

Sequential Learning

Lecture 4.5 : Summary of the first four courses

Rémy Degenne
(remy.degenne@inria.fr)



Université
de Lille



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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...]

Markov Decision Process

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- ▶ $R = (\nu_{(s,a)})_{(s,a) \in \mathcal{S} \times \mathcal{A}}$ 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)

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** π^* 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^π, V^*, Q^* .

Properties of V^π and T^π :

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

Similar properties for Q^π, V^*, Q^* .

Goals

Policy evaluation

Given a policy π , return V^π (or Q^π)

Often : find a “good enough” approximation of V^π .

Finding the best policy

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

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

Often : find a policy π which is “close enough” to π^* .

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^{π_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^π (under suitable conditions on α , 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 θ_{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.