

# Sequential Learning

## Lecture 4.5 : Summary of the first four courses

Rémy Degenne  
(remy.degenne@inria.fr)



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# Markov Decision Process

A MDP is parameterized by a tuple  $(\mathcal{S}, \mathcal{A}, R, P)$  where

- ▶  $\mathcal{S}$  is the **state space**
- ▶  $\mathcal{A}$  is the **action space**
- ▶  $R = (\nu_{(s,a)})_{(s,a) \in \mathcal{S} \times \mathcal{A}}$  where  $\nu_{(s,a)} \in \Delta(\mathbb{R})$  is the **reward distribution** for the state-action pair  $(s, a)$
- ▶  $P = (p(\cdot|s, a))_{(s,a) \in \mathcal{S} \times \mathcal{A}}$  where  $p(\cdot|s, a) \in \Delta(\mathcal{S})$  is the **transition kernel** associated to the state-action pair  $(s, a)$

In each (discrete) **decision time**  $t = 1, 2, \dots$ , a learning agent

- ▶ selects an **action**  $a_t$  based on his current **state**  $s_t$  (or possibly all the previous observations),
- ▶ gets a **reward**  $r_t \sim \nu_{(s_t, a_t)}$
- ▶ makes a transition to a **new state**  $s_{t+1} \sim p(\cdot|s_t, a_t)$

[Bellman 1957, Howard 1960, Blackwell 70s...]

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**Goal** : (made more precise later) select actions so as to maximize some notion of **expected cumulated rewards**

Mean reward of action  $a$  in state  $s$

$$r(s, a) = \mathbb{E}_{R \sim \nu_{(s,a)}}[R]$$

# Different Markov Decision Problems

**Overall goal** : learn the optimal policy  $\pi^*$  associated to some MDP parameterized by  $r(s, a)$  and  $p(\cdot|s, a)$  for  $(s, a) \in \mathcal{S} \times \mathcal{A}$ .

## Different contexts :

- 1 Small state space  $\mathcal{S}$ , known dynamics
- 2 Small state space  $\mathcal{S}$ , unknown dynamics
- 3 Large state space  $\mathcal{S}$ , known dynamics
- 4 Large state space  $\mathcal{S}$ , unknown dynamics

# Value and policy

## Value of a policy :

- ▶  $V^\pi(s) = \mathbb{E}^\pi \left[ \sum_{t=1}^{\infty} \gamma^{t-1} r_t \mid s_1 = s \right]$  and  
 $Q^\pi(s, a) = \mathbb{E}^\pi \left[ \sum_{t=1}^{\infty} \gamma^{t-1} r_t \mid s_1 = s, a_1 = a \right]$ .
- ▶  $V^*(s) = V^{\pi^*}(s) = \max_{\pi} V^\pi(s)$  and  
 $Q^*(s, a) = Q^{\pi^*}(s, a) = \max_{\pi} Q^\pi(s, a)$ .
- ▶  $V^\pi = \mathbb{E}_{a \sim \pi(s)} [Q^\pi(s, a)]$ .

## Greedy policy :

- ▶  $\text{greedy}(V) = \operatorname{argmax}_{a \in \mathcal{A}} \left( r(s, a) + \gamma \mathbb{E}_{s' \sim p(\cdot | s, a)} [V(s')] \right)$
- ▶  $\text{greedy}(Q) = \operatorname{argmax}_{a \in \mathcal{A}} Q(s, a)$
- ▶  $\pi^* = \text{greedy}(V^*)$  and  $\pi^* = \text{greedy}(Q^*)$  .

# Bellman equations and operators

The value of a policy satisfies a **Bellman equation**, written with the **Bellman operator**

$$V^\pi(s) = \mathbb{E}_{a \sim \pi(s)} [r(s, a) + \gamma \mathbb{E}_{s' \sim p(\cdot | s, a)} [V(s')]] , \\ V^\pi = T^\pi V^\pi .$$

Similar equations and operators for  $Q^\pi, V^*, Q^*$ .

**Properties of  $V^\pi$  and  $T^\pi$  :**

- ▶  $T^\pi$  is a  $\gamma$ -contraction
- ▶  $V^\pi$  is the unique fixed point
- ▶  $V_{n+1} = T^\pi V_n$  tends to  $V^\pi$ .

Similar properties for  $Q^\pi, V^*, Q^*$ .

# Goals

## Policy evaluation

Given a policy  $\pi$ , return  $V^\pi$  (or  $Q^\pi$ )

Often : find a “good enough” approximation of  $V^\pi$ .

## Finding the best policy

Find  $\pi^* = \operatorname{argmax}_\pi V^\pi = \operatorname{argmax}_\pi Q^\pi$ .

Property : there exists a deterministic  $\pi^*$ , given by  $\pi^* = \operatorname{greedy}(V^*)$  and  $\pi^* = \operatorname{greedy}(Q^*)$ .

Often : find a policy  $\pi$  which is “close enough” to  $\pi^*$ .

# Small MDP, known dynamics

**Solve the Bellman equation** for policy evaluation :  $V^\pi = (\mathbf{I} - \gamma\mathbf{P})^{-1}\mathbf{r}$ .

**Value iteration** for policy evaluation or finding the best policy :

- 1 Iterate  $V_{n+1} = T^\pi V_n$  (resp.  $V_{n+1} = T^* V_n$ )
- 2 Stop when  $\|V_{n+1} - T^\pi V_n\|$  is small

Then if we are iterating with  $T^*$  to find the best policy : return  $\pi = \text{greedy}(V_n)$ .

**Policy iteration** for finding the best policy :

- 1 Use policy evaluation to find  $V^{\pi_n}$
- 2 Perform policy improvement :  $\pi_{n+1} = \text{greedy}(V^{\pi_n})$ .

Both can also be performed with  $Q$  instead of  $V$ . (Advantages? Drawbacks?)



# Small MDP, unknown dynamics

Main ideas : Robbins-Monro estimation and temporal differences.

**TD(0)** for policy evaluation

- ▶  $\hat{V}(s_k) \leftarrow \hat{V}(s_k) + \alpha_{N(s_k)}(s_k) \delta(s_k)$  where  
 $\delta(s_k) = r_k + \gamma \hat{V}(s_{k+1}) - \hat{V}(s_k)$  and  $(r_k, s_{k+1}) = \text{step}(s_k, \pi)$ .

Parallel Robbins-Monro on each state.  $\hat{V}$  converges to  $V^\pi$  (under suitable conditions on  $\alpha$ , etc.).

**Q-Learning** for finding the best policy

- ▶  $Q(s, a) \leftarrow Q(s, a) + \alpha_{N(s,a)}(s, a) (r + \gamma \max_b Q(s', b) - Q(s, a))$   
where  $(r, s') = \text{step}(s, a)$
- ▶ Return  $Q, \pi = \text{greedy}(Q)$ .

Parallel Robbins-Monro on each state-action pair.  $Q$  converges to  $Q^*$ .  
Works for any behaviour policy, provided it explores enough.

Both are modified value iteration / policy iteration, with R-M and TD techniques to deal with unknown dynamics.

# Large MDP

**Function approximation.** Since  $\mathcal{F}(\mathcal{S}, \mathbb{R})$  is too large, introduce a (parametric) set of functions  $\mathcal{F}_V$  and look for best  $V$  in  $\mathcal{F}_V$ .

Ex : functions representable by a given neural network.

## Policy evaluation

- ▶ Minimize  $\text{MSVE}_\nu(V) = \mathbb{E}_{s \sim \nu} \left[ (V^\pi(s) - V(s))^2 \right]$ .
- ▶ Use TD(0) semi-gradient. Converges to  $\theta_{TD}$
- ▶ Or : estimate the solution directly with LSTD, using that  $A\theta_{TD} = b$  for some  $A, b$  (linear approximation). Variant for  $Q$  : LSTD-Q.

## Finding the best policy

- ▶ LSPI : policy iteration using LSTD-Q for policy evaluation
- ▶ Fitted Q-iteration : value iteration for  $Q$ , with regression to estimate  $T^*Q$  from samples
- ▶ Approximate Q-learning : use semi-gradient updates for  $Q$ .

And more to come, not value-based.