

Lecture 15 Reinforcement Learning

Outline

- Reinforcement Learning Problem
- Principle of optimality
- Markov Decision Process
- Monte-Carlo Methods
- Temporal Difference Learning

Any definition

RL brings a way of programming agents by reward and punishment without specifying *how* the task is to be achieved. (Kaelbling,1996)

 \rightarrow Based on trial-error interactions

 \rightarrow A set of problems rather than a set of techniques

Control Learning

Learning to choose actions

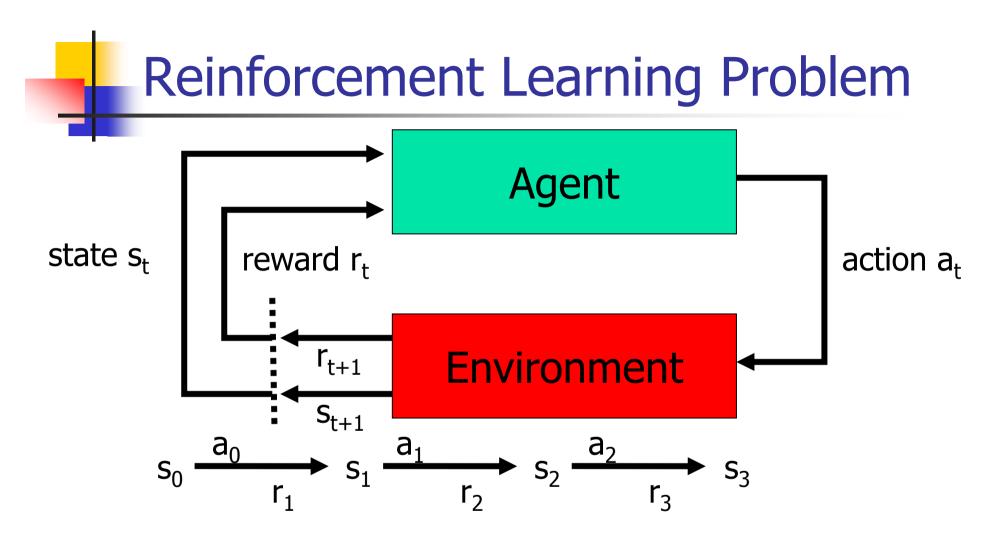
- Robot learning to dock to a battery station
- Learning to choose actions to optimize a factory output
- Learning to play Backgammon

Problem characteristics:

- Delayed reward
- No direct feedback (error signal) for good and bad actions
- Opportunity for active exploration
- Possibly that state is only partially observable

Learning to play Backgammon

- Immediate reward
 - +100 win
 - -100 loose
 - 0 for all other actions/states
- Trained by playing 1.5 million games against itself (Tesauro [1995])
- Now approximately equal to the best human player



Goal: Learn to choose actions a_t that maximize future rewards $r_1+\gamma r_2+\gamma^2 r_3+...$, where $0<\gamma<1$ is a discount factor

Models of Optimal Behavior:

Agent tries to maximize one of the following:

finite-horizon Model:

$$E(\sum_{t=0} r_t)$$
;

infinite-horizon discounted model:

average-reward model:

$$E(\sum_{t=0}^{\infty} \gamma^t r_t) \quad .$$

$$\lim_{h \to \infty} E(\frac{1}{h} \sum_{t=0}^{h} r_t) \quad .$$

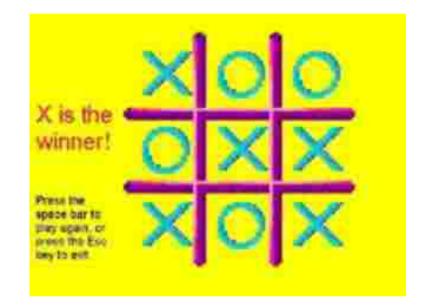
Example 1: Slot Machine

- State: configuration of slots
- Action: stopping time
- Reward: \$\$\$
- Problem: Find $\pi: S \rightarrow A$ that
- maximizes R



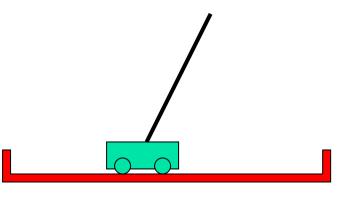
Example 2: Tic Tac Toe

- State: board
- Action: next move
- Reward: 1 for win, -1 for loss, 0 for draw
- Problem: Find π :S \rightarrow A that
- maximizes R



Example 3: Inverted Pendulum

- State: x(t), x'(t), q(t), q'(t)
- Action: Force F (bang bang)
- Reward: 1 for any step where pole balanced
- Problem: Find π :S \rightarrow A that
- maximizes R



Example 4: Mobile Robot

- State: location of robot, people
- Action: motion
- Reward: number of happy faces
- Problem: Find π :S \rightarrow A that
- maximizes R

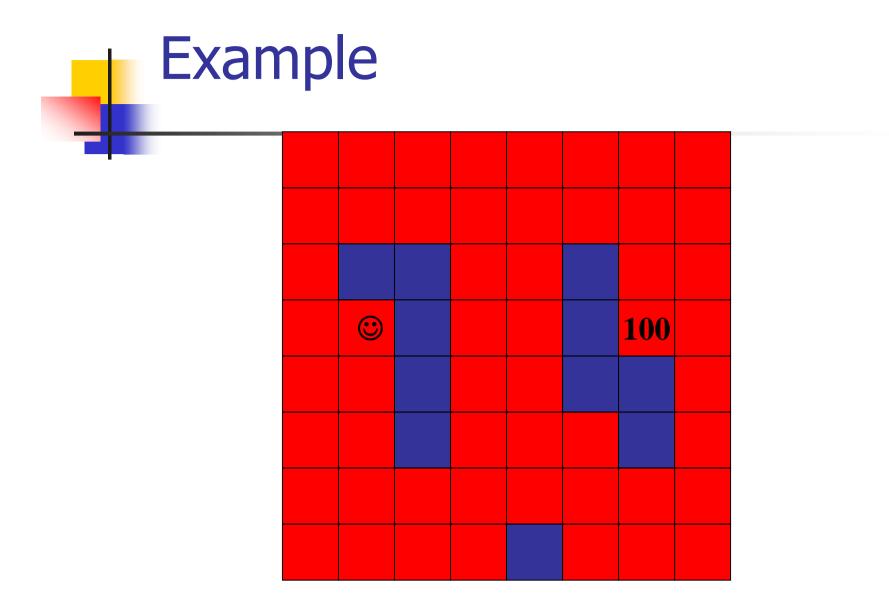


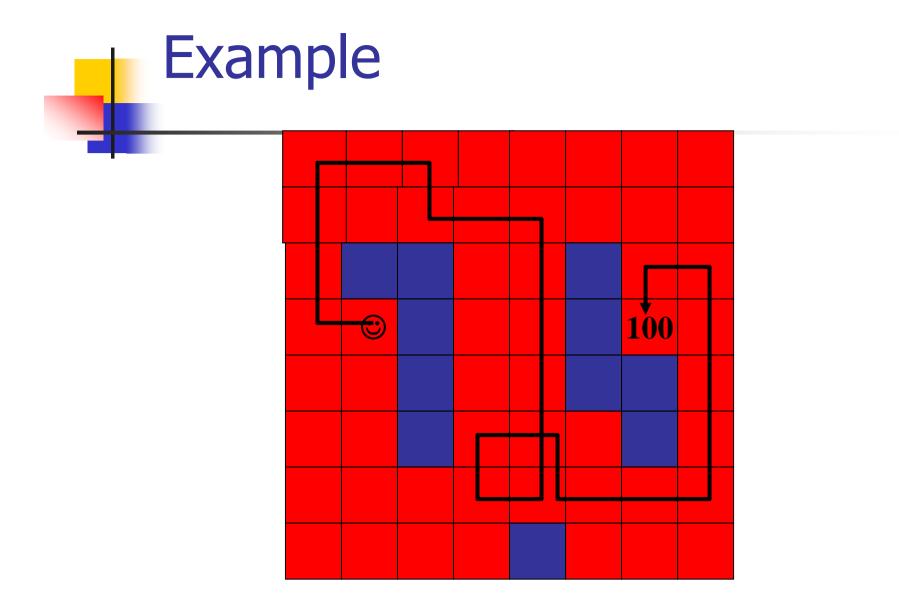
Common Features of RL Problems

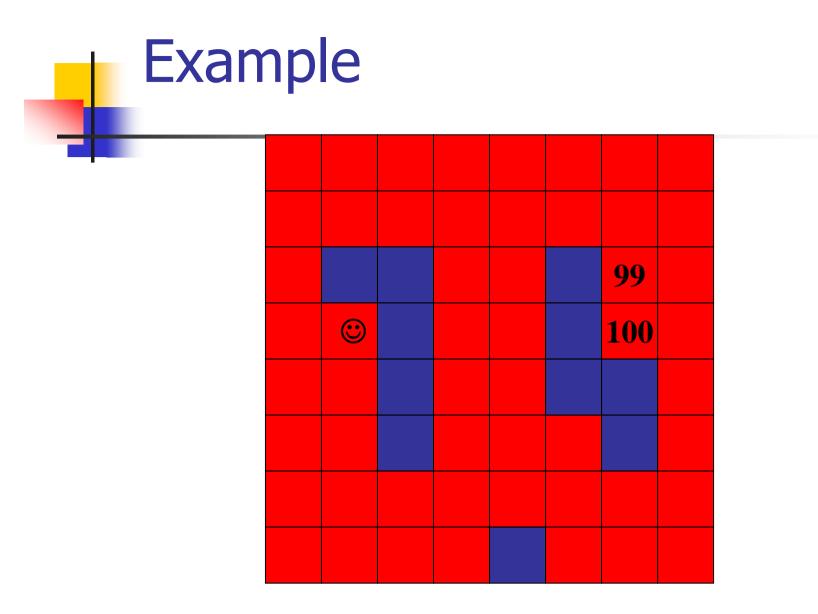
- Temporal Credit Assignment
- Exploration/Exploitation
- Planning and Acting under Uncertainty

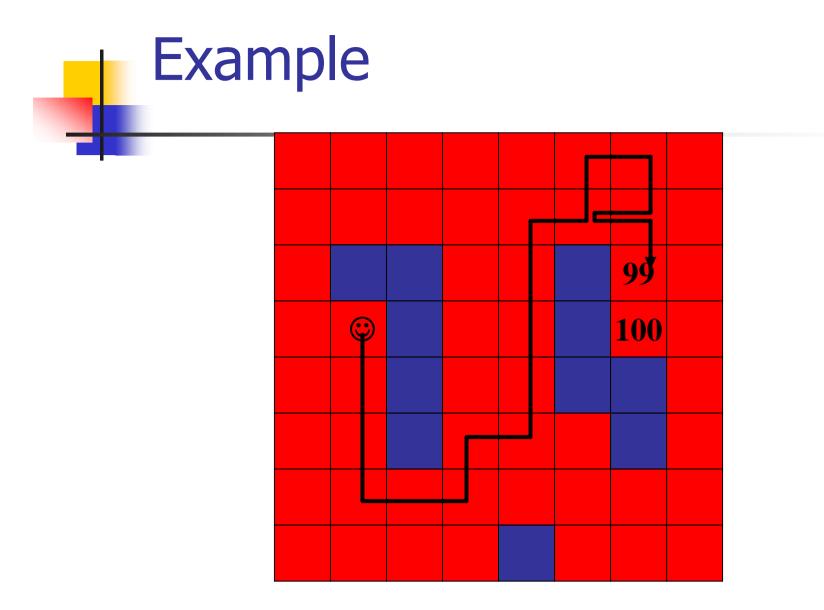
The Principle of Optimality

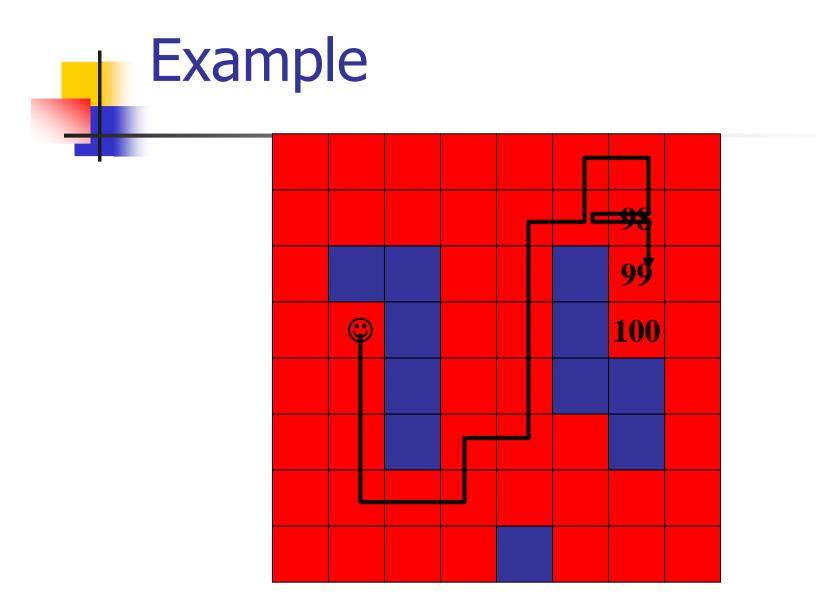
- "an optimal set of decisions has the property that whatever the first decision is, the remaining decisions must be optimal w/ respect to the outcome...of the first decision." –R. Bellman
- "if you don't do the best you have with what you happen to have got, you will never do the best you might have done with what you should have had." –R. Aris

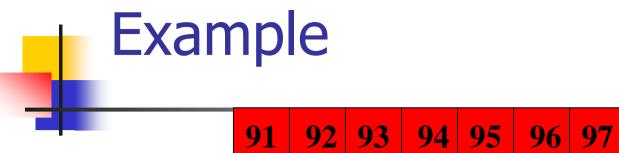




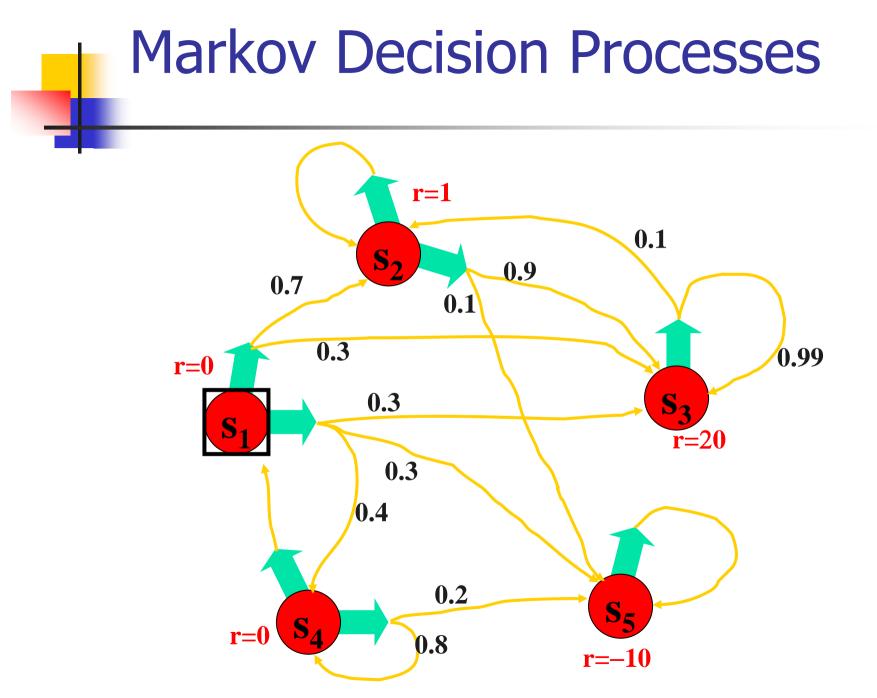








| 91 | 92 | 93 | 94 | 95 | 96 | 97 | 96 |
|----|-----------|----|----|-----------|----|-----------|-----------|
| 92 | 93 | 94 | 95 | 96 | 97 | 98 | 97 |
| 91 | | | 94 | 95 | | 99 | 98 |
| 90 | : | | 93 | 94 | | 100 | 99 |
| 89 | 88 | | 92 | 93 | | | 98 |
| 88 | 89 | | 91 | 92 | 93 | | 97 |
| 89 | 90 | 91 | 92 | 93 | 94 | 95 | 96 |
| 88 | 89 | 90 | 91 | | 93 | 94 | 95 |



Markov Decision Process (MDP)

- Finite set of states S
- Finite set of actions A
- At each time step the agent observes state s_t∈S and chooses action a_t∈A(s_t)
- Then receives immediate reward r_{t+1}
- And state changes to s_{t+1}
- Markov assumption : $s_{t+1} = \delta(s_t, a_t)$ and $r_{t+1} = r(s_t, a_t)$
 - Next reward and state only depend on *current* state s_t and action a_t
 - Functions $\delta(s_t, a_t)$ and $r(s_t, a_t)$ may be non-deterministic
 - Functions δ(s_t,a_t) and r(s_t,a_t) not necessarily known to agent

What's Different from HMMs?

- Controllable Markov chain (HMMs are not controllable)
- Equally, an FSM with probabilistic transitions and full state output
- Objective: Learn a policy (HMMs learn models)
- State fully observable (state in HMMs is only partially observable)

Why MDPs?

- Natural model of physical, apparently random phenomenom
- MDP's obey the principle of optimality (under suitable restrictions on the reward function)
 - The Bellman equation: (fixed policy)

 $Q(s,a) = r(s) + E_{s,a}[Q(s',a')]$

- The Bellman equation: $Q^*(s) = r(s) + \max_{a^*} E_s[Q^*(s', a^*)]$
- There is a (almost unique) optimal *deterministic* mapping from states to actions



Execute actions in the environment, observe results and

- Learn a policy $\pi_t(s,a) : S \to A$ from states $s_t \in S$ to actions $a_t \in A$ that maximizes the expected reward : $E[r_t + \gamma r_{t+1} + \gamma^2 r_{t+2} + ...]$ from any starting state s_t
- $0 < \gamma < 1$ is the discount factor for future rewards
- Target function is $\pi_t(s,a) : S \to A$
- But there are no direct training examples of the form <s,a>
- Training examples are of the form <<s,a>,r>

State Value Function

Consider deterministic environments, namely $\delta(s,a)$ and r(s,a) are deterministic functions of s and a.

For each policy $\pi(s,a) : S \rightarrow A$ the agent might adopt we define an evaluation function:

 $V^{\pi}(s) = r_t + \gamma r_{t+1} + \gamma^2 r_{t+2} + \ldots = \sum_{i=0} r_{t+i} \gamma^i$

where $r_t,\,r_{t+1},...$ are generated by following the policy π from start state s

Task: Learn the optimal policy π^* that maximizes V^{π}(s)

$$\pi^* = \operatorname{argmax}_{\pi} V^{\pi}(s) , \forall s$$

Action Value Function

 State value function denotes the reward for starting in state s and following policy π.

 $V^{\pi}(s) = r_{t} + \gamma r_{t+1} + \gamma^{2} r_{t+2} + ... = \sum_{i=0} r_{t+i} \gamma^{i}$

 Action value function denotes the reward for starting in state s, taking action a and following policy π afterwards.

 $Q^{\pi}(s,a) = r(s,a) + \gamma r_{t+1} + \gamma^2 r_{t+2} + ... = r(s,a) + \gamma V^{\pi}(\delta(s,a))$

Predicting the Value of a State

Key idea: Learn a Value Function

 $Q:SxA \rightarrow \Re$

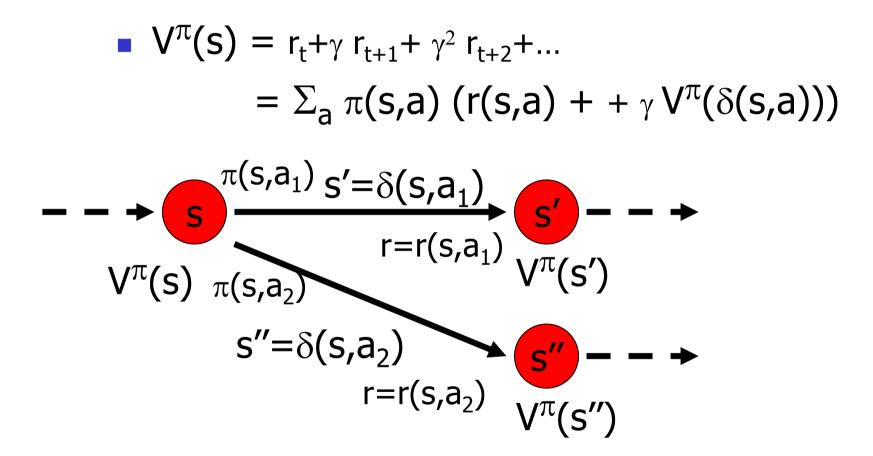
such that

$$Q(s,a) = \sum \gamma^t r_t$$

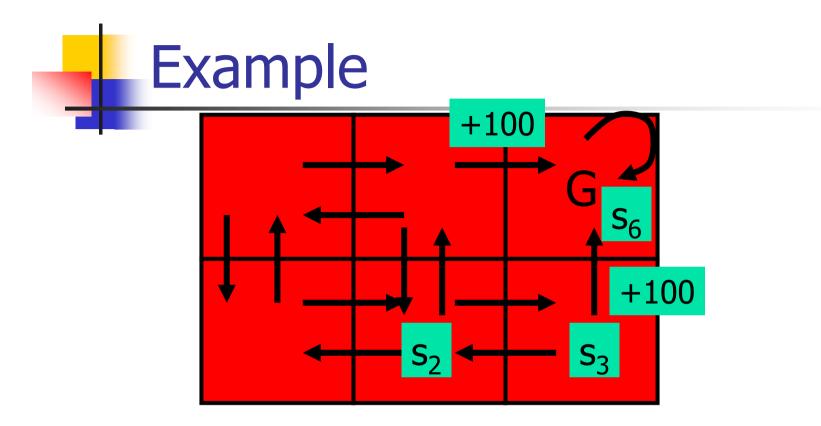
A mapping that predicts what each state *s* is worth, given we take action *a* (Sum of future rewards.)

- Why: Duality between value functions and policies
 - By knowing the optimal policy, we can calculate the value of each state
 - By knowing the optimal value-function, we can calculate the optimal move

Bellman Equation (Deterministic Case)



Set of |s| linear equations, solve it directly or by policy evaluation.



G: terminal state, upon entering G agent obtains a reward of +100, remains in G forever and obtains no further rewards

Compute V^{π}(s) for equil-probable policy π (s,a)=1/|a| V^{π}(s₃) = $\frac{1}{2} \gamma V^{\pi}(s_2) + \frac{1}{2} (100 + \gamma V^{\pi}(s_6))$

Iterative Policy Evaluation

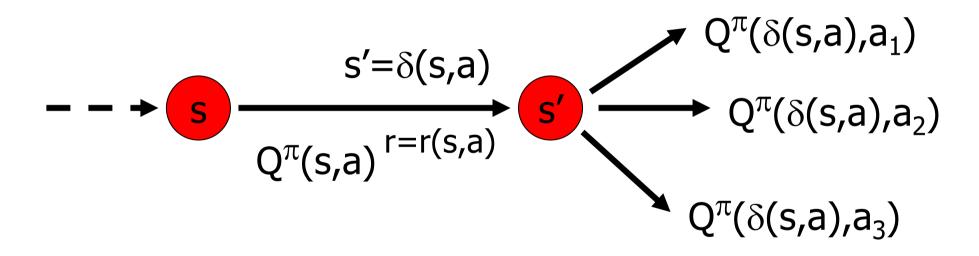
Instead of solving the Bellman equation directly one can use *iterative policy evaluation* by using the Bellman equation as an update rule.

$$V_{k+1}^{\pi}(s) = \Sigma_a \pi(s,a) (r(s,a) + \gamma V_k^{\pi}(\delta(s,a)))$$

The sequence $V_k^{\ \pi}$ is guaranteed to converge to V^{π}

Bellman Equation (Deterministic Case)

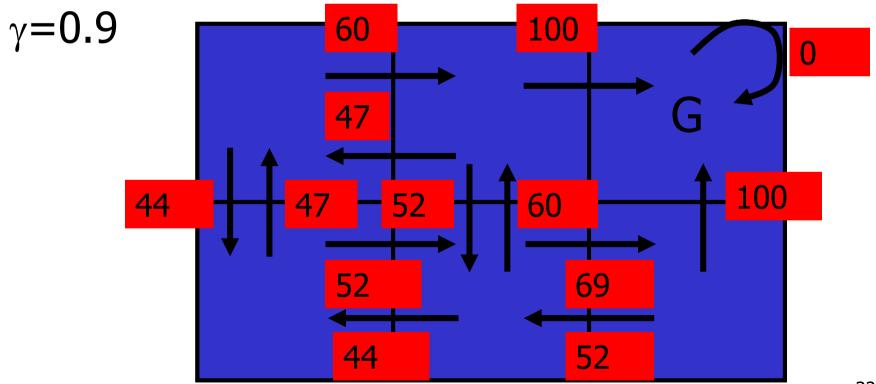
• $Q^{\pi}(s,a) = r(s,a) + \gamma \Sigma_{a'} \pi(\delta(s,a),a') Q^{\pi}(\delta(s,a),a')$



Iterative Policy Evaluation

Bellman equation as an update rule for action-value function:

 $Q_{k+1}^{\pi}(s,a) = r(s,a) + \gamma \Sigma_{a'} \pi(\delta(s,a),a') Q_k^{\pi}(\delta(s,a),a')$



Optimal Value Functions

Bellman optimality equations • $V^*(s) = \max_a Q^*(s,a)$ = $\max_a (r(s,a) + \gamma V^*(\delta(s,a)))$ • $Q^*(s,a) = r(s,a) + V^*(\delta(s,a)))$ = $r(s,a) + \gamma \max_{a'} Q^*(\delta(s,a),a')$

Policy Improvement

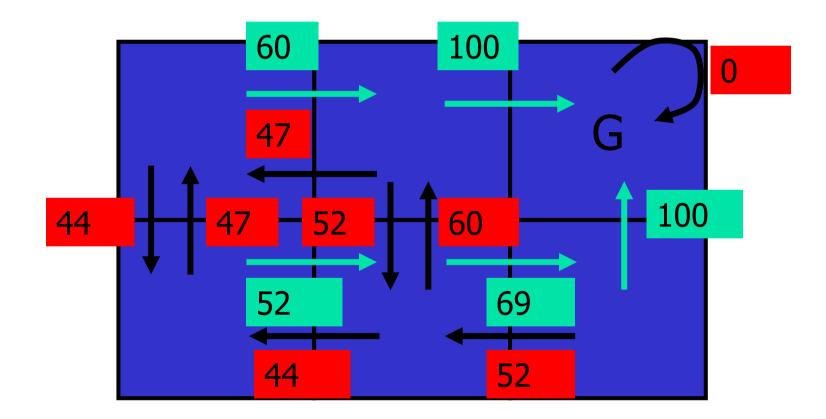
- Suppose we have determined the value function V^π for an arbitrary deterministic policy π.
- For some state s we would like to know if it is better to choose an action a≠π(s).
- Select a and follow the existing policy π afterwards gives us reward Q^π(s,a)
- If $Q^{\pi}(s,a) > V^{\pi}$ then a is obviously better than $\pi(s)$
- Therefore choose new policy π' as $\pi'(s) = \operatorname{argmax}_a Q^{\pi}(s,a) = \operatorname{argmax}_a r(s,a) + \gamma V^{\pi}(\delta(s,a))$

$$\pi(s,a)=1/|a| \begin{cases} \nabla^{\pi}=63 & \nabla^{\pi}=71 & \nabla^{\pi}=0 \\ \nabla^{\pi}=56 & \nabla^{\pi}=61 & \nabla^{\pi}=78 \end{cases} r=100$$

 $\pi'(s) = \operatorname{argmax}_a r(s,a) + \gamma V^{\pi}(\delta(s,a))$



 $\pi'(s) = \operatorname{argmax}_a Q^{\pi}(s,a)$



Generalized Policy Iteration

Intertwine *policy evaluation* with *policy improvement*

$$\pi_{0} \xrightarrow{E} V^{\pi_{0}} \xrightarrow{I} \pi_{1} \xrightarrow{E} V^{\pi_{1}} \xrightarrow{I} \pi_{2} \xrightarrow{E} V^{\pi_{2}} \xrightarrow{I} \cdots \xrightarrow{I} \pi^{*} \xrightarrow{E} V^{\pi^{*}}$$
evaluation
$$\overbrace{V \to V^{\pi}}^{\pi} \xrightarrow{\pi} \operatorname{greedy}(V) \xrightarrow{V}$$
improvement

Value Iteration (Q-Learning)

 Idea: do not wait for policy evaluation to converge, but improve policy after each iteration.

$$V_{k+1}^{\pi}(s) = \max_{a} (r(s,a) + \gamma V_{k}^{\pi}(\delta(s,a)))$$

or

$$Q_{k+1}^{\pi}(s,a) = r(s,a) + \gamma \max_{a'} Q_k^{\pi}(\delta(s,a),a')$$

Stop when $\forall_{s} | V_{k+1}^{\pi}(s) - V_{k}^{\pi}(s) | < \varepsilon$ or $\forall_{s,a} | Q_{k+1}^{\pi}(s,a) - Q_{k}^{\pi}(s,a) | < \varepsilon$

Non-Deterministic Case

 State transition function δ(s,a) no longer deterministic but probabilistic given by P(s'|s,a) = Pr{s_{t+1}=s'|s_t=s, a_t=a}

Transition probability that given a current state s and action a the next state is s'.

 Reward function r(s,a) no longer deterministic but probabilistic given by

$$R(s',s,a) = E\{r_{t+1}|s_t=s, a_t=a, s_{t+1}=s'\}$$

P(s'|s,a) and R(s',s,a) completely specify MDP.

Bellman Equation (Non-Deterministic Case)

- $Q^{\pi}(s,a) = \Sigma_{s'} P(s'|s,a) [R(s',s,a) + \gamma V^{\pi}(s')]$
- $V^{\pi}(s) = \Sigma_a \pi(s,a) \Sigma_{s'} P(s'|s,a) [R(s',s,a) + \gamma V^{\pi}(s')]$
- $Q^{\pi}(s,a) = \Sigma_{s'} P(s'|s,a) [R(s',s,a) + \gamma \Sigma_{a'} \pi(s',a') Q^{\pi}(s',a')]$

Bellman optimality equations:

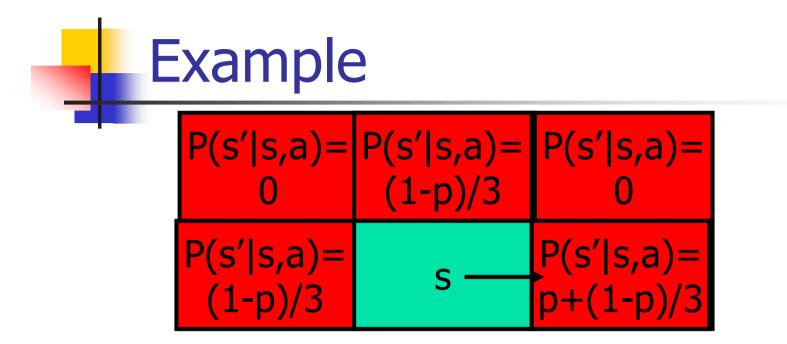
- $V^*(s) = \max_a \Sigma_{s'} P(s'|s,a) [R(s',s,a) + \gamma V^*(s')]$
- $Q^*(s,a) = \Sigma_{s'} P(s'|s,a) [R(s',s,a) + \gamma \max_{a'} Q^*(s',a')]$

Value Iteration (Q-Learning)

 $V_{k+1}^{\pi}(s) = \max_{a} \Sigma_{s'} P(s'|s,a) [R(s',s,a) + \gamma V_{k}^{\pi}(s')]$ or

 $Q_{k+1}^{\pi}(s,a) = \sum_{s'} P(s'|s,a) [R(s',s,a) + \gamma \max_{a'} Q_k^{\pi}(s',a')]$

Stop when $\forall_{s} |V_{k+1}^{\pi}(s) - V_{k}^{\pi}(s)| < \varepsilon$ or $\forall_{s,a} |Q_{k+1}^{\pi}(s,a) - Q_{k}^{\pi}(s,a)| < \varepsilon$



Now assume that actions a are nondeterministic, with probability p agent moves to the correct square indicated by a, with probability (1-p) agent moves to a random neighboring square.



Deterministic optimal value function

$$V^{*} = V^{*} = V^{*} = V^{*} = 0$$

$$100 \quad 0$$

$$V^{*} = V^{*} = V^{*} = V^{*} = 1$$

$$100 \quad 100$$

Non-deterministic optimal value function p=0.5

$$V^{*} = V^{*} = V^{*} = V^{*} = 0$$

$$77 \quad 90 \quad 0$$

$$V^{*} = V^{*} = V^{*} = V^{*} =$$

$$71 \quad 80 \quad 93$$

Reinforcement Learning

- What if the transition probabilities P(s'|s,a) and reward function R(s',s,a) are unknown?
- Can we still learn V(s), Q(s,a) and identify and optimal policy π(s,a)?
- The answer is yes. Consider the observed rewards r_t and state transitions s_{t+1}as training samples drawn from the true underlying probability functions R(s',s,a) and P(s'|s,a).
- Use approximate state V(s) and action value Q(s,a) functions

Model Free Methods

- Models of the environment:
- T: $S \times A \rightarrow \prod (S)$ and $R : S \times A \rightarrow R$
- Do we know them? Do we have to know them?
- Monte Carlo Methods
- Adaptive Heuristic Critic
- Q Learning

Monte Carlo Methods

- *Idea*:
- Hold statistics about rewards for each state
- Take the average
- This is the V(s)

- Based only on experience ☺
- Assumes episodic tasks
 B
- (Experience is divided into episodes and all episodes will terminate regardless of the actions selected.)
- Incremental in episode-by-episode sense not step-by-step sense.

Problem: Unvisited <s, a> pairs (problem of *maintaining exploration*)

For every <s, a> make sure that: P(<s, a> selected as a start state and action) >0

(Assumption of *exploring starts*)

Monte Carlo Method

- Initialize:
 - $\pi \leftarrow$ policy to be evaluated
 - $V(s) \leftarrow$ an arbitrary state-value function
 - Return(s) \leftarrow an empty list, for all s \in S
- Repeat forever
 - Generate an episode using π
 - For each state s appearing in the episode:
 - $R \leftarrow$ return following the first occurence of s
 - Append R to Returns(s)
 - $V(s) \leftarrow average(Returns(s))$

Monte Carlo Method

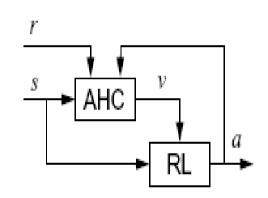
$$V^{\pi}=30$$
 $V^{\pi}=60$
 $V^{\pi}=0$
 $V^{\pi}=30$
 $V^{\pi}=40$
 $V^{\pi}=70$

 $V(s_t) \leftarrow 30 + 0.1 [0 + 0.9*0 + 0.9^2*100 - 30] = 35.1$

ADAPTIVE HEURISTIC CRITIC & TD(λ)

1

How the AHC learns, TD(0) algorithm:



$$V(s) := V(s) + \alpha(r + \gamma V(s') - V(s))$$

Architecture for the adaptive heuristic critic.

AHC : TD Algorithm



Q values in Value Iteration:

$$Q^*(s,a) = R(s,a) + \gamma \sum_{s' \in \mathcal{S}} T(s,a,s') \max_{a'} Q^*(s',a')$$

But we don't know R(s,a) and T(s,a,s')

Instead use the following :

$$Q(s, a) := Q(s, a) + \alpha(r + \gamma \max_{a'} Q(s', a') - Q(s, a))$$

Decayed α properly? Q values will converge. (Singh 1994)



- Simpler than AHC learning
- Q-Learning is exploration sensitive
- Analog to value iteration in MDP
- Most popular Model free learning algorithm

Temporal Difference Learning

- Monte-Carlo:
 - $V(s_t) \leftarrow V(s_t) + \alpha [R_t V(s_t)]$
 - target for V(s) : E {R_t | s_t=s}

Must wait until the end of the episode to update V

Temporal Difference (TD):

- $V(s_t) \leftarrow V(s_t) + \alpha [r_{t+1} + \gamma V(s_{t+1}) V(s_t)]$
- target for V(s) : E { $r_{t+1} + \gamma V(s_{t+1}) | s_t = s$ }
- TD method is *bootstrapping* by using the existing estimate of the next state $V(s_{t+1})$ for updating $V(s_t)$

TD(0) : Policy Evaluation

Initialize:

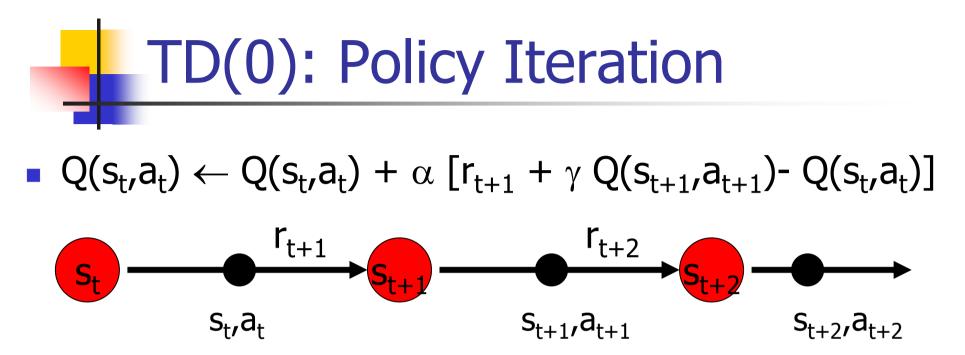
- $\pi \leftarrow$ policy to be evaluated
- V(s) ← an arbitrary state-value function

Repeat for each episode

- Initialize s
- Repeat for each step of episode
 - a \leftarrow action given by π for s
 - Take action a, observe reward r, and next state s'
 - $V(s) \leftarrow V(s_t) + \alpha [r + \gamma V(s') V(s)]$

• $S \leftarrow S'$

Until s is terminal



The update rule uses a quintuple of events $(s_t, a_t, r_{t+1}, s_{t+1}, a_{t+1})$, therefore called SARSA.

Problem: Unlike in the deterministic case we can not choose A completely greedy policy $\pi(s)=\max_a Q(s,a)$, as due to the unknown transition and reward functions $\delta(s,a)$ and r(s,a) we can not be sure if another action might eventually turn out to be better.

ε-greedy policy

- Soft policy: π(s,a) > 0 for all s∈S, a∈A(s)
 Non-zero probability off choosing every possible action
- ε-greedy policy: Most of the time with probability (1-ε) follow the optimal policy

 $\pi(s) = \max_a Q(s,a)$

but with probability e pick a random action:

 $\pi(s,a) \ge \epsilon/|A(s)|$

 Let ε→0 go to zero as t→∞ for example ε=1/t so that ε-greedy policy converges to the optimal deterministic policy

SARSA Policy Iteration

Initialize Q(s,a) arbitrarily:

- Repeat for each episode
- Initialize s
- Choose a from using ϵ -greedy policy derived from Q
- Repeat for each step of episode
 - Take action a, observe reward r, and next state s'
 - $Q(s,a) \leftarrow Q(s,a) + \alpha [r + \gamma Q(s',a') Q(s,a)]$
 - s ← s', a ← a'
- Until s is terminal

Q-Learning (Off-Policy TD)

- Approximates the optimal value functions V*(s) or Q*(s,a) independent of the policy being followed.
- The policy determines which state-action pairs are visited and updated

 $Q(s_t, a_t) \leftarrow Q(s_t, a_t) + \alpha [r_{t+1} + \gamma \max_{a'} Q(s_{t+1}, a') - Q(s_t, a_t)]$

Q-Learning Off-Policy Iteration

Initialize Q(s,a) arbitrarily:

- Repeat for each episode
- Initialize s
- Choose a from s using ε -greedy policy derived from Q
- Repeat for each step of episode
 - Take action a, observe reward r, and next state s'
 - $Q(s,a) \leftarrow Q(s,a) + \alpha [r + \gamma \max_{a'}Q(s',a') Q(s,a)]$

■ S ← S′

Until s is terminal

TD versus Monte-Carlo

- So far TD uses one-step look-ahead but why not use 2-steps or n-steps look-ahead to update Q(s,a).
- 2-step look-ahead
- $Q(s_t, a_t) \leftarrow Q(s_t, a_t) + \alpha [r_{t+1} + \gamma r_{t+2} + \gamma^2 Q(s_{t+2}, a_{t+2}) Q(s_t, a_t)]$
- N-step look-ahead
- $\begin{aligned} Q(\mathbf{s}_{t}, \mathbf{a}_{t}) \leftarrow Q(\mathbf{s}_{t}, \mathbf{a}_{t}) + \alpha \left[\mathbf{r}_{t+1} + \gamma \mathbf{r}_{t+2} + \dots + \gamma^{n-1} \mathbf{r}_{t+n} + \gamma^{n} \\ Q(\mathbf{s}_{t+n}, \mathbf{a}_{t+n}) Q(\mathbf{s}_{t}, \mathbf{a}_{t}) \right] \end{aligned}$
- Monte-Carlo method

 $\begin{aligned} Q(s_t, a_t) \leftarrow Q(s_t, a_t) + \alpha \; [r_{t+1} + \gamma \; r_{t+2} + \ldots + \gamma^{N-1} \; r_{t+N} - Q(s_t, a_t)] \\ \text{(compute total reward } \mathsf{R}^{\mathsf{T}} \; \text{at the end of the episode)} \end{aligned}$

Temporal Difference Learning

Drawback of one-step look-ahead:

 Reward is only propagated back to the successor state (takes long time to finally propagate to the start state)

$$\begin{array}{c} \overbrace{s_{t}} \stackrel{r=0}{\longrightarrow} \overbrace{s_{t+1}} \stackrel{r=0}{\longrightarrow} \overbrace{s_{t+2}} \dots \stackrel{r=0}{\longrightarrow} \overbrace{s_{T-1}} \stackrel{r=100}{\longrightarrow} \overbrace{s_{T}} \stackrel{r=100}{\Longrightarrow} \stackrel{r=100}{\underbrace{s_{T}} \stackrel{r=10}{\underbrace{s_{T}} \stackrel{r=1}{\underbrace{s_{T}} \stackrel{r=10}{\underbrace{s_{T}} \stackrel{r=1}{\underbrace{s_{T}} \stackrel{r=1}{\underbrace{s_{T$$

Monte-Carlo Method

Drawback of Monte-Carlo method

- Learning only takes place after an episode terminated
- Performs a random walk until goal state is discovered for the first time as all state-action values seem equal
- It might take long time to find the goal state by random walk
- TD-learning actively explores state space if each action receives a small default penalty



Idea: blend TD learning with Monte-Carlo method Define:

$$R_{t}^{(1)} = r_{t+1} + \gamma V_{t}(s_{t+1})$$

$$R_{t}^{(2)} = r_{t+1} + \gamma r_{t+2} + \gamma^{2} V_{t}(s_{t+2})$$

. . .

$$\begin{split} \mathsf{R}_t^{(n)} &= \mathsf{r}_{t+1} + \gamma \ \mathsf{r}_{t+2} + \ldots + \gamma^{n-1} \mathsf{r}_{t+n} + \gamma^n \mathsf{V}_t(\mathsf{s}_{t+n}) \\ \text{The quantity } \mathsf{R}_t^{(n)} \text{ is called the n-step return at} \\ \text{time t.} \end{split}$$

TD(λ)

- n-step backup : $V_t(s_t) \leftarrow V_t(s_t) + \alpha [R_t^{(n)}-V_t(s_t)]$
- $TD(\lambda)$: use average of n-step returns for backup

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

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

(if s_T is a terminal state)

The weight of the n-step return decreases with a factor of $\boldsymbol{\lambda}$

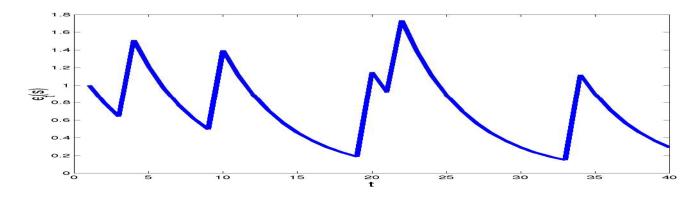
TD(0): one-step temporal difference method TD(1) : Monte-Carlo method

Eligibility Traces

Practical implementation of $TD(\lambda)$:

With each state-action pair associate an eligibility trace $e_t(s,a)$ On each step, the eligibility trace for all state-action pairs decays by a factor $\gamma\lambda$ and the eligibility trace for the one state and action visited on the step is incremented by 1.

$$\begin{split} e_t(s,a) &= \gamma \lambda \ e_{t-1}(s,a) + 1 & \text{ if } s = s_t \text{ and } a = a_t \\ &= \gamma \lambda \ e_{t-1}(s,a) & \text{ otherwise} \end{split}$$



On-line TD(λ)

Initialize Q(s,a) arbitrarily and e(s,a)=0 for all s,a: Repeat for each episode

- Initialize s,a
- Repeat for each step of episode
 - Take action a, observe r, s'
 - Choose a' from s' using policy derived from Q (ε-greedy)
 - $\delta \leftarrow r + \gamma Q(s',a') Q(s,a)$
 - e(s,a) ← e(s,a) +1

For all s,a:

- $Q(s,a) \leftarrow Q(s,a) + \alpha \delta e(s,a)$
- $e(s,a) \leftarrow \gamma \lambda e(s,a)$
- s ← s', a ← a'
- Until s is terminal

Function Approximation

- So far we assumed that the action value function Q(s,a) is represented as a table.
- Limited to problems with a small number of states and actions
- For large state spaces a table based representation requires large memory and large data sets to fill them accurately
- Generalization: Use any supervised learning algorithm to estimate Q(s,a) from a limited set of action value pairs
 - Neural Networks (Neuro-Dynamic Programming)
 - Linear Regression
 - Nearest Neighbors

Function Approximation

 Minimize the mean-squared error between training examples Q_t(s_t, a_t) and the true value function Q^π(s, a)

 $\sum_{s \in S} P(s) [Q^{\pi}(s,a) - Q_t(s_{t,\prime}a_t)]^2$

- Notice that during policy iteration P(s) and Q^π(s,a) change over time
- Parameterize Q^π(s,a) by a vector θ=(θ₁,..., θ_n) for example weights in a feed-forward neural network

Stochastic Gradient Descent

 Use gradient descent to adjust the parameter vector θ in the direction that reduces the error for the current example

 $\theta_{t+1} = \theta_t + \alpha \left[Q^{\pi}(s_t, a_t) - Q_t(s_t, a_t) \right] \nabla_{\theta t} Q_t(s_t, a_t)$

 The target output q_t of the t-th training example is not the true value of QP but some approximation of it.

$$\theta_{t+1} = \theta_t + \alpha \left[v_t - Q_t(s_t, a_t) \right] \nabla_{\theta_t} Q_t(s_t, a_t)$$

■ Still converges to a local optimum if $E\{v_t\}=Q^{\pi}(s_t,a_t)$ if $a \rightarrow 0$ for $t \rightarrow \infty$