Lecture 4 and 5: Modern reinforcement learning—LSTD and policy gradients

Reinforcement Learning and Decision Making MLSALT7, Lent 2016

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Once again, the value function V can be defined as the unique fixed point of the Bellman operator, $V^{\pi} = T^{\pi}V^{\pi}$

$$(\mathfrak{I}^{\pi} V)(x) = \mathbf{r}(x) + \gamma \int_{\mathfrak{X}} \mathbf{P}(x'|x, \pi(x)) \, V(x') \, dx'$$

or more concisely, as

 $\mathfrak{T}^{\pi} V = r + \gamma P^{\pi} V.$

Instead of computing V^{π} for every state (which may not even be possible!) we will approximate this with a weighted combination of features $\phi : \mathcal{X} \to \mathbb{R}^k$, i.e.

$$V^{\pi}(\mathbf{x}) \approx \boldsymbol{\varphi}(\mathbf{x})^{\top} \boldsymbol{\theta}$$

Let's return to the definition of the value function, by way of the Bellman operator, and apply it to the approximator:

$$\mathfrak{T}^{\pi} \big[\varphi(x)^{\top} \theta \big] = \mathfrak{r}(x) + \gamma P^{\pi} \big[\varphi(x)^{\top} \theta \big]$$

What's wrong with applying the Bellman operator to this approximation?

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What's wrong with applying the Bellman operator to this approximation? More concretely: does there exist a θ^* such that $\Im^{\pi}[\varphi(x)^{\top}\theta] = \varphi(x)^{\top}\theta^*$? Let's go back to what it means to approximate some function. We want to find the best approximation under some norm,

$$f(x) \approx \varphi(x)^\top \theta^* \text{ where } \theta^* = \mathop{\arg\min}_{\theta} \|\varphi(x)^\top \theta - f(x)\|$$

However $\mathfrak{T}^{\pi} \Phi w$ may not necessarily lie in the span of Φ . Instead, we will introduce a projection operator Π such that

 $\Pi v = \Phi \arg\min_{u \in \mathbb{R}^k} \|\Phi u - v\|^2.$

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We can then define the following fixed point:

 $\Phi w = \Pi \mathfrak{T}^{\pi} \Phi w.$



First, the Bellman operator is a γ -contraction, i.e. for any y, z

$$\|\mathfrak{T}^{\pi}\mathbf{y} - \mathfrak{T}^{\pi}z\| \leqslant \gamma \|\mathbf{y} - z\|$$

and the projection operator is non-expansive, i.e. for any \boldsymbol{y}

 $\|\Pi y\|\leqslant \|y\|.$

Combining these two means that $\Pi \mathfrak{T}^{\pi}$ is a γ -contraction, and due to the Banach fixed-point theorem there exists a unique fixed point $\hat{v} = \Pi \mathfrak{T}^{\pi} \hat{v}$.

Note! This does not mean that there is a unique solution $\Phi w = \hat{v}$.

We can write the fixed point as the following:

$$w = \underset{u \in \mathbb{R}^{k}}{\arg\min} \|\Phi u - (r + \gamma P^{\pi} \Phi w)\|^{2}$$
$$= (\Phi^{T} \Phi)^{-1} \Phi^{T} (r + \gamma P^{\pi} \Phi w)$$
$$\vdots$$
$$= \underbrace{(\Phi^{T} (\Phi - \gamma P^{\pi} \Phi))^{-1}}_{A^{-1}} \underbrace{\Phi^{T} r}_{b}$$

The previous approach required forming the entire feature matrix Φ and also required the transition model P^{π} .

Instead we will assume samples $(x_i, \mathfrak{a}_i, x_i')$ generated on-policy and form

$$\hat{\Phi} = \begin{bmatrix} \varphi(x_1)^T \\ \vdots \\ \varphi(x_m)^T \end{bmatrix}, \ \hat{\Phi}' = \begin{bmatrix} \varphi(x_1')^T \\ \vdots \\ \varphi(x_m')^T \end{bmatrix}, \ \hat{r} = \begin{bmatrix} r(x_1) \\ \vdots \\ r(x_m) \end{bmatrix}$$

Solving for the fixed point is then given by

$$w = (\hat{\Phi}^{\mathsf{T}}(\hat{\Phi} - \gamma \hat{\Phi}'))^{-1} \hat{\Phi}^{\mathsf{T}} \hat{r}$$

The previous method is model-free for policy evaluation, but in order to improve the policy (moving from LSTD to LSPI) we would need a model.

Instead, learn the Q-function,

$$\begin{split} & Q^{\pi}(x, \mathfrak{a}) = \mathbb{E}\Big[\sum_{t=0}^{\infty} \gamma^{t} r(x_{t}) \Big| x_{0} = x, \mathfrak{a}_{0} = \mathfrak{a}, \pi \Big] \\ & \pi^{new}(x) = \operatorname*{arg\,max}_{\mathfrak{a}} Q^{\pi}(x, \mathfrak{a}) \end{split}$$

Learning the Q-function

Define the Bellman operator as

 $\mathsf{T}^{\pi}Q=r+\gamma\mathsf{P}\mathsf{H}^{\pi}Q$

where H^{π} is called Π^{π} in (Lagoudakis and Parr), and PH^{π} basically describes the probability of transitioning from $(x, a) \rightarrow (x', a')$.

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Now, when we move to the empirical version we have samples (x, a, x') not necessarily drawn on-policy. We construct

$$\hat{\Phi} = \begin{bmatrix} \varphi(x_1, \alpha_1)^T \\ \vdots \\ \varphi(x_m, \alpha_m)^T \end{bmatrix}, \ \hat{\Phi}' = \begin{bmatrix} \varphi(x'_1, \pi(x'_1))^T \\ \vdots \\ \varphi(x'_m, \pi(x'_m))^T \end{bmatrix}, \ \hat{r} = \begin{bmatrix} r(x_1) \\ \vdots \\ r(x_m) \end{bmatrix}.$$

The solution for w is the same as before.

Rather than indirectly parameterizing a policy via its value function we can directly parameterize the policy

 $\pi_{\theta}(a|x)$

with parameters θ . Note that now we'll go back to assuming a stochastic policy (we'll see why shortly) and return to the value of a full trajectory,

$$J(\boldsymbol{\theta}) = \mathbb{E}_{\boldsymbol{\theta}} \Big[\sum_{t=0}^{k} \gamma^{t} r(\boldsymbol{X}_{t}) \Big]$$

For a discrete space we can define policies of the form:

$$\pi_{\theta}(a|x) = \frac{\exp(\theta_{ax})}{\sum_{a'} \exp(\theta_{a'x})}$$

Or a common policy for continuous spaces is:

$$\pi_{\theta}(a|x) = Kx + m \quad \mathrm{for} \ \theta = (K,m)$$

A little notation

Let's let $\tau=(x_0,\ldots,x_k,a_0,\ldots,a_k)$ denote a single trajectory. How can we write the probability of $\tau?$

$$p_{\theta}(\tau) = p(x_0)\pi_{\theta}(a_0|x_0) \prod_{n=1}^{k} p(x_n|x_{n-1}, a_{n-1})\pi_{\theta}(a_n|x_n)$$

We can also write the reward for a trajectory as

$$R(\tau) = \sum_{t=0}^{k} \gamma^{t} r(x_{t})$$

for which our objective becomes

$$\mathbf{J}(\boldsymbol{\theta}) = \mathbb{E}_{\boldsymbol{\theta}}[\mathbf{R}(\boldsymbol{\tau})]$$

Now we can expand our gradient as follows:

$$\begin{split} \nabla J(\theta) &= \nabla \int R(\tau) \, p_{\theta}(\tau) d\tau \\ &= \int R(\tau) \, \nabla p_{\theta}(\tau) d\tau \\ &= \int R(\tau) \, p_{\theta}(\tau) \, \nabla \log p_{\theta}(\tau) d\tau \qquad \qquad \text{because } \nabla \log f = \nabla f/f \\ &= \int R(\tau) \, p_{\theta}(\tau) \Big[\sum_{n=0}^{k} \nabla \log \pi_{\theta}(a_{n}|x_{n}) \Big] \end{split}$$

- We need only sample trajectories
- We need to know the policy and its gradient, but that's fine because we decide on that

- MC approx. to the gradient can be quite noisy
- Looking more closely we can see that the gradient consists of summing over terms

 $\nabla \log \pi_{\theta}(A_n | X_n) r(X_t)$

- for t<n rewards at time t cannot be affected by actions that come after it
- we end up with expectations

$$\mathbb{E}_{\theta}[\nabla \log \pi_{\theta}(A_n | X_n)] \mathbb{E}_{\theta}[r(X_t)] = 0$$
Expectation of the score is zero

Exploting independence cont'd

$$R(\tau) \left[\sum_{n=0}^{k} \nabla \log \pi_{\theta}(A_{n} | X_{n}) \right]$$
$$= \sum_{n=0}^{k} \nabla \log \pi_{\theta}(A_{n} | X_{n}) \sum_{t=0}^{k} \gamma^{t} r(X_{t})$$

Expand the term inside the expectation...

Under expectation we can eliminate rewards for *t*<*n*

$$= \sum_{n=0}^{k} \nabla \log \pi_{\theta}(A_{n}|X_{n}) \sum_{t=n}^{k} \gamma^{t} r(X_{t})$$

Call this $R_{n}(\tau)$

Including those terms would only add noise!

Baselines

- Let $R_n(\cdot)$ be the sum of rewards after step n
- We can now write

$$\nabla J(\theta) \approx \frac{1}{N} \sum_{i=1}^{N} \sum_{n=0}^{k} R_n(\tau^{(i)}) \nabla \log \pi_{\theta}(a_n^{(i)} | x_n^{(i)})$$

for *N* sample trajectories $\tau^{(i)}$

- Following this gradient coincides with
 REINFORCE [Williams, 1992]
 - GPOMDP [Baxter and Bartlett, 2001]

 For the same reason that we eliminated rewards with *t*<*n*, we can write the reward as

$$R_n(\tau^{(i)}) = \sum_{t=n}^k r(x_t^{(i)}) - b_t$$

Variance of gradient depends on magnitude of rewards.

- for **baseline** quantities independent of the *n*th action
 - can depend on previous rewards! (say the average)
 - this baseline can then be selected to reduce variance
 - see [Greensmith et al., 2001], [Riedmiller, et al, 2008]