L16: Markov decision processes: exact offline solution

AIMA4e: Chapter 17.2.1–17.2.3 Kochenderfer: 7.2–7.7 (more detailed alternative to AIMA4e) and 7.8 (not covered in AIMA)

What you should know after this lecture

How to find optimal policies for MDPs!

- Value iteration
- Policy iteration
- Linear programming
- Linear quadratic regulators (for a class of continuous MDPs)

Probabilistic sequential decision-making

Probabilistic transitions Atomic, discrete states Full observability Solution is a policy



- Agent can observe current state completely and correctly
- World dynamics are probabilistic and known to the agent
- Agent selects actions to maximize expected summed rewards over time
- Agent plans on-line to select next action based on current state (but still potentially thinking about a longer horizon)

Solution strategies for MDPs

Two main categories

- <u>Online</u> action selection given current state s₀ via some form of search: last lecture
- Offline solution to derive a complete policy π that can be executed online very efficiently : **this lecture**

We will focus on the infinite-horizon discounted case with finite (and non-huge) S and A. Goal will be to compute Q^{*} or π^* for all s.

Recall from last time

$$\begin{split} V^{*}(s) &= \max_{a} \sum_{s'} P(s' \mid s, a) \left[R(s, a, s') + \gamma V^{*}(s') \right] \\ Q^{*}(s, a) &= \sum_{s'} P(s' \mid s, a) \left[R(s, a, s') + \gamma \max_{a'} Q^{*}(s', a') \right] \\ Q^{*}(s, a) &= \sum_{s'} P(s' \mid s, a) \left[R(s, a, s') + \gamma V^{*}(s') \right] \\ V^{*}(s) &= \max_{a} Q^{*}(s, a) \\ \pi^{*}(s) &= \operatorname*{argmax}_{a} Q^{*}(s, a) \end{split}$$

Policy evaluation

An important sub-problem: given policy π , what is the value of executing it?

$$V_{\pi}(s) = \sum_{s'} P(s' \mid s, \pi(s)) [R(s, \pi(s), s') + \gamma V_{\pi}(s')]$$

Note that:

- MDP + policy is a Markov chain
- Can define an iterative algorithm

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 \begin{array}{ll} \text{PolicyEvaluation}(\pi, \mathcal{S}, \mathcal{A}, \mathsf{T}, \mathsf{R}, \gamma, \varepsilon) \\ 1 & V_{\pi}(s) = 0 \text{ for } s \in \mathcal{S} \\ 2 & \textbf{while True:} \\ 3 & \textbf{for } s \in \mathcal{S} \\ 4 & V_{\pi, \text{new}}(s) = \sum_{s'} \mathsf{P}(s' \mid s, \pi(s)) \left[\mathsf{R}(s, \pi(s), s') + \gamma V_{\pi}(s')\right] \\ 5 & \textbf{if } |V_{\pi} - V_{\pi, \text{new}}| < \varepsilon \\ 6 & \textbf{return } V_{\pi, \text{new}} \\ 7 & V_{\pi} = V_{\pi, \text{new}} \end{array}
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- Observe that values are defined by a set of |S| linear equations in |S| unknowns (the $V_{\pi}(s)$ values). Let
 - V_{π} be a vector of the V_{π} values, for each state
 - T_{π} be a transition matrix, where $T_{ij} = P(s_j | s_i, \pi(s_i))$
 - + R be a reward vector (simplified to match standard treatment), where $R_i = R(s_i, \pi(s_i))$

Then

$$V_{\pi} = R + \gamma T_{\pi} V_{\pi}$$
$$(I - \gamma T_{\pi}) V_{\pi} = R$$
$$V_{\pi} = (I - \gamma T_{\pi})^{-1} R$$

Unfortunately, there's not such an easy solution for finding the optimal value function, because the max operations make the system non-linear.

Finding optimal policy: Stupidest possible algorithm

- Given finite s and a, there are finitely many policies! We could enumerate them, but how would we decide which one is best?
- Because of the Markov property (future depends only on s_t), there's a helpful theorem
- For any MDP there exists at least one deterministic optimal policy π^{*} such that for all other policies π,

 $V_{\pi^*}(s) \geqslant V_{\pi}(s)$

• This means we don't have to worry that some policies might be good in some parts of the state space and others good in other parts—there is at least one policy that's as good as the best policy at all states!

Finding optimal policy: Value iteration

Not so easy when we don't know the policy! System of equations isn't linear any more (it has max operations in it.)

$$V^*(s) = \max_{a} \sum_{s'} P(s' \mid s, a) \left[R(s, a, s') + \gamma V^*(s') \right]$$

ValueIteration($S, A, T, R, \gamma, \varepsilon$)

 $\begin{array}{ll} 1 & V(s) = 0 \text{ for } s \in \$ \\ 2 & \text{while True:} \\ 3 & \text{ for } s \in \$ \\ 4 & V_{new}(s) = \max_{a} \sum_{s'} \mathsf{P}(s' \mid s, a) \left[\mathsf{R}(s, a, s') + \gamma \mathsf{V}(s')\right] \\ 5 & \text{ if } |V_{\pi} - V_{new}| < \varepsilon \\ 6 & \text{ return } V_{new} \\ 7 & V = V_{new} \end{array}$

where $|V_1 - V_2| = \max_s |V_1(s) - V_2(s)|$.

Finding optimal policy: Q-Value iteration

For acting, it is more useful to have the Q(s, a) values.

$$Q^{*}(s, a) = \sum_{s'} P(s' \mid s, a) \left[R(s, a, s') + \gamma \max_{a'} Q^{*}(s', a') \right]$$

 $QValueIteration(\$, \mathcal{A}, \mathsf{T}, \mathsf{R}, \gamma, \varepsilon)$

$$\begin{array}{ll} 1 & Q(s,a) = 0 \text{ for } s \in \mathbb{S}, \, a \in \mathcal{A} \\ 2 & \text{while True:} \\ 3 & \text{ for } s \in \mathbb{S}, \, a \in \mathcal{A} \\ 4 & Q_{new}(s,a) = \sum_{s'} P(s' \mid s,a) \left[R(s,a,s') + \gamma \max_{a'} Q(s',a') \right] \\ 5 & \text{ if } |Q - Q_{new}| < \varepsilon \\ 6 & \text{ return } Q_{new} \\ 7 & Q = Q_{new} \end{array}$$

where $|Q_1-Q_2|=max_{s,\alpha}|Q_1(s,\alpha)-Q_2(s,\alpha)|.$

Cool facts about value iteration

- Guaranteed to converge to Q*
- Max-norm error $|Q Q^*|$ decreases monotonically per iteration
- Can initialize to any value
- When initialized to 0, iterations are finite-horizon value functions.
- Can execute "in place" (don't need a separate Q_{new})
- Can randomly pick (s, a) to update, rather than doing it systematically
- Serves as the basis for Q-learning
- If $|Q Q_{new}| < \varepsilon$ then $|Q Q^*| < \varepsilon \gamma / (1 \gamma)$
- Define greedy policy with respect to value function $\pi_Q(s) = argmax_a Q(s, a)$. Then if $|Q(s, a) Q^*(s, a)| < \varepsilon$, $|V_{\pi_Q} V^*| < 2\varepsilon$.

Gridworld domain¹

- Simple grid world with a goal state with reward 1 and a bad state with reward -100
- Actions move in the desired direction with probability 0.8, in one of the perpendicular directions with probability 0.1
- Taking an action that would bump into a wall leaves agent where it is.

0	0	0	1	Action = north
0		0	-100	P = 0.8
0	0	0	0	$P = 0.1 \checkmark P = 0.1$

Gridworld value iteration

Running value iteration with $\gamma=0.9$

One iteration

0	0	0.72	1.81
0		0	-99.91
0	0	0	0

• Five iterations

0.809	1.598	2.475	3.745
0.268		0.302	-99.59
0	0.034	0.122	0.004

Gridworld value iteration

Running value iteration with $\gamma = 0.9$

• Ten iterations

2.686	3.527	4.402	5.812
2.021		1.095	-98.82
1.390	0.903	0.738	0.123

1000 iterations

5.470	6.313	7.190	8.669
4.802		3.347	-96.67
4.161	3.654	3.222	1.526

Gridworld value iteration

Running policy iteration with $\gamma = 0.9$

• Resulting policy after 1000 iterations



Policy iteration

Actually, it usually happens that $\pi_Q = \pi^*$ long before Q is close to Q^{*}. So doing value iteration until convergence might be too much work. Let's try working explicitly in the space of policies without enumerating them all!

PolicyIteration $(S, A, T, R, \gamma, \epsilon)$ 1 $\pi(s) = a$ for $s \in S$ and an arbitrary $a \in A$ while True: 2 3 $Q = Q_{\pi}$ // Policy evaluation $\pi' = \pi_0$ // Greedy policy wrt Q 4 5 if $\pi = \pi'$: 6 return π 7 $\pi = \pi'$

- Worst-case complexity is bad, but often very good in practice.
- Interesting combinations of policy and value iteration (e.g., don't completely solve policy evaluation step—use some iterations of iterative policy evaluation instead.)

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Gridworld policy iteration

Running policy iteration with $\gamma=$ 0.9, initialized with policy $\pi(s)=North$

• One iteration

0.418	0.884	2.331	6.367
0.367		-8.610	-105.7
-0.168	-4.641	-14.27	-85.05

• Two iterations

5.414	6.248	7.116	8.634
4.753		2.881	-102.7
2.251	1.977	1.849	-8.701

Gridworld policy iteration

Running policy iteration with $\gamma=$ 0.9, initialized with policy $\pi(s)=North$

• Three iterations (converged)

5.470	6.313	7.190	8.669
4.803		3.347	-96.67
4.161	3.654	3.222	1.526

Gridworld results

- Approximation of value function
 - Policy iteration: exact value function after three iterations
 - Value iteration: after 100 iterations, $\|V V^*\|_2 = 7.1 \times 10^{-4}$
- Calculation of optimal policy
 - Policy iteration: three iterations
 - Value iteration: 12 iterations

In other words, value iteration converges to optimal policy long before it converges to correct value in this MDP (but, this property is highly MDP-specific)

Policy iteration or value iteration?

- Policy iteration requires fewer iterations that value iteration, but each iteration requires solving a linear system instead of just applying Bellman backup.
- In practice, policy iteration is often faster, especially if the transition probabilities are structured (e.g., sparse) to make solution of linear system efficient.
- <u>Modified policy iteration</u> (Puterman and Shin, 1978) solves linear system approximately, using backups very similar to value iteration, and often performs better than either value or policy iteration.

Linear programming

- Define variables V_i for the optimal value of state i.
- Minimize

$$\sum_{i} V_i$$

• Subject to, for all $s \in |S|$, $a \in |A|$

$$V_{i} \geq \sum_{s_{j}} P(s_{j} \mid s_{i}, a) [R(s_{i}, a, s_{j}) + \gamma V_{j}]$$

Linear programming

- Why do we minimize a weighted combination of the values? Shouldn't we maximize value?
- Value functions V that satisfy the constraints are <u>upper bounds</u> on the optimal value function V*

 $V(s) \geqslant V^*(s) \ \, \forall s$

• Minimizing value ensures that we choose the <u>lowest upper</u> bound

$$\min_{V} V(s) = V^*(s) \quad \forall s$$

Only known algorithm with worst-case time complexity polynomial in |S| and |A|. The complexity of policy iteration has a term that depends on $1/(1 - \gamma)$. In practice other methods are usually more efficient.

Linear quadratic regulator

There is no general-purpose solution method for continuous state and action MDPs. But there's an interesting and useful special case when the dynamics are linear-Gaussian (like the systems we did Kalman filtering on) and the rewards are quadratic.

- $S = \mathbb{R}^n$ (often called X)
- $\mathcal{A} = \mathbb{R}^{m}$ (often called controls U)
- $s' | s, a = T_s s + T_a a + W$ where T_s is an $n \times n$ matrix, T_a is an $n \times m$ matrix, and W is drawn from a zero-mean finite-variance Gaussian.
- $R(s, a) = s^T R_s s + a^T R_a a$ where R_s is $n \times n$ and positive semidefinite and R_a is $m \times m$ and positive definite. This penalizes states and actions that deviate from 0. You need to define your states s so that 0 is a desired state and actions a so that 0 is a desired action.

Linear quadratic regulator: solutions

Good to know that these exist, but we won't study them:

- In finite-horizon case, there is a value iteration method that finds the exact optimal control sequence, using the <u>dynamic Riccati</u> equation
- Even cooler, in the infinite-horizon (non-discounted) case, there is a stationary optimal policy of the form a = Fs where F is a fixed matrix (found by solving for a fixed-point of the Riccati equation).
- Even more cool, these same basic things work out in continuous time, where $\dot{s} = T_s s + T_a a + W$.

Next time

• Approximate value and policy iteration via reinforcement learning