Lecture 3 and 4: Dynamic programming and reinforcement learning

Reinforcement Learning and Decision Making MLSALT7, Lent 2016

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Dynamic programming for control

Recall in the last lecture that we wrote the value function recursively as

$$\nu^{\pi}(x) = r(x) + \gamma \int p^{\pi}(z|x) \, \nu^{\pi}(z) \, dz$$

which for discrete/tabular MDPs can be written as

$$\mathbf{v}^{\pi} = \mathbf{r} + \gamma \mathbf{P}^{\pi} \mathbf{v}^{\pi}.$$

Alternatively we can introduce the Bellman operator \mathbb{T}^{π} , the discrete version of which can be applied to any vector \mathbf{v} resulting in

$$\mathfrak{T}^{\pi}\boldsymbol{\nu} = \mathbf{r} + \gamma \mathbf{P}^{\pi}\boldsymbol{\nu}$$

for which $\mathbf{v}^{\pi} = \mathbf{T}^{\pi} \mathbf{v}^{\pi}$ is the unique fixed point.

The Bellman operator is a γ -contraction

We can look at the max-norm $\|\boldsymbol{\nu}\|_{\infty} = \max_i \nu_i$ of the Bellman operator:

$$\begin{split} \|\mathfrak{I}^{\pi}\boldsymbol{\nu}_{1} - \mathfrak{I}^{\pi}\boldsymbol{\nu}_{2}\|_{\infty} &= \|\mathbf{r} + \gamma \mathbf{P}^{\pi}\boldsymbol{\nu}_{1} - \mathbf{r} - \gamma \mathbf{P}^{\pi}\boldsymbol{\nu}_{2}\|_{\infty} \\ &= \gamma \|\mathbf{P}^{\pi}(\boldsymbol{\nu}_{1} - \boldsymbol{\nu}_{2})\|_{\infty} \\ &\leqslant \gamma \|\boldsymbol{\nu}_{1} - \boldsymbol{\nu}_{2}\|_{\infty} \end{split}$$

For $\gamma \in [0, 1)$ this shows that \mathbb{T}^{π} is a contraction under this norm and as a result due to the Banach fixed-point theorem, \mathbb{T}^{π} has a unique fixed-point.

This states that given any two vectors applying the operator to them will take them both closer towards <u>something</u>. That <u>something</u> can be found by repeatedly applying the contraction:

$$\mathbf{v}^{\mathfrak{i}+1} = \mathfrak{T}^{\pi} \mathbf{v}^{\mathfrak{i}}$$

which we have defined as the value function!

In slightly more detail we can look at the max-norm comparing the iterative procedure and the fixed-point:

$$\begin{split} \|\boldsymbol{\nu}^{i+1} - \boldsymbol{\nu}^{\pi}\|_{\infty} &= \|\mathcal{T}^{\pi} \boldsymbol{\nu}^{i} - \mathcal{T}^{\pi} \boldsymbol{\nu}^{\pi}\|_{\infty} \qquad \text{(the def'n of } \mathcal{T}^{\pi} \text{ and } \boldsymbol{\nu}^{\pi}) \\ &\leqslant \gamma \|\boldsymbol{\nu}^{i} - \boldsymbol{\nu}^{\pi}\|_{\infty} \qquad \text{(the } \gamma\text{-contraction)} \\ &\vdots \\ &\leqslant \gamma^{i+1} \|\boldsymbol{\nu}^{0} - \boldsymbol{\nu}^{\pi}\|_{\infty} \to 0 \end{split}$$

Let's say we have an arbitrary value function $\nu.$ How should we act optimally under this value function?

We can define a policy:

$$\pi'(x) = \operatorname*{arg\,max}_{a} \left[r_x + \gamma \mathbf{P}_{ax}^\top \boldsymbol{\nu} \right]$$

If $\mathbf{v} = \mathbf{v}^{\pi}$ is value function associated with policy π , do π and π' coincide?

No! The equation above looks similar to the Bellman operator introduced earlier, but because of the arg max above, $\nu_x^{\pi'} \ge \nu_x^{\pi}$ for all $x \in \mathcal{X}$.

The Optimal Bellman operator and value iteration

Define the optimal Bellman operator \mathfrak{T}^* as

$$(\mathfrak{T}^*\boldsymbol{\nu})_{\boldsymbol{x}} = \max_{\boldsymbol{\alpha}} \big[\mathbf{r}_{\boldsymbol{x}} + \gamma \mathbf{P}_{\boldsymbol{\alpha}\boldsymbol{x}}^{\top} \boldsymbol{\nu} \big].$$

This can similarly be shown to be a γ -contraction which has as its fixed-point the optimal value function $v^* = T^* v^*$.

The iterative procedure:

- **1** start from some arbitrary v
- **2** and iterate $\mathbf{v} \leftarrow \Upsilon^* \mathbf{v}$ until convergence

is known as value iteration, due to Bellman (1957). Convergence can be shown in exactly the same way as the convergence for \Im^{π}

What does the policy look like in this case? How would things change if the reward was defined as $r_{\alpha xz}$?

Backward induction

Value iteration is also known as backward induction. Why?

Let's assume that rather than considering an infinite-horizon problem we want to solve

$$a_0 = \operatorname*{arg\,max}_{\alpha} \mathbb{E}_{a_{1:T}, x_{1:T}} \Big[\sum_{t=1}^{I} r(x_t) \Big| x_0 \Big]$$

How could we use backward induction to solve this?

The key is in the name. We can start by solving the problem optimally at step T:

. . .

$$\begin{split} \nu_{\mathsf{T}}(x) &= \max_{a} \left[r_{x} + \mathbf{P}_{ax}^{\top} \mathbf{r} \right] & \text{ and performing induction} \\ \nu_{\mathsf{T}-1}(x) &= \max_{a} \left[r_{x} + \mathbf{P}_{ax}^{\top} \boldsymbol{\nu}_{\mathsf{T}} \right] & \end{split}$$

until ultimately $a_0 = \arg \max v_1$. This can be written recursively as $v_t = \mathcal{T}^* v_{t+1}$ where $v_{T+1} = r$.

We have already shown how to compute the value of a policy by finding the fixed-point $v^{\pi} = T^{\pi}v^{\pi}$. This is known as policy evaluation.

But now given the value of a policy we an improve upon that value (as shown in an earlier slide) by setting

$$\pi(\mathbf{x}) \leftarrow \arg \max_{\alpha} \left[\underbrace{\mathbf{r}_{\mathbf{x}} + \mathbf{P}_{\alpha \mathbf{x}}^{\top} \boldsymbol{\nu}^{\pi}}_{Q^{\pi}(\mathbf{x}, \alpha)} \right]$$

This is known as policy improvement. The function Q^{π} is often known as a state/action value-function and is not strictly necessary here since it follows directly from v^{π} .

After we have evaluated and updated the policy π how can we use this to find the optimal policy?

This interleaved process of policy evaluation and improvement is known as policy iteration:

- 1 Initialize ν arbitrarily
- **2** Iterate:

$$\begin{array}{l} \bullet \ \pi(\mathbf{x}) \leftarrow \arg \max_{\mathbf{a}} \left[\mathbf{r}_{\mathbf{x}} + \mathbf{P}_{a\mathbf{x}}^{\top} \mathbf{v} \right] \\ \bullet \ \mathbf{v} \leftarrow \lim_{\mathbf{n} \to \infty} (\mathfrak{I}^{\pi})^{\mathbf{n}} \mathbf{v} \end{array}$$

item most of these techniques can be gathered as generalized policy iteration where

- some improvement is made to the value function $\boldsymbol{\nu}$, and
- the policy π is updated to be greedy with respect to ν



Reinforcement learning

The approaches introduced previously, based on dynamic programming, relied on knowing the model p(z|x, a) in order to perform the necessary integrals.

- now we will only assume that we can sample z conditioned on (x, a)
- further we will only assume we can sample a single z!

The temporal difference (TD) error

Let's return to the recursive definition of the value function:

$$v^{\pi}(\mathbf{x}) = \mathbf{r}(\mathbf{x}) + \gamma \sum_{z} \mathbf{p}^{\pi}(z|\mathbf{x}) v^{\pi}(z)$$

if we can sample from $\mathbf{p}^{\pi}(z|\mathbf{x})$ then this can be approximated as

$$\approx r(x) + \gamma \frac{1}{N} \sum_i \nu^{\pi}(z^i) \quad \mathrm{for} \ z^i \sim p^{\pi}(\cdot|x)$$

but we will only assume we can take a single sample. IE we have no "reset switch" that will allow us to draw multiple samples from a single x. Once we've taken an action and moved to z that's it.

$$pprox \mathbf{r}(\mathbf{x}) + \gamma \mathbf{v}^{\pi}(z) \quad \text{for } z \sim \mathbf{p}^{\pi}(\cdot | \mathbf{x})$$

The difference between the left- and right-side is known as the temporal difference error. It should be zero in expectation!

TD for policy evaluation

Temporal difference methods use the TD-error (essentially) as a noisy gradient (think stochastic approximation)

Given a policy π and an arbitrary value function \mathbf{v} we can compute \mathbf{v}^{π} by iterating:

- **1** sample the next state $z \sim p^{\pi}(\cdot|\mathbf{x})$
- 2 update the value function,



3 and repeat: $\mathbf{x} \leftarrow \mathbf{z}$

But this is just policy evaluation. Can we combine this with policy improvement?

In order to apply TD for control we first have to note that we must directly learn Q rather than a value function. Why?

For DP-based methods policy improvement is

$$\pi(\mathbf{x}) = \operatorname*{arg\,max}_{a} Q(\mathbf{x}, \mathbf{a}) = \operatorname*{arg\,max}_{a} r_{\mathbf{x}} + \mathbf{P}_{a\mathbf{x}}^{\top} \boldsymbol{\nu}$$

only ν is needed because integration with respect to $p^{\pi}(z|x)$ can be performed. This doesn't work for RL.

Finally: the integration above also allows us to consider all outcomes of taking action a. For RL we must instead explore by injecting noise into our action selection.

On-policy TD: SARSA

- () initialize Q(x, a) arbitrarily
- 2 select action a arbitrarily
- 3 iterate
 - **1** sample $x' \sim p(\cdot|x, a)$
 - 2 choose action \mathfrak{a}' from Q " $\varepsilon\text{-greedily}$ ",

$$\mathfrak{a}' = \begin{cases} \arg\max_{\mathfrak{a}'} Q(\mathfrak{x}',\mathfrak{a}') & \text{w.p. } 1-\varepsilon, \\ \text{Uniform}(\mathcal{A}) & \text{otherwise.} \end{cases}$$

3 update the value function

$$Q(x, a) \leftarrow Q(x, a) + \alpha \big[r_x + \gamma \underbrace{Q(x', a')}_{-Q(x, a)} - Q(x, a) \big]$$

"on-policy" because α^\prime used here

Note: the funny name stands for "State-Action-Reward-State-Action"

- () initialize Q(x, a) arbitrarily
- 2 select action \mathfrak{a} arbitrarily
- 3 iterate
 - $\textbf{0} \text{ sample } x' \sim p(\cdot|x,a)$
 - 2 choose action \mathfrak{a}' from Q " \mathfrak{c} -greedily",
 - **3** update the value function

$$Q(x, a) \leftarrow Q(x, a) + \alpha \big[r_x + \gamma \max_{\substack{a'' \\ \text{"off-policy" due to } a''}} Q(x', a'') - Q(x, a) \big]$$

What is the difference between on- and off-policy methods?

- updates are performed without using actions selected by the exploratory policy for off-policy
- Q-learning can not care about rewards it gets while it's learning
 - we WILL choose bad actions \mathfrak{a}' due to the noisy exploration
 - but this does not affect the value function learned by Q-learning due to the max

no noise, but the optimal path passes by a cliff with high penalty for falling



Cliff-world results

