τ is very large, all actions with roughly same probability of being selected. When τ is low, almost certainty of selecting the action with higher Q-value.

- A hot topic of research
- We want to explore efficiently the state space
- A lot of other more complex mechanisms based on criteria
 - ► Less explored state, action pairs
 - ► Higher changes in value of state action pair
 - ▶ Bases on recency of last exploration
 - Uncertainty on estimation of values
 - Error in an agent's ability to predict the consequence of action (curiosity)
 - ▶ ...

Monte Carlo policy control

```
Initialize \pi and Q randomly:
repeat
  Generate trial using exploration method based on \pi
  for each s, a in trial do
     R \leftarrow return following the first occurrence of s
     Q(s,a) \leftarrow Q(s,a) + \alpha(R - Q(s,a))
  end for
  for each s, a in trial do
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  end for
until true
```

Monte Carlo policy control

```
Initialize \pi and {\it Q} randomly:
```

repeat

Generate trial using exploration method on greedy policy derived from

Q values

for each s, a in trial do

$$R \leftarrow$$
 return following the first occurrence of s, a

$$Q(s,a) \leftarrow Q(s,a) + \alpha(R - Q(s,a))$$

end for

until false

Temporal Differences methods: Q-learning

Temporal Differences policy evaluation

- Monte-Carlo methods compute expectation of Long-term-Reward averaging the return of several trials.
- Average is done after termination of the trial.
- We saw in previous lecture that Bellman equation also allow to estimate expectation of Long-term-Reward

$$Q^{\pi}(s,a) = \mathbb{E}_{\pi}[R_t|S_t = s, A_t = a]$$

= $\mathbb{E}_{\pi}[r_{t+1} + \gamma Q^{\pi}(S_{t+1}, \pi(S_{t+1}))|S_t = s, A_t = a]$

• Computing expectations with world model:

$$Q^{\pi}(s,a) = \sum_{s'} P^{a}_{ss'} \left(r(s') + \gamma Q^{\pi}(s',\pi(s')) \right)$$

Temporal Differences policy evaluation

- How to get rid of the world-model?
- Q-value function and averaging, like in the case of MC

$$Q(S_t, a) \leftarrow Q(S_t, a) + \alpha(R_t(s_t) - Q(S_t, a))$$

• But now substitute R_t with Bellman equation:

$$Q(S_t, a) \leftarrow Q(S_t, a) + \alpha \left[r_{t+1} + \gamma Q(S_{t+1}, \pi(S_{t+1})) - Q(S_t, a) \right]$$

or

$$Q(S_t, a) \leftarrow (1 - \alpha)Q(S_t, a) + \alpha \left[r_{t+1} + \gamma Q(S_{t+1}, \pi(S_{t+1}))\right]$$

• This is called bootstrapping

Temporal Differences policy evaluation

Temporal Differences policy evaluation

```
Given \pi initialize Q randomly: 

repeat s \leftarrow initial state of episode 

repeat a \leftarrow \pi(s) Take action a and observe s' and r Q(s,a) \leftarrow Q(s,a) + \alpha \left(r + \gamma Q(s',\pi(s')) - Q(s,a)\right) s \leftarrow s' until s is terminal until convergence
```

MC and TD comparison

- ullet Goal: learn Q^{π} online from experience under policy π
- Incremental every-visit Monte-Carlo
 - ▶ Update value Q(s, a) toward actual return R_t

$$Q(s_t, a_t) \leftarrow Q(s_t, a_t) + \alpha(R_t - Q(s_t, a_t))$$

- Simplest temporal-difference learning algorithm: TD(0)
 - ▶ Update value $Q(s_t, a_t)$ toward estimated return $r_{t+1} + \gamma V(s_{t+1})$

$$Q(s_t, a_t) \leftarrow Q(s_t, a_t) + \alpha(r_{t+1} + \gamma V(s_{t+1}) - Q(s_t, a_t))$$

- Remember that $Q(s_t, \pi(s_t)) = V(s_t)$
- $r_{t+1} + \gamma V(s_{t+1})$ is called the TD target
- $\delta_t = r_{t+1} + \gamma V(s_{t+1}) Q(s_t, a_t)$ 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
 - ▶ 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

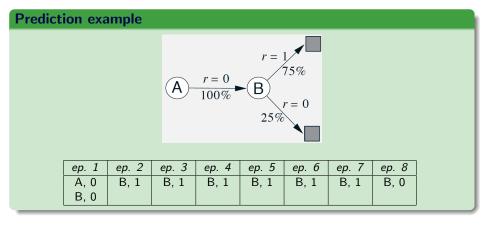
Bias/variance trade-off

- Return $R_t = r_{t+1} + r_{t+2} + \ldots + \gamma^{T-1} r_T$ is unbiased estimate of $V^{\pi}(S_t)$
- ullet True TD target $r_{t+1} + V^\pi(s_{t+1})$ is unbiased estimate of $V^\pi(s_t)$
- TD target $R_{t+1} + V(s_{t+1})$ 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

Temporal Differences policy evaluation

- On-line learning: evaluation is embedded in generation of the experience.
- It can be applied to non-episodic tasks
- You don't need to end episode to learn
- Like in MC you don't need the World model
- In practice faster: Takes profit on Markovian property

Temporal Differences policy evaluation



Find V(A) and V(B) from these episodes using TD(0) and MC.

MC vs TD control

- Temporal-difference (TD) learning has several advantages over Monte-Carlo (MC)
 - Lower variance
 - Online
 - Incomplete sequences
- Can we use it for policy learning?
- Natural idea: use TD instead of MC in our control loop
 - ▶ Apply TD to Q(S, A)
 - Use ϵ -greedy policy improvement
 - Update every time-step

Temporal Differences policy learning

Q-learning: Temporal Differences policy learning

```
Given \pi initialize Q randomly:
repeat
   s \leftarrow \text{initial state of episode}
   repeat
      Set a using f.i. \epsilon-greedy strategy on \pi
      Take action a and observe s' and r
      Q(s, a) \leftarrow Q(s, a) + \alpha (r + \gamma Q(s', \pi(s')) - Q(s, a))
      \pi(s) = \arg \max_{a \in A} Q(s, a) // \text{ ties randomly broken}
      s \leftarrow s'
   until s is terminal
until false
```

Temporal Differences extended

Temporal Differences extended

Bootstrapping in Bellman equation is done from next state:

$$V_{(1)}^{\pi}(s) = \mathbb{E}_{\pi}[R_t|S_t = s]$$

$$= \mathbb{E}_{\pi}[r_{t+1} + \gamma r_{t+2} + \gamma^2 r_{t+3} + \dots |S_t = s]$$

$$= \mathbb{E}_{\pi}[r_{t+1} + \gamma V^{\pi}(S_{t+1})|S_t = s]$$

• But we can obtain estimation from 2 steps in the future also:

$$V_{(2)}^{\pi}(s) = \mathbb{E}_{\pi}[R_{t}|S_{t} = s]$$

$$= \mathbb{E}_{\pi}[r_{t+1} + \gamma r_{t+2} + \gamma^{2} r_{t+3} + \dots |S_{t} = s]$$

$$= \mathbb{E}_{\pi}[r_{t+1} + \gamma r_{t+2} + \gamma^{2} (r_{t+3} + \dots))|S_{t} = s]$$

$$= \mathbb{E}_{\pi}[r_{t+1} + \gamma r_{t+2} + \gamma^{2} R_{t+2}|S_{t} = s]$$

$$= \mathbb{E}_{\pi}[r_{t+1} + \gamma r_{t+2} + \gamma^{2} V^{\pi}(S_{t+2})|S_{t} = s]$$

Temporal Differences extended

• In general we could extend that to the *n-steps estimator of long-term* reward.

$$\begin{aligned} V_{(n)}^{\pi}(s) &= \mathbb{E}_{\pi}[R_{t}|S_{t} = s] \\ &= \mathbb{E}_{\pi}[r_{t+1} + \gamma r_{t+2} + \ldots + \gamma^{n-1}r_{n} + \gamma^{n}r_{n+1} \dots |S_{t} = s] \\ &= \mathbb{E}_{\pi}\left[\sum_{k=0}^{n} \gamma^{k} r_{t+k+1} + \gamma^{n} V^{\pi}(S_{t+n})|S_{t} = s\right] \end{aligned}$$

Temporal Differences extended: n-step estimators

- All estimators of expectation are valid, but different bias and variance.
- Which one to use?
- Any of them is Ok at the end, but different learning speed with different value of *n*.
- Implementation of the algorithm is easy. For each episode
 - Execute n actions, keep rewards
 - Apply update

$$V^{\pi}(S_t) = \alpha V^{\pi}(S_t) + (1 - \alpha) \sum_{k=0}^{n} \gamma^k r_{t+k+1} + \gamma^n V^{\pi}(S_{t+n})$$

Temporal Differences n-steps policy evaluation

All store and access operations (for S_t and R_t) can take their index mod n

Temporal Differences n-steps policy evaluation

```
Given \pi and n, initialize Q randomly:
for each episode do
   s \leftarrow initial state of episode, and T \leftarrow \infty
   for t = 0, 1, 2... do
        if t < T then
           Take action a \leftarrow \pi(s) and observe and store s' and r
           If s' is terminal \leftarrow t+1
        end if
        \tau \leftarrow t - n + 1
       if \tau \geq 0 then
R \leftarrow \sum_{i=\tau+1}^{\min(\tau+n,T)} \gamma^{i-\tau-1} r_i
           If \tau + n < T then: R \leftarrow R + \gamma^n V(s_{\tau+n})
            Q(s,a) \leftarrow Q(s,a) + \alpha (R - Q(s,a))
        end if
        s \leftarrow s'
   end for
end for
```

Temporal Differences n-steps policy evaluation

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

Temporal Differences extended $TD(\lambda)$

- Another option. Instead of using one estimator, update using an average of them
- ullet For practical purposes, use a geometric average $(0 \le \lambda \le 1)$

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

Can be rewritten for episodes as:

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

- Unifies different algorithms:
 - ▶ When $\lambda = 0$ we have TD(0), the standard Q-learning method
 - ▶ When $\lambda = 1$ we have the standard MC method
- In general for other values of λ we use a smart incremental implementation using **eligibility traces** (chapter 12, Sutton book)

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- What?
- You will understand in next slides.
- Let's go now to discuss the concepts of on-line and off-line policy learning.

On-line vs. Off-line policy learning: Sarsa

Off-policy vs. On-policy learning

- When learning value functions of a policy, we sample *using the policy* to estimate them
- In Q-learning, the method tries to learn the value function of the optimal policy (V*) when in fact samples are obtained from different policy $(\epsilon$ -greedy policy)
- A subtle point with implications about the convergence of the algorithms to the optimal solution
- We'll do the following distinction:
 - On-policy learning: When learning the value function V^{π} of the current policy π
 - **On-policy learning:** When Learning the value function V^{π} using another policy π'

Off-policy vs. On-policy learning

- In this sense, Q-learning is an example of off-policy learning.
- Policy for which we learn values:

$$\pi^*(s) = \arg\max_{a} Q * (s, a)$$

• Sampling policy (ϵ -greedy policy):

$$\pi(a|s) = egin{cases} \epsilon/m + 1 - \epsilon, & ext{if } a = rg \max_{a' \in \mathcal{A}} Q(s, a') \ \epsilon/m, & ext{otherwise} \end{cases}$$

where $m = |\mathcal{A}(s)|$

Off-policy vs. On-policy learning

- Trivia. What about MC learning?
- It's on-policy or off-policy learning?

Updating action-value functions with SARSA

- Let's try to implement an on-policy learning version of Q-learning
- Sarsa: on-policy TD(0) learning
- In Q-learning, update equation was:

$$Q(s,a) \leftarrow Q(s,a) + \alpha(r + \gamma) \underbrace{\max_{a'} Q(s',a')}_{Q(s',a')} - Q(s,a))$$

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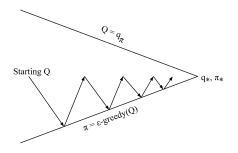
$$Q(s,a) \leftarrow Q(s,a) + \alpha(r + \gamma) \underbrace{\max_{a'} Q(s',a')}_{a'} - Q(s,a))$$

Update equation is sarsa:

$$Q(s,a) \leftarrow Q(s,a) + \alpha(r + \gamma) \qquad Q(s',a') \qquad -Q(s,a)$$

Now a is action selected

On-policy control with SARSA



Every time-step:

Policy evaluation with Sarsa, $Q pprox q_\pi$

Policy improvement with ϵ -greedy policy improvement.

SARSA algorithm for on-policy control

SARSA: on-line learning

```
Initialize Q(s, a), \forall s \in S, a \in A(s), arbitrarily, and
Q(terminal - state,) = 0
for each episode do
  Choose initial state s
  Choose a from s using policy derived from Q (e.g., \epsilon-greedy)
  for each step of episode do
     Execute action a, observe r, s'
     Choose a' from s' using policy derived from Q (e.g., \epsilon-greedy)
     Q(s, a) \leftarrow Q(s, a) + \alpha(R + \gamma Q(s', a') - Q(s, a))
     s \leftarrow s' : a \leftarrow a'
  end for
end for
```

SARSA algorithm for on-policy control

It can be proved that Sarsa converges to the optimal policy under the following conditions:

- Greedy in the Limit of Infinite Exploration (GLIE):
 - ► All state—action pairs are explored infinitely many times:

$$\lim_{t\to\infty} N_k(s,a) = \infty$$

▶ The policy converges on a greedy policy (f.i. ϵ decreases inversely proportional to the number of experiences)

$$\lim_{t \to \infty} \pi_t(a|s) = rg \max_{a'} Q^{\pi_t}(s, a')$$

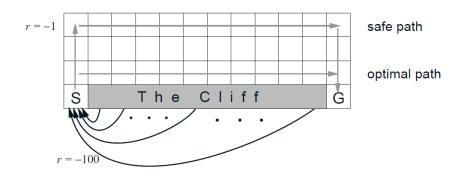
2 Robbins-Monro sequence of step-sizes α_t

$$\sum_{t=1}^{\infty} \alpha_t = \infty$$

$$\sum_{t=1}^{\infty} \alpha_t^2 < \infty$$

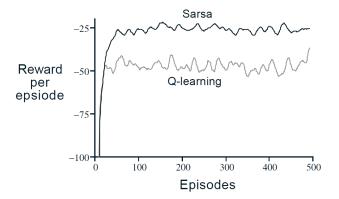
Practical differences btw Sarsa and Q-learning

Cliff-Walk example ($\epsilon = 0.1$):



Practical differences btw Sarsa and Q-learning

Cliff-Walk example: Reward during learning



As ϵ decreases, sarsa tends to Q-learning

Practical differences btw Sarsa and Q-learning

- In the cliff-walking task:
 - Q-learning: learns optimal policy along edge
 - ► Sarsa: learns a safe non-optimal policy away from edge
- ϵ -greedy algorithm
 - For $\epsilon = 0$ SARSA performs better online
 - For $\epsilon o 0$ gradually, both converge to optimal

Expected Sarsa

- In sarsa, we use the current policy to estimate returns
- However, we can do better: Expected Sarsa
- In sarsa, each episode uses one sample of action taken by the policy.
- ... but we know the policy probabilities to select one action (f.i. in ϵ -greedy procedure), so we can use it.
- Sarsa update:

$$Q(s, a) \leftarrow Q(s, a) + \alpha(r + \gamma Q(s', a') - Q(s, a))$$

Expected sarsa update:

$$Q(s,a) \leftarrow Q(s,a) + \alpha(r + \gamma \sum_{a} \pi(a'|s')Q(s',a') - Q(s,a))$$

• Same convergence guarantees and less variance than original Sarsa

Expected SARSA

Expected SARSA

```
Initialize Q(s, a), \forall s \in S, a \in A(s), arbitrarily, and
Q(terminal - state,) = 0
for each episode do
  Choose initial state s
  Choose a from s using policy derived from Q (e.g., \epsilon-greedy)
  for each step of episode do
     Execute action a, observe r, s'
     Choose a' from s' using policy derived from Q (e.g., \epsilon-greedy)
     Q(s,a) \leftarrow Q(s,a) + \alpha(r + \gamma \sum_{a} \pi(a'|s')Q(s',a') - Q(s,a))
     s \leftarrow s' : a \leftarrow a'
  end for
end for
```

- Return to Temporal Differences extended
- Very good to estimate values for a given policy
- More difficult to apply to control, because trace of states visited follow one policy but we are estimating another one
- Remember this conclusions?

- Return to Temporal Differences extended
- Very good to estimate values for a given policy
- More difficult to apply to control, because trace of states visited follow one policy but we are estimating another one
- Remember this conclusions?
- n-steps estimators and $TD(\lambda)$ can be easily implemented to control for Sarsa and Extended Sarsa, because they are on-policy learning methods.

$Sarsa(\lambda)$

```
Initialize Q(s, a) arbitrarily
loop
    e(s, a) = 0, for all s, a
    Initialize s, a
    repeat
         Take action a, observe r, s'
         Choose a' from s' using policy derived from Q (e.g., \epsilon-greedy)
         \delta \leftarrow r + \gamma Q(s', a') - Q(s, a)
         e(s, a) \leftarrow e(s, a) + 1
         for all s, a do
              Q(s, a) \leftarrow Q(s, a) + \alpha \delta e(s, a)
              e(s, a) \leftarrow \gamma \lambda e(s, a)
         end for
         s \leftarrow s' : a \leftarrow a':
    until s is terminal
end loop
```

Temporal Differences with Sarsa

ullet Benefits of temporal differences using larger n-step than $\mathsf{TD}(0)$



• In general, faster propagation of rewards and, so, faster learning.

- In off-policy learning is more difficult to implement n-steps methods
- However, n-steps still can be used in off-policy learning (f.i Q-learning). The trick is to use a dynamic n. When an exploratory action is taken then stop the trace of action from which to update Peng's $Q(\lambda)$.



General Off-policy learning

- Evaluate target policy $\pi(a|s)$ to compute $V^{\pi}(s)$ or $Q^{\pi}(s,a)$
- ... while following behavior policy $\mu(a|s)$

$$\{s_1, a_1, r_2, \ldots, s_T\} \sim \mu$$

- Why is this important?
 - ► Learn from observing humans or other agents
 - ▶ Re-use experience generated from old policies $\pi_1, \pi_2, \dots, \pi_{t-1}$
 - ► Learn about *optimal* policy while following *exploratory* policy
 - ▶ Learn about *multiple* policies while following *one* policy

Importance Sampling

 Estimate the expectation of a different distribution w.r.t. the distribution used to draw samples

$$\mathbb{E}_{x \sim p} [f(x)] = \sum_{x \sim p} p(x) f(x)$$

$$= \sum_{x \sim q} q(x) \frac{p(x)}{q(x)} f(x)$$

$$= \mathbb{E}_{x \sim q} \left[\frac{p(x)}{q(x)} f(x) \right]$$

$$\approx \frac{1}{T} \sum_{t=1}^{T} \frac{p(x^t)}{q(x^t)} f(x^t)$$

where data is sampled using q distribution. That means, we can estimate $\mathbb{E}_{x \sim p}[f(x)]$ using distribution q instead of p

Importance Sampling for off-policy Monte Carlo

- Application of IS to MC
- ullet Use returns generated from μ to evaluate π
- ullet Weight return R_t according to similarity between policies multiplying importance sampling corrections along whole episode

$$R_t^{\pi} = rac{\pi(a_t|s_t)}{\mu(a_t|s_t)} rac{\pi(a_{t+1}|s_{t+1})}{\mu(a_{t+1}|s_{t+1})} \dots rac{\pi(a_T|s_T)}{\mu(a_T|s_T)} R_t^{\mu}$$

Update value towards corrected return

$$Q(s_t, a_t) \leftarrow Q(s_t, a_t) + \alpha(R_t^{\pi} - Q(s_t, a_t))$$

where action are taken following μ

- Caution:
 - ▶ Cannot use if μ is zero where π is non–zero
 - Importance sampling can dramatically increase variance (choose μ wisely)

Importance Sampling for off-policy TD(0)

- Application of IS to Sarsa
- Off–Policy IS Sarsa comes from Bellman expectation equation for Q(s, a):

$$Q^{\pi}(s,a) = \mathbb{E}_{\pi} [r_{t+1} + \gamma Q^{\pi}(s_{t+1}, s_{a+1}) | s = s_t, a = a_t]$$

$$= r(s,a) + \gamma \sum_{s \in S} P_{s,a}^{s'} \sum_{a' \in A} \pi(a'|s') Q^{\pi}(s', a')$$

$$= r(s,a) + \gamma \sum_{s \in S} P_{s,a}^{s'} \sum_{a' \in A} \mu(a'|s') \frac{\pi(a'|s')}{\mu(a'|s')} Q^{\pi}(s', a')$$

$$= \mathbb{E}_{\mu} \left[r_{t+1} + \gamma \frac{\pi(a'|s')}{\mu(a'|s')} Q^{\pi}(s_{t+1}, s_{a+1}) | s = s_t, a = a_t \right]$$

Importance Sampling for off-policy TD(0)

• Off-policy update for Q-learning IS

$$Q^{\pi}(s,a) = Q^{\pi}(s,a) + \alpha \left[r + \gamma \frac{\pi(a'|s')}{\mu(a'|s')} Q^{\pi}(s',a') - Q^{\pi}(s,a) \right]$$

- Advantages over MC IS
 - Significantly lower variance because only one IS weight to multiply needed
 - ► Can change the policy at each time step!
- Like in MC IS, $\mu(a|s) > 0$ when $\pi(a|s) > 0$