Lecture 3: Tabular solution methods of RL

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Outline

- Introduction to reinforcement learning (RL)
- Tabular solution methods of RL
- Approximate solution methods of RL

Reference: Reinforcement Learning: An Introduction. <u>Richard S. Sutton</u> and <u>Andrew G. Barto</u>. Second Edition. MIT Press, Cambridge, MA, 2018

Tabular Solution Methods

Tabular Solution Methods

- Multi-Armed Bandits
- Finite Markov Decision Process
- Dynamic Programming
- Monte Carlo Methods
- Temporal Difference Learning

O Set up

- Action A_t at time t is $a \in \{a_1, a_2, ..., a_K\}$
- Obtain a random reward R_t following unknown distribution p_{A_t}

O Goal

Maximize the total reward over time, i.e.,

$$R_1 + R_2 + R_3 + \dots, R_n + \dots$$

as a consequence of

$$A_1, A_2, A_3, \ldots, A_n, \ldots$$

• Assuming independence, and remove randomness by expectation, it is sufficient to maximize

$$\mathbb{E}[R_1|A_1] + \mathbb{E}[R_2|A_2] + \mathbb{E}[R_3|A_3] + ..., \mathbb{E}[R_n|A_n] + ...$$

which is equivalent to maximizing $\mathbb{E}[R_t|A_t]$ by choosing A_t

O New goal

• Maximize $\mathbb{E}[R_t | A_t]$ by choosing A_t

O True value function

• $q_*(a) := \mathbb{E}[R_t | A_t = a]$ tells us how to choose the action A_t

$$A_t = \underset{a}{\operatorname{arg max}} q_*(a) = \underset{a}{\operatorname{arg max}} \mathbb{E}[R_t | A_t = a]$$

to maximize $\mathbb{E}[R_t|A_t]$

O Estimated value function

• Find $Q_t(a) \approx q_*(a)$ through interactions

$$A_0, R_0, A_1, R_1, \ldots, A_n, R_n, \ldots$$

New goal: find $Q_t(a) \approx q_*(a)$ through interactions

$$A_0, R_0, A_1, R_1, \dots, A_n, R_n, \dots$$

Action-value methods

O Estimation

$$Q_t(a) := \frac{\text{sum of rewards when a taken prior to } t}{\text{number of times } a \text{ taken prior to } t} = \frac{\sum_{i=1}^{t-1} R_i \cdot \mathbf{1}_{A_i = a}}{\sum_{i=1}^{t-1} \mathbf{1}_{A_i = a}}$$

O Greedy selection

$$A_t = \arg\max Q_t(a)$$

- Each action will be explored at least once
- "Local maximizer", if the rewards for arg max $q_*(a)$ happen to be smaller than they should be
- No exploration to escape "local maximizers"

New goal: find $Q_t(a) \approx q_*(a)$ through interactions

$$A_0, R_0, A_1, R_1, \dots, A_n, R_n, \dots$$

Action-value methods

O Estimation

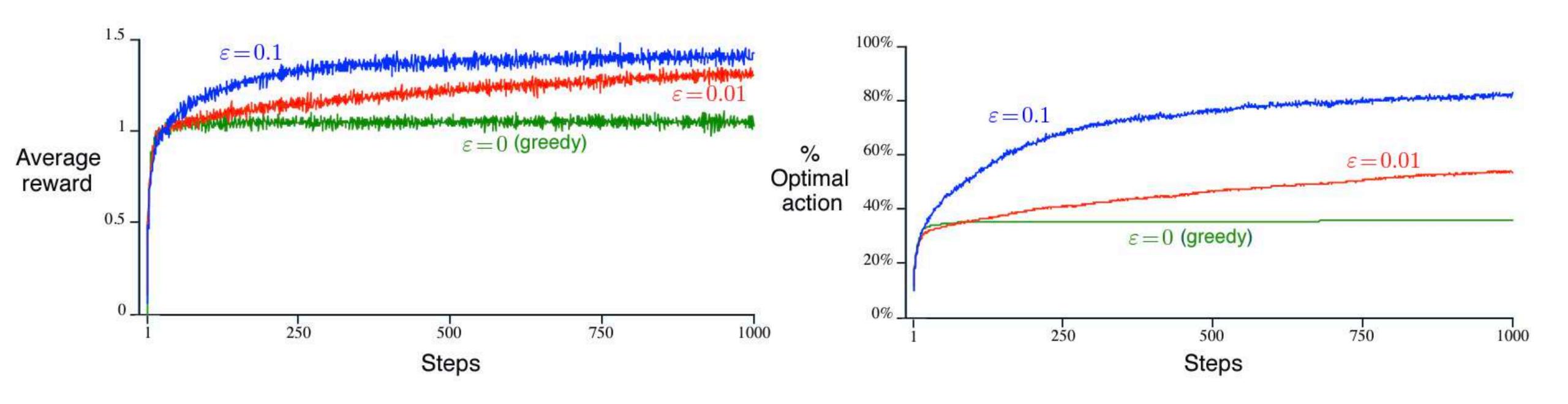
$$Q_t(a) := \frac{\text{sum of rewards when a taken prior to } t}{\text{number of times } a \text{ taken prior to } t} = \frac{\sum_{i=1}^{t-1} R_i \cdot \mathbf{1}_{A_i = a}}{\sum_{i=1}^{t-1} \mathbf{1}_{A_i = a}}$$

 \circ ε -greedy selection

$$A_{t} \sim \pi_{t}(a) = \begin{cases} 1 - \varepsilon + \frac{\varepsilon}{K}, & \text{if } a = \arg\max Q_{t}(a) \\ \frac{\varepsilon}{K}, & \text{otherwise} \end{cases}$$

- Enhance exploration via setting $\varepsilon > 0$ to avoid "local maximizer"
- $\pi_t(a)$ converges to a Dirac delta distribution center at arg max $Q_t(a)$ as $\varepsilon \to 0$
- $\varepsilon \to 0$ when $t \to \infty$ to reduce redundant exploration

Action Value Methods



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O Action-value methods

Direct estimation

$$Q_t(a) := \frac{\text{sum of rewards when a taken prior to } t}{\text{number of times } a \text{ taken prior to } t} = \frac{\sum_{i=1}^{t-1} R_i \cdot \mathbf{1}_{A_i = a}}{\sum_{i=1}^{t-1} \mathbf{1}_{A_i = a}}$$

Incremental implementation for computational efficiency

Suppose we only have one action, then

$$Q_{n+1} = \frac{1}{n} \sum_{i=1}^{n} R_i = \frac{1}{n} \left(R_n + \sum_{i=1}^{n-1} R_i \right) = \frac{1}{n} \left(R_n + (n-1) \frac{1}{n-1} \sum_{i=1}^{n-1} R_i \right)$$
$$= \frac{1}{n} \left(R_n + (n-1)Q_n \right) = \frac{1}{n} \left(R_n + nQ_n - Q_n \right) = Q_n + \frac{1}{n} \left[R_n - Q_n \right]$$

O Action-value methods

Incremental implementation when we have only one action

$$Q_{n+1} = Q_n + \frac{1}{n} [R_n - Q_n]$$

• General rule true for other cases

$$NewEstimate \leftarrow OldEstimate + StepSize \left[Target - OldEstimate\right]$$

$$Q_t(a) = Q_t(a) + \frac{1}{N_t(a)} [R_t - Q_t(a)]$$

O Action-value methods

Incremental update rule in general

$$NewEstimate \leftarrow OldEstimate + StepSize \left[Target - OldEstimate\right]$$

A simple bandit algorithm

```
Initialize, for a=1 to k:
Q(a) \leftarrow 0
N(a) \leftarrow 0
Loop forever:
A \leftarrow \begin{cases} \operatorname{argmax}_a Q(a) & \text{with probability } 1-\varepsilon \\ \operatorname{a random action} & \text{with probability } \varepsilon \end{cases}
R \leftarrow bandit(A)
N(A) \leftarrow N(A) + 1
Q(A) \leftarrow Q(A) + \frac{1}{N(A)} [R - Q(A)]
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O Action-value methods

• Incremental update rule for non-stationary reward probability with StepSize as a constant α

$$NewEstimate \leftarrow OldEstimate + StepSize \left[Target - OldEstimate \right]$$

• Resulting in a weighted average of past rewards emphasizing recent rewards

$$Q_{n+1} = Q_n + \alpha \Big[R_n - Q_n \Big]$$

$$= \alpha R_n + (1 - \alpha) Q_n$$

$$= \alpha R_n + (1 - \alpha) [\alpha R_{n-1} + (1 - \alpha) Q_{n-1}]$$

$$= \alpha R_n + (1 - \alpha) \alpha R_{n-1} + (1 - \alpha)^2 Q_{n-1}$$

$$= \alpha R_n + (1 - \alpha) \alpha R_{n-1} + (1 - \alpha)^2 \alpha R_{n-2} + \cdots + (1 - \alpha)^{n-1} \alpha R_1 + (1 - \alpha)^n Q_1$$

$$= (1 - \alpha)^n Q_1 + \sum_{i=1}^n \alpha (1 - \alpha)^{n-i} R_i.$$

• More general, time dependent StepSize $\alpha_t(a)$

Action-value methods

O ε -greedy selection: no preference for non-greedy actions

$$A_{t} \sim \pi_{t}(a) = \begin{cases} 1 - \varepsilon + \frac{\varepsilon}{K}, & \text{if } a = \arg\max Q_{t}(a) \\ \frac{\varepsilon}{K}, & \text{otherwise} \end{cases}$$

O Upper-Confidence Bound selection: preference based on uncertainty

$$A_t = \arg\max \left[Q_t(a) + c\sqrt{\frac{\ln t}{N_t(a)}} \right]$$

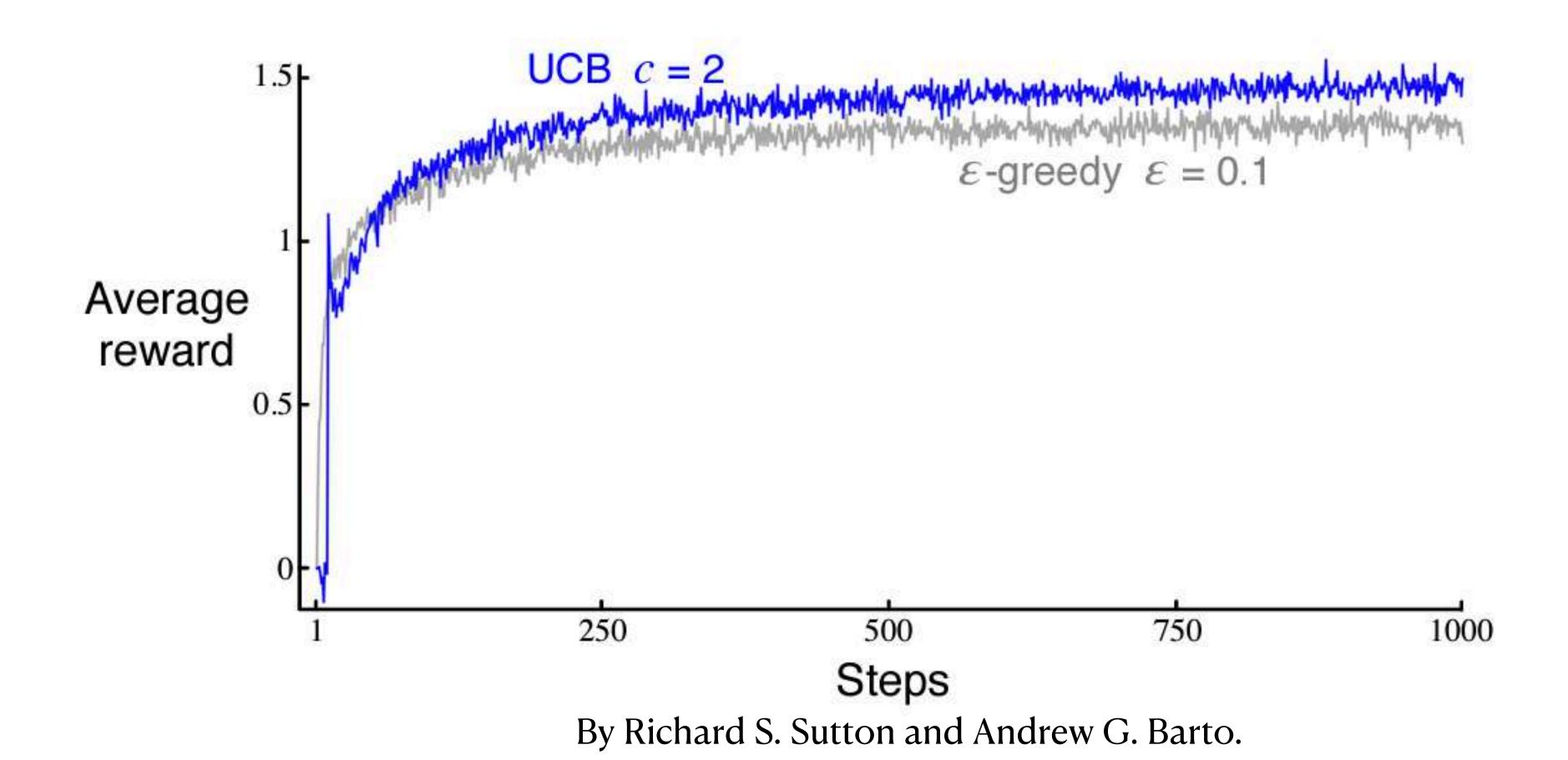
Action-value methods

O Upper-Confidence Bound selection: preference based on uncertainty

$$A_t = \arg\max \left[Q_t(a) + c\sqrt{\frac{\ln t}{N_t(a)}} \right]$$

- Enhance exploration via adding $c\sqrt{\frac{\ln t}{N_t(a)}}$ to avoid "local maximize"
- Smaller $N_t(A_t)$, larger uncertainty on A_t , more preference to use $c\sqrt{\frac{\ln t}{N_t(A_t)}}$ to choose A_t
- Larger $N_t(A_t)$, smaller uncertainty on A_t , more preference to use $Q_t(A_t)$ to choose A_t
- $c\sqrt{\frac{\ln t}{N_t(a)}} \to 0$ when $t \to \infty$ to reduce redundant exploration

ε -greedy selection v.s. Upper-Confidence Bound selection



O Action-value methods:

- Maximize the expected reward $\mathbb{E}[R_t|A_t]$ by choosing A_t appropriately
- Construct $Q_t(a)$ converging to $q_*(a)$ to choose A_t

O Policy-gradient methods:

- A probability distribution $\pi_t(a)$ as a policy to determine A_t without estimating Q_t
- $\pi_*(a) := \delta_{\arg\max_{\bar{a}} q_*(\bar{a})}(a)$ is the optimal policy to determine A_t
- Update $\pi_t(a)$ to approach $\pi_*(a)$
- The updating rule is equivalent to stochastic gradient accent to maximize $\mathbb{E}_{A_t \sim \pi_t} \left[\mathbb{E}[R_t | A_t] \right]$ as a functional of $\pi_t(a)$

O Policy-gradient methods

- Find $\pi_t(a) \to \pi_*(a)$ that maximizes $\mathbb{E}_{A_t \sim \pi_t} \left[\mathbb{E}[R_t | A_t] \right]$
- Action selection and modeling of $\pi_t(a)$

$$\Pr\{A_t = a\} := \frac{e^{H_t(a)}}{\sum_{b=1}^K e^{H_t(b)}} := \pi_t(a)$$

where $H_t(a)$ is the preference for choosing $A_t = a$ at time t

- $\mathbb{E}_{A_t \sim \pi_t} \left[\mathbb{E}[R_t | A_t] \right]$ is a functional of $H_t(a)$
- Find $H_t(a) \to H_*(a)$ that maximizes $\mathbb{E}_{A_t \sim \pi_t} \left[\mathbb{E}[R_t | A_t] \right]$

O Policy-gradient methods

- Find $H_t(a) \to H_*(a)$ that maximizes $\mathbb{E}_{A_t \sim \pi_t} \left[\mathbb{E}[R_t | A_t] \right]$
- Update $H_t(a)$ in a sense of the stochastic gradient accent for $\mathbb{E}_{A_t \sim \pi_t} \left[\mathbb{E}[R_t | A_t] \right]$ via

$$H_{t+1}(A_t) := H_t(A_t) + \alpha(R_t - \bar{R}_t)(1 - \pi_t(A_t))$$

and

$$H_{t+1}(a) := H_t(a) - \alpha (R_t - \bar{R}_t) \pi_t(a)$$
 for $a \neq A_t$

where \bar{R}_t is the average reward up to but not including time t.

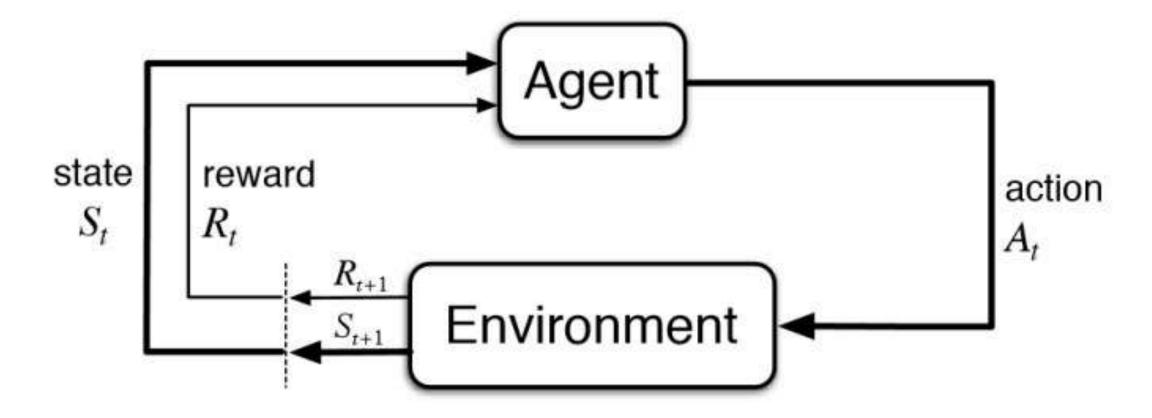
• Then as $t \to \infty$,

$$H_t \to H_* \Leftrightarrow \pi_t \to \pi_*$$

that maximizes $\mathbb{E}_{A_t \sim \pi_t} \left[\mathbb{E}[R_t | A_t] \right]$

Summary

- O Setup: Single state and K actions
- O Methods:
 - Value function method: find $Q_t(a) \to q_*(a)$ to maximize $\mathbb{E}[R_t|A_t]$ by choosing $A_t \to \arg\max_a q_*(a)$
 - Policy-gradient method: find $H_t \to H_* \Leftrightarrow \pi_t \to \pi_*$ that maximizes $\mathbb{E}_{A_t \sim \pi_t} \left[\mathbb{E}[R_t | A_t] \right]$
- O Looking for $\pi_*(a) := \delta_{\arg\max_{\bar{a}} q_*(\bar{a})}(a)$ is equivalent to identifying $\arg\max_{\bar{a}} q_*(a)$
- O Initialization and hyper-parameters are important!



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- Goal: Learn how to take actions in order to maximize reward via iteraction
- Components: environment in a state $S_t \in \mathcal{S}$ returning a reward $R_t \in \mathcal{R}$, agent taking actions $A_t \in \mathcal{A}$
- Interaction: $S_0, A_0, R_0, S_1, A_1, R_1, S_2, A_2, R_2, \dots$
- Markov decision process with a probability distribution:

$$p(s', r | s, a) := \Pr\{S_t = s', R_t = r | S_{t-1} = s, A_{t-1} = a\}$$

- Interaction: $S_0, A_0, R_0, S_1, A_1, R_1, S_2, A_2, R_2, \dots$
- The probability distribution p defines the MDP and is the most basic unknown target:

$$p(s', r | s, a) := \Pr\{S_t = s', R_t = r | S_{t-1} = s, A_{t-1} = a\}$$

• State-transition probability:

$$p(s'|s,a) := \Pr\{S_t = s'|S_{t-1} = s, A_{t-1} = a\} = \sum_{r \in \mathcal{R}} p(s',r|s,a)$$

• Expected rewards for state-action pairs:

$$r(s, a) := \mathbf{E}\{R_t | S_{t-1} = s, A_{t-1} = a\} = \sum_{r \in \mathcal{R}} r \sum_{s' \in \mathcal{S}} p(s', r | s, a)$$

• Expected rewards for state-action-next-state triples:

$$r(s, a, s') := \mathbf{E}\{R_t | S_{t-1} = s, A_{t-1} = a, S_t = s'\} = \sum_{r \in \mathcal{R}} r \frac{p(s', r | s, a)}{p(s' | s, a)}$$

- Interaction: $S_0, A_0, R_0, S_1, A_1, R_1, S_2, A_2, R_2, \dots$
- Goal: maximize the expected discounted return

$$G_t := R_{t+1} + \gamma R_{t+2} + \gamma^2 R_{t+3} + \dots = \sum_{k=0}^{\infty} \gamma^k R_{t+k+1} = R_{t+1} + \gamma G_{t+1}$$

- γ is a parameter in [0,1] called the discount rate
- R_t and R_{t+1} both depends on S_k for $k \le t-1$
- Application types:
 - 1) episodes with a termination time T and a state state s_+
 - 2) continuing tasks without termination

- Interaction: $S_0, A_0, R_0, S_1, A_1, R_1, S_2, A_2, R_2, \dots$
- Question: how to choose actions to maximize the expected discounted return G_t ?

$$G_t := R_{t+1} + \gamma R_{t+2} + \gamma^2 R_{t+3} + \dots = \sum_{k=0}^{\infty} \gamma^k R_{t+k+1} = R_{t+1} + \gamma G_{t+1}$$

- What we have learned in the case of a single state?
 - Value function method: find $Q_t(a) \to q_*(a)$ to maximize $\mathbb{E}[R_t|A_t]$ by choosing $A_t \to \arg\max_a q_*(a)$
 - Policy-gradient method: find $H_t \to H_* \Leftrightarrow \pi_t \to \pi_*$ that maximizes $\mathbb{E}_{A_t \sim \pi_t} \left[\mathbb{E}[R_t | A_t] \right]$
 - Consider $\mathbb{E}[R_t|A_t]$ instead of G_t because R_t 's are independent when A_t 's are given
- For MDP, we can also use policy and value function

- Interaction: $S_0, A_0, R_0, S_1, A_1, R_1, S_2, A_2, R_2, \dots$
- Question: how to choose actions to maximize the expected discounted return G_t ?

$$G_t := R_{t+1} + \gamma R_{t+2} + \gamma^2 R_{t+3} + \dots = \sum_{k=0}^{\infty} \gamma^k R_{t+k+1} = R_{t+1} + \gamma G_{t+1}$$

- Policy in MDP: a conditional probability distribution $\pi(a \mid s)$ for $A_t = a$ and $S_t = s$
- Value functions in MDP following the policy π
 - Value function of states: $v_{\pi}(s) := \mathbb{E}_{\pi}[G_t | S_t = s] = \mathbb{E}_{\pi}\left[\sum_{k=0}^{\infty} \gamma^k R_{t+k+1} \middle| S_t = s\right]$
 - Value function of state-action pair:

$$q_{\pi}(s, a) := \mathbb{E}_{\pi}[G_t | S_t = s, A_t = a] = \mathbb{E}_{\pi} \left[\sum_{k=0}^{\infty} \gamma^k R_{t+k+1} \middle| S_t = s, A_t = a \right]$$

• Bellman equation for the policy π

$$\begin{aligned} v_{\pi}(s) &:= \mathbb{E}_{\pi}[G_{t} \mid S_{t} = s] \\ &= \mathbb{E}_{\pi}[R_{t+1} + \gamma G_{t+1} \mid S_{t} = s] \\ &= \sum_{a} \pi(a \mid s) \sum_{s'} \sum_{r} p(s', r \mid s, a) \left[r + \gamma \mathbb{E}_{\pi}[G_{t+1} \mid S_{t+1} = s'] \right] \\ &= \sum_{a} \pi(a \mid s) \sum_{s'} \sum_{r} p(s', r \mid s, a) \left[r + \gamma v_{\pi}(s') \right] \end{aligned}$$
for all $s \in \mathcal{S}$

- $v_{\pi}(s)$ is the unique solution (i.e., fixed point) of its Bellman equation
- Given π , a fixed point iteration of the Bellman equation returns $v_{\pi}(s)$

- Interaction: $S_0, A_0, R_0, S_1, A_1, R_1, S_2, A_2, R_2, \dots$
- Question: how to choose actions to maximize the expected discounted return G_t ?
- Policy: $\pi(a \mid s)$ and Value functions following the policy π : $v_{\pi}(s)$ and $q_{\pi}(s, a)$
- How to compare policies?

$$\pi \ge \pi'$$
 if and only if $v_{\pi}(s) \ge v_{\pi'}(s)$ for all $s \in \mathcal{S}$

- Any optimal policy π_* ?
 - $\max_{\pi} v_{\pi}(s)$ for all $s \in \mathcal{S}$ share the same optimizer π_*

choose $\pi_*(a \mid s)$ to be the best probability for each fixed s

• π_* may not be unique

- Interaction: $S_0, A_0, R_0, S_1, A_1, R_1, S_2, A_2, R_2, \dots$
- Question: how to choose actions to maximize the expected discounted return G_t ?
- Policy: $\pi(a \mid s)$ and Value functions following the policy π : $v_{\pi}(s)$ and $q_{\pi}(s, a)$
- How to compare policies?

$$\pi \ge \pi'$$
 if and only if $v_{\pi}(s) \ge v_{\pi'}(s)$ for all $s \in \mathcal{S}$

- The optimal policy π_* share the same
 - optimal state value function $v_*(s) := \max_{\pi} v_{\pi}(s)$ for all $s \in \mathcal{S}$
 - optimal state-action value function $q_*(s,a) := \max_{\pi} q_{\pi}(s,a)$ for all $s \in \mathcal{S}$ and $a \in \mathcal{A}(s)$

The optimal value functions

- $v_*(s) := \max_{\pi} v_{\pi}(s)$ and $q_*(s, a) := \max_{\pi} q_{\pi}(s, a)$ determine each other
- $q_*(s, a) := \max_{\pi} \mathbb{E}_{\pi}[G_t | S_t = s, A_t = a]$ $= \mathbb{E}_{\pi_*}[R_{t+1} + \gamma G_{t+1} | S_t = s, A_t = a]$ $= \mathbb{E}_{\pi_*}[R_{t+1} + \gamma v_*(S_{t+1}) | S_t = s, A_t = a]$
- Furthermore, $v_*(S_{t+1}) = \max_{a'} q_*(s', a')$
- Hence, $q_*(s, a) = \mathbb{E}_{\pi_*}[R_{t+1} + \gamma \max_{a'} q_*(s', a') | S_t = s, A_t = a]$ $= \sum_{s',r} p(s', r | s, a)[r + \gamma \max_{a'} q_*(s', a')]$
- The above is the Bellman equation of π_* for computing $q_*(s,a)$ via a fixed point iteration

The optimal value functions

•
$$v_*(s) := \max_{\pi} v_{\pi}(s)$$
 and $q_*(s, a) := \max_{\pi} q_{\pi}(s, a)$ determine each other

•
$$v_*(s) = \max_{a \in \mathcal{A}(s)} q_{\pi_*}(s, a)$$

$$= \max_{a \in \mathcal{A}(s)} \mathbb{E}_{\pi_*}[G_t | S_t = s, A_t = a]$$

$$= \max_{a \in \mathcal{A}(s)} \mathbb{E}_{\pi_*}[R_{t+1} + \gamma G_{t+1} | S_t = s, A_t = a]$$

$$= \max_{a \in \mathcal{A}(s)} \mathbb{E}_{\pi_*}[R_{t+1} + \gamma v_*(S_{t+1}) | S_t = s, A_t = a]$$

$$= \max_{a \in \mathcal{A}(s)} \sum_{s', r} p(s', r | s, a)[r + \gamma v_*(s')]$$

• The above is the Bellman equation of π_* for computing $v_*(s)$ via a fixed point iteration

Summary

- Introduced MDP
- Introduced policy
- Introduced value functions
- The optimal policy is associated with the optimal value functions
- Identify optimal value functions via fixed point iterations of the Bellman equation
- Derive ε -greedy policy using value functions,

e.g.,
$$A_t = \begin{cases} \arg\max_a q_*(S_t, a) & \text{with probability } 1 - \varepsilon \\ \text{uniform random action with probability } \varepsilon \end{cases}$$

Model-free method

Overview

- Model-based method: p(s', r | s, a) is known
- Expense computation
- Interesting theoretically
- Help us to understand some model-free methods
- Assume termination state s_+ and let $\mathcal{S}^+ := s_+ \cup \mathcal{S}$
- The environment is assumed to be a finite MDP for simplicity here

Computation motivation: optimal value function evaluation

- Model-based method: p(s', r | s, a) is known
- The optimal value function can derive the optimal policy
- The optimal value function satisfies the Bellman optimality equation:

$$v_*(s) = \max_{a} \mathbb{E}[R_{t+1} + \gamma v_*(S_{t+1}) | S_t = s, A_t = a]$$

$$= \max_{a} \sum_{s',r} p(s', r | s, a)[r + \gamma v_*(s')]$$

or

$$q_*(s, a) = \mathbb{E}[R_{t+1} + \gamma \max_{a'} q_*(S_{t+1}, a') | S_t = s, A_t = a]$$

$$= \sum_{s', r} p(s', r | s, a) [r + \gamma \max_{a'} q_*(s', a')]$$

for all $s \in \mathcal{S}$, $a \in \mathcal{A}(s)$, and $s' \in \mathcal{S}^+$

• Use the above equations to derive iterative methods to compute the optimal value functions

Computation Step 1: policy evaluation

- Model-based method: p(s', r | s, a) is known
- Goal: given a policy π , evaluate its value in v_{π}
- Observation:

$$v_{\pi}(s) := \mathbb{E}_{\pi}[G_{t} | S_{t} = s]$$

$$:= \mathbb{E}_{\pi}[R_{t+1} + \gamma G_{t+1} | S_{t} = s]$$

$$= \mathbb{E}_{\pi}[R_{t+1} + \gamma v_{\pi}(S_{t+1}) | S_{t} = s]$$

$$= \sum_{a} \pi(a | s) \sum_{s',r} p(s', r | s, a) [r + \gamma v_{\pi}(s')]$$

• Iterative procedure for the fixed point v_{π} ; then $v_k \to v_{\pi}$

$$v_{k+1}(s) = \mathbb{E}_{\pi}[R_{t+1} + \gamma v_k(S_{t+1}) | S_t = s]$$

$$= \sum_{a} \pi(a | s) \sum_{s',r} p(s', r | s, a) [r + \gamma v_k(s')]$$

Computation Step 1: policy evaluation

Iterative Policy Evaluation, for estimating $V \approx v_{\pi}$

```
Input \pi, the policy to be evaluated
Algorithm parameter: a small threshold \theta > 0 determining accuracy of estimation
Initialize V(s), for all s \in \mathbb{S}^+, arbitrarily except that V(terminal) = 0
```

Loop:

 $\Delta \leftarrow 0$

$$\begin{aligned} & \text{Loop for each } s \in \mathbb{S}: \\ & v \leftarrow V(s) \\ & V(s) \leftarrow \sum_{a} \pi(a|s) \sum_{s',r} p(s',r|s,a) \big[r + \gamma V(s') \big] \\ & \Delta \leftarrow \max(\Delta,|v-V(s)|) \end{aligned}$$
 until $\Delta < \theta$

Computation Step 2: policy improvement

- Goal: improve a policy π based on its value in v_{π}
- Observation: π' below is better than or as good as π

$$\pi'(s) := \arg \max_{a} q_{\pi}(s, a)$$

$$= \arg \max_{a} \mathbb{E}[R_{t+1} + \gamma v_{\pi}(S_{t+1}) | S_{t} = s, A_{t} = a]$$

$$= \arg \max_{a} \sum_{s', r} p(s', r | s, a) [r + \gamma v_{\pi}(s')]$$

- Iterative procedure: use π' as a new improved policy
- When π' is as good as π , then $\pi' = \pi = \pi_*$, because ν_* is the unique fixed point of the Bellman equation

Complete algorithm

Policy Iteration (using iterative policy evaluation) for estimating $\pi \approx \pi_*$

1. Initialization

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

2. Policy Evaluation

Loop:

$$\Delta \leftarrow 0$$

Loop for each $s \in S$:

$$v \leftarrow V(s)$$

$$V(s) \leftarrow \sum_{s',r} p(s',r|s,\pi(s)) [r + \gamma V(s')]$$

$$\Delta \leftarrow \max(\Delta, |v - V(s)|)$$

until $\Delta < \theta$ (a small positive number determining the accuracy of estimation)

3. Policy Improvement

$$policy\text{-}stable \leftarrow true$$

For each $s \in S$:

$$old\text{-}action \leftarrow \pi(s)$$

$$\pi(s) \leftarrow \operatorname{arg\,max}_a \sum_{s',r} p(s',r|s,a) [r + \gamma V(s')]$$

If $old\text{-}action \neq \pi(s)$, then $policy\text{-}stable \leftarrow false$

If policy-stable, then stop and return $V \approx v_*$ and $\pi \approx \pi_*$; else go to 2

Complete algorithm

Policy Iteration (using iterative policy evaluation) for estimating $\pi \approx \pi_*$

1. Initialization

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

2. Policy Evaluation

 $\Delta \leftarrow 0$

Loop:

Greedy evaluation:
$$V(s) \leftarrow \max_{a} \sum_{s',r} p(s',r|s,a)[r+\gamma V(s')]$$

Loop for each $s \in S$:

$$v \leftarrow V(s)$$

$$V(s) \leftarrow \sum_{s',r} p(s',r|s,\pi(s)) [r + \gamma V(s')]$$

$$\Delta \leftarrow \max(\Delta, |v - V(s)|)$$

until $\Delta < \theta$ (a small positive number determining the accuracy of estimation)

3. Policy Improvement $policy\text{-}stable \leftarrow true$

Greedy policy:
$$\pi(s) = \arg\max_{a} \sum_{s',r} p(s',r|s,a)[r+\gamma V(s')]$$

For each $s \in S$:

$$old\text{-}action \leftarrow \pi(s)$$

$$\pi(s) \leftarrow \operatorname{arg\,max}_a \sum_{s',r} p(s',r|s,a) [r + \gamma V(s')]$$

If $old\text{-}action \neq \pi(s)$, then $policy\text{-}stable \leftarrow false$

If policy-stable, then stop and return $V \approx v_*$ and $\pi \approx \pi_*$; else go to 2

Complete algorithm: fixed point iteration via the Bellman optimality equation

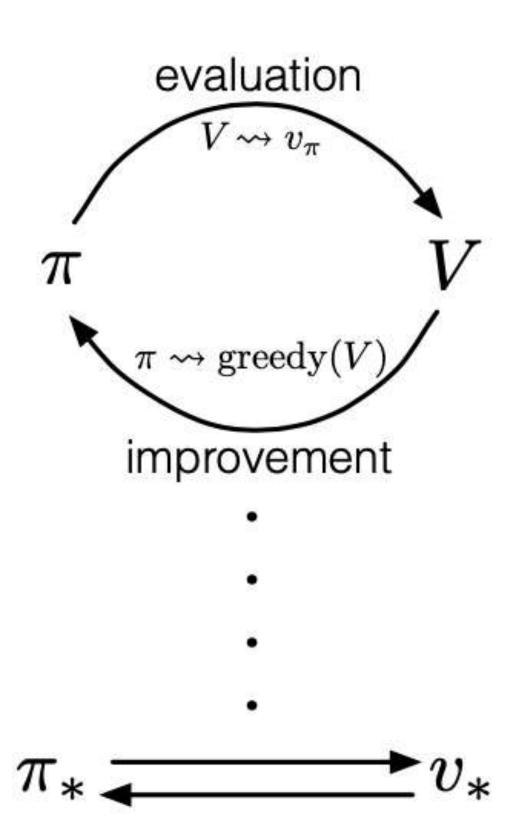
Generalized Policy Iteration

Observation

- Policy iteration for the optimal policy two simultaneous, interacting processes; each process takes many iterations
- Value iteration for the optimal policy two simultaneous, interacting processes; each process takes one step

Question

• Can we make it more flexible?



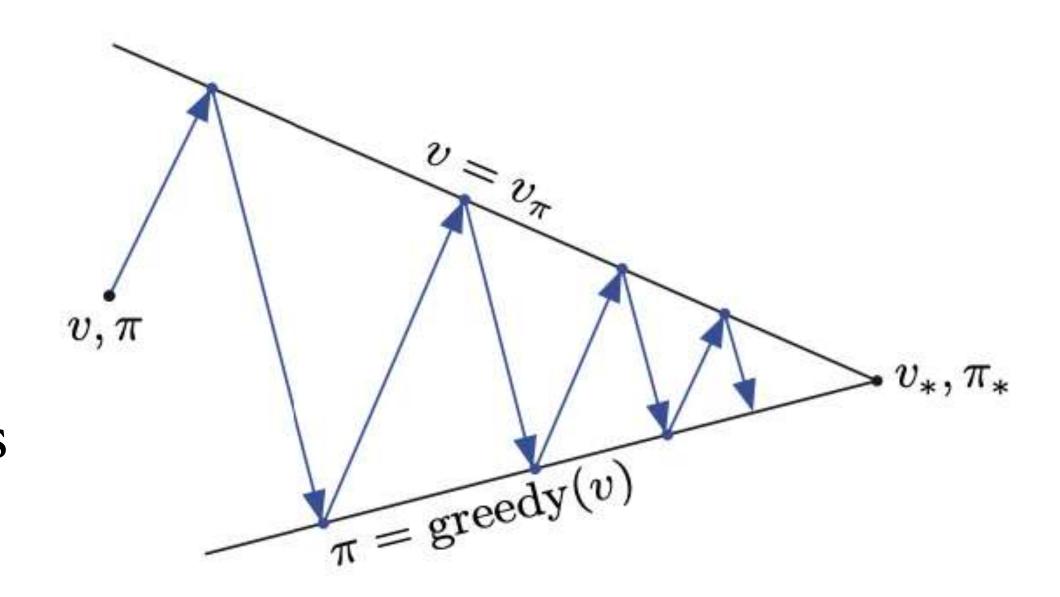
Generalized Policy Iteration

Question

• Can we make it more flexible?

Answer

- Yes
- Randomly or alternatively choose a subset of states to carry out these two processes
- O(1) steps in each process
- This is called the generalized policy iteration (GPI)



Summary

- Model-based method
- Introduced policy iteration method to compute the optimal policy
- Introduced value iteration method to compute the optimal policy
- Generalized policy iteration (GPI)
- Expensive computation

Overview

- Model-free method: p(s', r | s, a) is unknown
- Assume samples from p(s', r | s, a) are available
- Cheaper computation
- Offline episode-by-episode update instead of online step-by-step update
- Interesting in real applications
- The environment is assumed to be a finite MDP for simplicity here

- Basic idea: $\mathbb{E}[f(x)] \approx E_n := \frac{1}{n} \sum_i f(x_i)$
- Law of large number: $E_n \to \mathbb{E}[f(x)]$ as $n \to \infty$
- Incremental implementation: $E_{n+1} = \frac{nE_n + f(x_{n+1})}{n+1}$

Algorithms to be discussed

- O Policy value evaluation
 - State value function $v_{\pi}(s)$ can help to determine a good policy if model is known, e.g., model-based method DP
 - State action value function $q_{\pi}(s,a)$ directly help to determine policies without a model, e.g., model-free method MC
- O Policy improvement

First visit MC for policy evaluation: average of the returns following first visits to s

```
First-visit MC prediction, for estimating V \approx v_{\pi}
Input: a policy \pi to be evaluated
Initialize:
    V(s) \in \mathbb{R}, arbitrarily, for all s \in S
    Returns(s) \leftarrow \text{an empty list, for all } s \in S
Loop forever (for each episode):
    Generate an episode following \pi: 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 to Returns(S_t)
              V(S_t) \leftarrow \operatorname{average}(Returns(S_t))
```

Every visit MC for policy evaluation: average of the returns following every visit to s

Every

```
First visit MC prediction, for estimating V \approx v_{\pi}
Input: a policy \pi to be evaluated
Initialize:
     V(s) \in \mathbb{R}, arbitrarily, for all s \in S
     Returns(s) \leftarrow \text{an empty list, for all } s \in S
Loop forever (for each episode):
     Generate an episode following \pi: 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 to Returns(S_t)
              V(S_t) \leftarrow \operatorname{average}(Returns(S_t))
```

First visit MC for state-action evaluation: average of the returns following first visits to (s,a)

```
First-visit MC prediction, for estimating V \approx v_{\pi}
                                                                 q(s,a) \approx q_{\pi}(s,a)
Input: a policy \pi to be evaluated
Initialize:
     V(s) \in \mathbb{R}, arbitrarily, for all s \in \mathcal{S} q(s,a) \in \mathbb{R}, arbitrarily, for all s \in \mathcal{S} and a \in \mathcal{A}(s)
    Returns(s) \leftarrow an empty list, for all s \in \mathcal{S} Returns(s, a) \leftarrow an empty list, for all s \in \mathcal{S} and a \in \mathcal{A}(s)
Loop forever (for each episode):
     Generate an episode following \pi: 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, A_t) appears in (S_0, A_0), ..., (S_{t-1}, A_{t-1})
         Unless S_t appears in S_0, S_1, \ldots, S_{t-1}:
             Append G to Returns(S_t) Append G to Returns(S_t, A_t)
              -V(S_t) \leftarrow \operatorname{average}(Returns(S_t))
                                                                                  average(Returns(S_t, A_t))
                                                                  q(S_t, A_t) \leftarrow
```

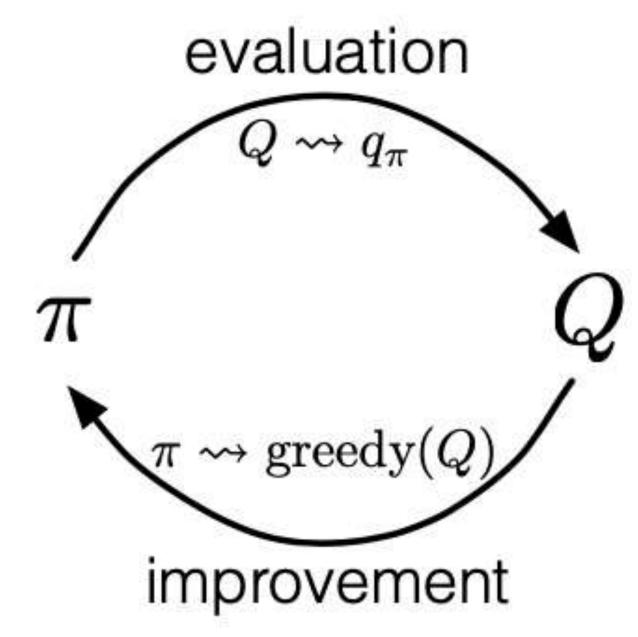
Every visit MC for state-action evaluation: average of the returns following every visit to (s,a)

Every

```
First visit MC prediction, for estimating V \approx v_{\pi}
                                                                q(s,a) \approx q_{\pi}(s,a)
Input: a policy \pi to be evaluated
Initialize:
     V(s) \in \mathbb{R}, arbitrarily, for all s \in \mathcal{S} q(s,a) \in \mathbb{R}, arbitrarily, for all s \in \mathcal{S} and a \in \mathcal{A}(s)
    Returns(s) \leftarrow an empty list, for all s \in \mathcal{S} Returns(s, a) \leftarrow an empty list, for all s \in \mathcal{S} and a \in \mathcal{A}(s)
Loop forever (for each episode):
     Generate an episode following \pi: 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 to Returns(S_t) Append G to Returns(S_t, A_t)
              -V(S_t) \leftarrow \operatorname{average}(Returns(S_t))
                                                                                 average(Returns(S_t, A_t))
                                                                  q(S_t, A_t) \leftarrow
```

Generalized Policy Iteration (GPI): Expensive

Assumptions: 1) Exploring start; 2) Infinitely many episodes



State-action value evaluation as before: $\pi_k(s) \to q_{\pi_k}(s, a)$

Greedy policy improvement: $\pi_{k+1}(s) = \arg \max_{a} q_{\pi_k}(s, a)$

$$\pi_0 \xrightarrow{\scriptscriptstyle \mathrm{E}} q_{\pi_0} \xrightarrow{\scriptscriptstyle \mathrm{I}} \pi_1 \xrightarrow{\scriptscriptstyle \mathrm{E}} \pi_1 \xrightarrow{\scriptscriptstyle \mathrm{E}} q_{\pi_1} \xrightarrow{\scriptscriptstyle \mathrm{I}} \pi_2 \xrightarrow{\scriptscriptstyle \mathrm{E}} \cdots \xrightarrow{\scriptscriptstyle \mathrm{I}} \pi_* \xrightarrow{\scriptscriptstyle \mathrm{E}} q_*$$

Generalized Policy Iteration (GPI): Cheap

Assumptions: 1) Exploring start; 2) Infinitely many episodes

```
Monte Carlo ES (Exploring Starts), for estimating \pi \approx \pi_*
Initialize:
     \pi(s) \in \mathcal{A}(s) (arbitrarily), for all s \in \mathcal{S}
     Q(s,a) \in \mathbb{R} (arbitrarily), for all s \in \mathcal{S}, a \in \mathcal{A}(s)
     Returns(s, a) \leftarrow \text{empty list, for all } s \in \mathcal{S}, \ a \in \mathcal{A}(s)
Loop forever (for each episode):
     Choose S_0 \in \mathcal{S}, A_0 \in \mathcal{A}(S_0) randomly such that all pairs have probability > 0
     Generate an episode from S_0, A_0, following \pi: S_0, A_0, R_1, \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 the pair S_t, A_t appears in S_0, A_0, S_1, A_1, \ldots, S_{t-1}, A_{t-1}:
                Append G to Returns(S_t, A_t)
               Q(S_t, A_t) \leftarrow \operatorname{average}(Returns(S_t, A_t))
               \pi(S_t) \leftarrow \operatorname{argmax}_a Q(S_t, a)
```

How to avoid exploring starts?

Two situations

- O *On-policy* method: attempt to evaluate or improve the policy that is used to make decisions
 - e.g., all methods that we have learned so far
 - Change greedy policy $\pi(s)$ to ε -greedy policy $\pi(a \mid s) > 0$ for all s and a
 - One policy has to balance exploration (not greedy) and exploitation (greedy)
- O Off-policy method: evaluate or improve a policy different from that used to generate the data

•

How to avoid exploring starts (on-policy)?

```
On-policy first-visit MC control (for \varepsilon-soft policies), estimates \pi \approx \pi_*
Algorithm parameter: small \varepsilon > 0
Initialize:
    \pi \leftarrow an arbitrary \varepsilon-soft policy
    Q(s,a) \in \mathbb{R} (arbitrarily), for all s \in \mathcal{S}, a \in \mathcal{A}(s)
     Returns(s, a) \leftarrow \text{empty list, for all } s \in \mathcal{S}, \ a \in \mathcal{A}(s)
Repeat forever (for each episode):
    Generate an episode following \pi: S_0, A_0, R_1, \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 the pair S_t, A_t appears in S_0, A_0, S_1, A_1, \ldots, S_{t-1}, A_{t-1}:
              Append G to Returns(S_t, A_t)
              Q(S_t, A_t) \leftarrow \text{average}(Returns(S_t, A_t))
              A^* \leftarrow \operatorname{arg\,max}_a Q(S_t, a)
                                                                                     (with ties broken arbitrarily)
              For all a \in \mathcal{A}(S_t):
                      \pi(a|S_t) \leftarrow \begin{cases} 1 - \varepsilon + \varepsilon/|\mathcal{A}(S_t)| & \text{if } a = A^* \\ \varepsilon/|\mathcal{A}(S_t)| & \text{if } a \neq A^* \end{cases}
```

By Richard S. Sutton and Andrew G. Barto.

How to avoid exploring starts?

Two situations

- O On-policy method: attempt to evaluate or improve the policy that is used to make decisions
 - One policy has to balance exploration (not greedy) and exploitation (greedy)
- O Off-policy method: evaluate or improve a policy different from that used to generate the data
 - Use two policies
 - Behavior policy encourages exploration and generate samples, e.g., stochastic policy $\pi_b(a \mid s) > 0$ for all s and a
 - Target policy encourages exploitation and approaches to the optimal policy, e.g., deterministic policy $\pi(s)$

•

How to avoid exploring starts (off-policy)?

Off-policy method:

- Algorithm: use behavior policy $\pi_b(a \mid s)$ (non-greedy) to generate data to estimate target policy $\pi(s)$ (greedy)
- We cannot completely believe data and need to process them!

How to avoid exploring starts (off-policy)?

Question: How can we process the data?

Answer: Important sampling

- A general technique for estimating expected values under one distribution given samples from another
- e.g., estimating the mean of a Bernoulli random variable X_p from the samples of another Bernoulli random variable X_q
- $Pr(X_p = 1) = p = 1 Pr(X_p = 0)$ and $E[X_p] = (1 p) * 0 + p * 1 = p$
- $Pr(X_q = 1) = q = 1 Pr(X_q = 0)$ and $E[X_q] = q$

How to avoid exploring starts (off-policy)?

Important sampling

- Goal: estimate $E[X_p]$ using the samples of X_q : $\{x_i\}_{i=1}^n$
- Law of large number: $E[X_q] \approx \frac{1}{n} \sum_{i=1}^n x_i$
- e.g., there should be about $nq x_i$'s as 1, and $(1 q)n x_i$'s as 0
- How can we make the average of these samples look like the average of the samples of X_p ?
- If we had the samples of X_p , there should be about np numbers as 1, and (1-p)n numbers as 0
- Hence, we modify $\{x_i\}_{i=1}^n$ to obtain $\{\bar{x}_i\}_{i=1}^n$ as $\bar{x}_i = \begin{cases} \frac{p}{q}x_i, & \text{if } x_i = 1\\ \frac{1-p}{1-q}x_i & \text{if } x_i = 0 \end{cases}$
- Finally, $\mathrm{E}[X_p] \approx \frac{1}{n} \sum_{i=1}^n \bar{x}_i = \text{the average of } nq \text{ samples of } 1 * \frac{p}{q} \text{ and } n(1-q) \text{ samples of } 0 * \frac{1-p}{1-q} \approx p = \mathrm{E}[X_p]$

How to avoid exploring starts (off-policy)?

Important sampling

- Goal: estimate $E[X_p]$ using the samples of X_q : $\{x_i\}_{i=1}^n$
- How can we make the average of these samples look like the average of the samples of X_p ?
- Hence, we modify $\{x_i\}_{i=1}^n$ to obtain $\{\bar{x}_i\}_{i=1}^n$ as $\bar{x}_i = \frac{\Pr[X_p = x_i]}{\Pr[X_q = x_i]} x_i$
- Then $E[X_p] \approx \frac{1}{n} \sum_{i=1}^n \bar{x}_i$

How to avoid exploring starts (off-policy)?

Off-policy method:

- Algorithm: use behavior policy $\pi_b(a \mid s)$ (non-greedy) to generate data to estimate target policy $\pi(s)$ (greedy)
- Use important sampling to process the data

 $\Pr[G_t]$ under π_h

•
$$v_{\pi_b}(s) = E_{\pi_b}[G_t | S_t = s]$$

•
$$v_{\pi}(s) = \mathrm{E}_{\pi}[G_{t} | S_{t} = s] = \mathrm{E}_{\pi_{b}}[\rho_{t:T-1}G_{t} | S_{t} = s]$$
 with T as the termination time and
$$\rho_{t:T-1} = \Pi_{k=t}^{T-1} \frac{\pi(A_{k} | S_{k})}{\pi_{b}(A_{k} | S_{k})} = \frac{\Pr[A_{t}, S_{t+1}, A_{t+1}, ..., S_{T}] \text{ under } \pi}{\Pr[A_{t}, S_{t+1}, A_{t+1}, ..., S_{T}] \text{ under } \pi_{b}}$$

$$- \Pr[G_{t}] \text{ under } \pi$$

•

How to avoid exploring starts (off-policy)?

Off-policy method:

- Most steps are the same as the on-policy methods discussed previously
- Use behavior policy $\pi_b(a \mid s)$ (non-greedy) to generate data to estimate target policy $\pi(s)$ (greedy)
- Use the formula $v_{\pi}(s) = \mathrm{E}_{\pi_b}[\rho_{t:T-1}G_t \,|\, S_t = s]$ with $\rho_{t:T-1} = \Pi_{k=t}^{T-1} \frac{\pi(A_k \,|\, S_k)}{\pi_b(A_k \,|\, S_k)}$ to evaluate and improve the policy $\pi(s)$
- Similarly, $q_{\pi}(s, a) := \mathbb{E}_{\pi_b}[\rho_{t+1:T-1}G_t | S_t = s, A_t = a]$

How to avoid exploring starts (off-policy)?

```
Off-policy MC control, for estimating \pi \approx \pi_*
Initialize, for all s \in \mathcal{S}, a \in \mathcal{A}(s):
     Q(s, a) \in \mathbb{R} (arbitrarily)
     C(s,a) \leftarrow 0
     \pi(s) \leftarrow \operatorname{argmax}_a Q(s, a) (with ties broken consistently)
Loop forever (for each episode):
     b \leftarrow \text{any soft policy}
     Generate an episode using b: S_0, A_0, R_1, \ldots, S_{T-1}, A_{T-1}, R_T
     G \leftarrow 0
     W \leftarrow 1
     Loop for each step of episode, t = T-1, T-2, \ldots, 0:
          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)]
          \pi(S_t) \leftarrow \operatorname{argmax}_a Q(S_t, a) (with ties broken consistently)
          If A_t \neq \pi(S_t) then exit inner Loop (proceed to next episode)
          W \leftarrow W \frac{1}{b(A_t|S_t)}
```

Summary

- Model-free method
- Introduced policy iteration method to compute the optimal policy
- Introduced value iteration method to compute the optimal policy
- Generalized policy iteration (GPI)
- Introduced on-policy and off-policy for exploring starts (without discounted rewards)
- Exploring starts with discounted rewards not discussed

Temporal-Difference Learning

Temporal-Difference (TD) Learning Overview

- One central and novel idea of RL
- Combining DP and MC
- Learn from raw data without model: model-free
- Update estimates using learned models: model-based
- Policy evaluation and improvement based on generalized policy iteration (GPI)

Main idea

$$v_{\pi}(s) := E_{\pi}[G_{t} | S_{t} = s]$$

$$= E_{\pi}[R_{t+1} + \gamma G_{t+1} | S_{t} = s]$$

$$= E_{\pi}[R_{t+1} + \gamma v_{\pi}(S_{t+1}) | S_{t} = s]$$

• MC update:

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

 G_t is available only when an episode ends

• TD update:

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

- MC: R_{t+1} is available right at the next time step, hence, online, fully incremental, and no need for discounts
- DP: $V(S_{t+1})$ is the currently learned model

Tabular TD(0) for estimating v_{π}

```
Input: the policy \pi to be evaluated
Algorithm parameter: step size \alpha \in (0, 1]
Initialize V(s), for all s \in S^+, arbitrarily except that V(terminal) = 0
Loop for each episode:
   Initialize S
   Loop for each step of episode:
       A \leftarrow \text{action given by } \pi \text{ for } S
       Take action A, observe R, S'
       V(S) \leftarrow V(S) + \alpha [R + \gamma V(S') - V(S)]
       S \leftarrow S'
   until S is terminal
```

Main idea

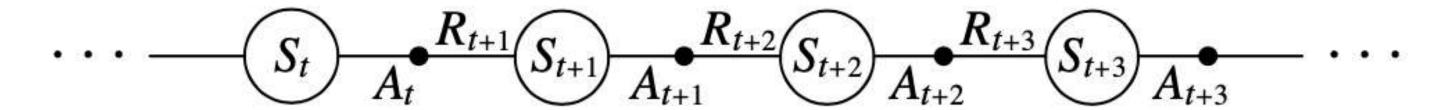
O MC methods:

- on-policy updates using one ε -greedy policy
- off-policy updates using important sampling with two policies

O TD update:

- on-policy updates for policies: Sarsa with one policy
- off-policy updates for policies: Q-learning with two policies

Sarsa: on-policy updates



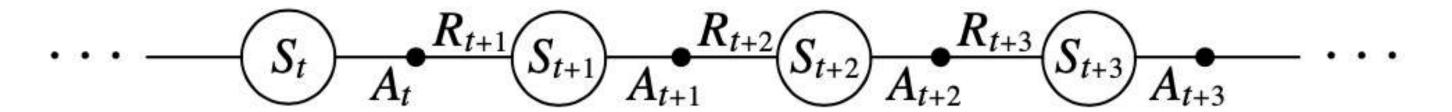
• Value function update:

$$Q(S_t, A_t) \leftarrow Q(S_t, A_t) + \alpha [R_{t+1} + \gamma Q(S_{t+1}, A_{t+1}) - Q(S_t, A_t)]$$

- Use ε -greedy policy with $\varepsilon = \frac{1}{t}$
- Converges with probability 1 to q_* if all (s, a)-pairs are visited an infinite number of times

```
Sarsa (on-policy TD control) for estimating Q \approx q_*
Algorithm parameters: step size \alpha \in (0,1], small \varepsilon > 0
Initialize Q(s,a), for all s \in S^+, a \in A(s), arbitrarily except that Q(terminal, \cdot) = 0
Loop for each episode:
   Initialize S
   Choose A from S using policy derived from Q (e.g., \varepsilon-greedy)
   Loop for each step of episode:
                                                                                       results of the same policy
       Take action A, observe R, S'
       Choose A' from S' using policy derived from Q (e.g., \varepsilon-greedy)
      Q(S,A) \leftarrow Q(S,A) + \alpha \left[ R + \gamma Q(S',A') - Q(S,A) \right]
       S \leftarrow S'; A \leftarrow A';
   until S is terminal
```

Q-learning: off-policy updates



• Value function update:

$$Q(S_t, A_t) \leftarrow Q(S_t, A_t) + \alpha [R_{t+1} + \gamma \max_{a} Q(S_{t+1}, a) - Q(S_t, A_t)]$$

- Use one ε -greedy policy (prefer exploration) to generate data
- Use another greedy policy (prefer exploitation) to update value function
- Converges with probability 1 to q_* since all (s, a)-pairs are visited an infinite number of times

 $Q \approx q_*$

Q-learning (off-policy TD control) for estimating $\pi \approx \pi_*$

Algorithm parameters: step size $\alpha \in (0, 1]$, small $\varepsilon > 0$

Initialize Q(s,a), for all $s \in S^+$, $a \in A(s)$, arbitrarily except that $Q(terminal, \cdot) = 0$

Loop for each episode:

Initialize S

Loop for each step of episode:

Choose A from S using policy derived from Q (e.g., ε -greedy)

Take action A, observe R, S'

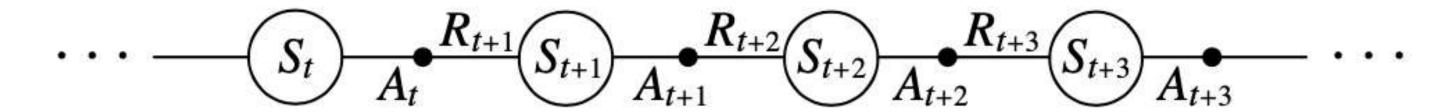
$$Q(S,A) \leftarrow Q(S,A) + \alpha \left[R + \gamma \max_{a} Q(S',a) - Q(S,A) \right]$$

$$S \leftarrow S'$$

until S is terminal

results of different policies

Expected Sarsa: on-policy updates



• Value function update:

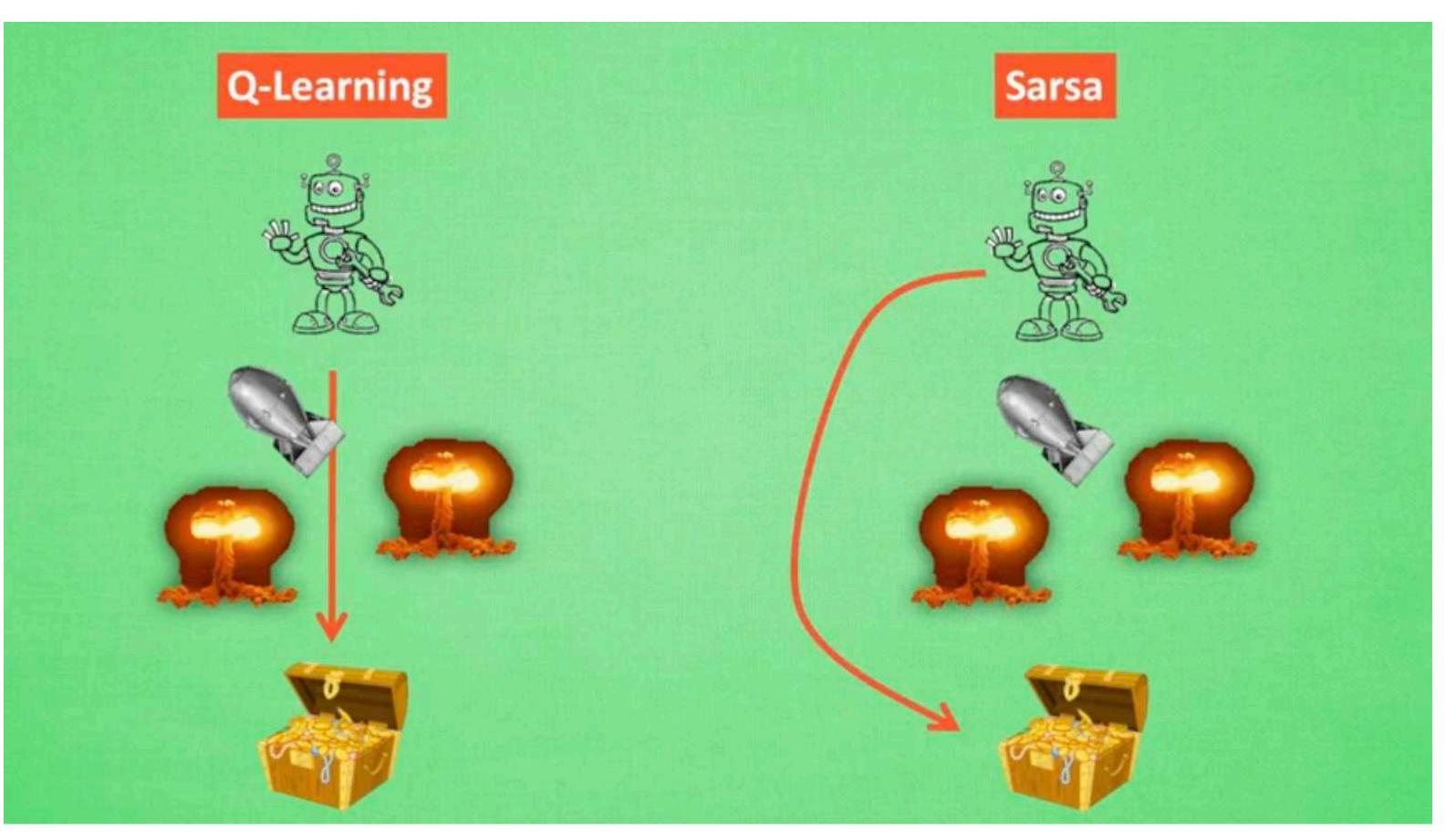
$$Q(S_t, A_t) \leftarrow Q(S_t, A_t) + \alpha [R_{t+1} + \gamma \mathbb{E}_{\pi}[Q(S_{t+1}, A_{t+1}) | S_{t+1}] - Q(S_t, A_t)]$$

$$= Q(S_t, A_t) + \alpha [R_{t+1} + \gamma \sum_{a} \pi(a | S_{t+1}) Q(S_{t+1}, a) - Q(S_t, A_t)]$$

- Sarsa: use ε -greedy policy π to choose A_{t+1} randomly given S_{t+1}
- Expected Sarsa: use the expectation to eliminate the variance of the randomness

Comparison

- Sarsa more conservative
- Q-learning more aggressive



By Mofanpy

Temporal-Difference (TD) Learning Comparison

- Expected Sarsa v.s. Sarsa more stable, working for a wide range of step size α , but more expensive
- Expected Sarsa v.s. Q-learning more conservative

Maximization Bias: bias towards positive value function estimation

- o $\max_{a \in \mathcal{A}(s)} Q(s, a)$ is used in TD learning
- O Example:
 - A single state s with many actions
 - $q_*(s, a) = 0$ for most a's
 - $Q(s, a) \neq 0$ for most a
 - $\max_{a \in \mathcal{A}(s)} Q(s, a)$ has a bias towards actions such that Q(s, a) > 0

How to avoid the maximum bias?

- Idea: Use two policies and alternatively update
- Use $A^* = \arg \max_a Q_1(a)$ to take an action
- Use $Q_2(A^*) = Q_2(\arg\max_a Q_1(a))$ to provide model estimate to update Q_1
- Example:

$$Q_1(S_t, A_t) \leftarrow Q_1(S_t, A_t) + \alpha [R_{t+1} + \gamma Q_2(S_{t+1}, \arg \max_{a} Q_1(S_{t+1}, a)) - Q_1(S_t, A_t)]$$

• Similarly, in the next round, use Q_2 to take action to update Q_1

How to avoid the maximum bias?

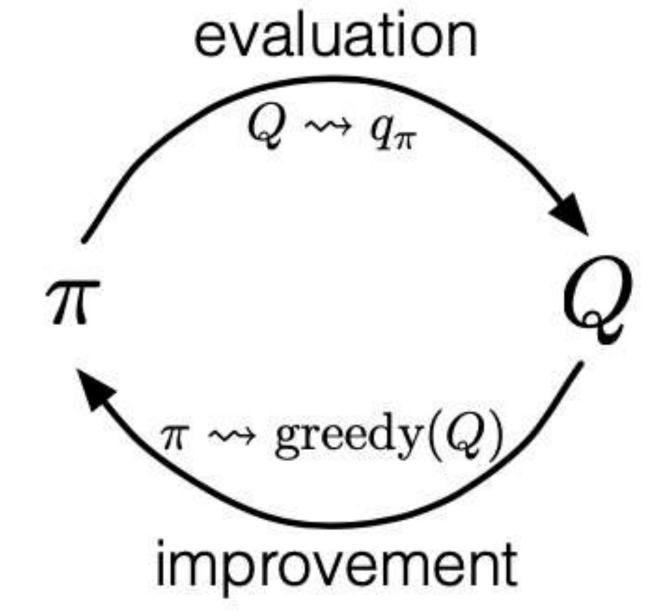
- Idea: Use two policies and alternatively update
- Use $A^* = \arg \max_{a} Q_1(a)$ to take an action
- Use $Q_2(A^*) = Q_2(\arg\max_a Q_1(a))$ to provide model estimate to update Q_1
- arg max $Q_1(a)$ has a bias towards actions with positive values in Q_1
- arg max $Q_1(a)$ has no bias towards actions with positive values in Q_2
- If Q_1 and Q_2 have no bias, then the new Q_2 has no bias
- Similarly, if Q_1 and Q_2 have no bias, then the new Q_1 has no bias

Double Q-learning to avoid the maximum bias

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Double Q-learning, for estimating Q_1 \approx Q_2 \approx q_*
Algorithm parameters: step size \alpha \in (0, 1], small \varepsilon > 0
Initialize Q_1(s,a) and Q_2(s,a), for all s \in S^+, a \in A(s), such that Q(terminal, \cdot) = 0
Loop for each episode:
   Initialize S
   Loop for each step of episode:
       Choose A from S using the policy \varepsilon-greedy in Q_1 + Q_2
       Take action A, observe R, S'
       With 0.5 probability:
          Q_1(S,A) \leftarrow Q_1(S,A) + \alpha \Big(R + \gamma Q_2(S', \operatorname{argmax}_a Q_1(S',a)) - Q_1(S,A)\Big)
       else:
          Q_2(S, A) \leftarrow Q_2(S, A) + \alpha \Big(R + \gamma Q_1(S', \operatorname{argmax}_a Q_2(S', a)) - Q_2(S, A)\Big)
       S \leftarrow S'
   until S is terminal
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Generalized Policy Iteration (GPI): Expensive

On-policy (expected) Sarsa with ε -greedy policies to avoid exploring starts; Off-policy (double) Q-Leaning with ε -greedy policies to avoid exploring starts;



State-action value evaluation as before: $\pi_k(s) \to q_{\pi_k}(s, a)$

Greedy policy improvement: $\pi_{k+1}(s) = \arg \max_{a} q_{\pi_k}(s, a)$

$$\pi_0 \xrightarrow{\mathrm{E}} q_{\pi_0} \xrightarrow{\mathrm{I}} \pi_1 \xrightarrow{\mathrm{E}} q_{\pi_1} \xrightarrow{\mathrm{I}} \pi_2 \xrightarrow{\mathrm{E}} \cdots \xrightarrow{\mathrm{I}} \pi_* \xrightarrow{\mathrm{E}} q_*$$

Summary

- Combining model-free and model-based
- Generalized policy iteration (GPI) as in MC and DP
- Introduced value iteration method to compute the optimal policy: Sarsa and Q-learning
- Introduced off-policy double Q-learning to avoid the maximum bias
- Can use on-policy and off-policy for exploring starts as in MC