Chapter 3

Control

3.1 Motivation

In the last chapter, we covered Markov Decision Processes (MDPs) where the state and action spaces were finite. We showed that, if we know the state transitions, we can calculate the optimal policy using efficient polynomial-time algorithms (Value Iteration and Policy Iteration).

What about the case where state and action spaces are infinite? This doesn't necessarily mean that they need to be super "complex", in a heuristic sense. They could simply be continuous, such as choosing the angle to tilt your steering wheel, or the force on a joint of a robotic drone.

Iterative in computation time, not samples. (Running for loop to compute some quantity as quickly as possible.)

Example 3.1.1: CartPole

Consider a pole balanced on a cart. The state consists of just four continuous values:

- 1. The position of the cart;
- 2. The velocity of the cart;
- 3. The angle of the pole;
- 4. The angular velocity of the pole.

The *control* we apply is a force on the cart, moving it left and right.

Goal: To stabilize the cart around an ideal state s^* .

If you've ever tried to balance a pen upright in the palm of your hand, this is essentially the same problem!

How do we formulate the cost function $c: \mathcal{S} \times \mathcal{A} \to \mathcal{S}$ for this problem? We want

something that we can easily generalize to other sorts of tasks. One way we can do so is simply by choosing a cost function that is *quadratic* in its arguments:

$$c(s_t, a_t) = a_t^{\top} R a_t + (s_t - s^{\star})^{\top} Q(s_t - s^{\star}). \tag{3.1}$$

For some intuition into this expression, consider the simple case where a_t and s_t are one-dimensional, so $c(s_t, a_t) = ra_t^2 + q(s_t - s^*)^2$.

This expression has the nice property that for any continuous cost function, we can write out its *Taylor approximation* around a given point. Then, we can think of the quadratic cost function we've written above as the second-order approximation to that smooth cost function.

Remark 3.1.1: Quadratic forms

If this notation is unfamiliar to you, we recommend checking out this video from Khan Academy!

Remark 3.1.2: Notation

We call actions *controls*. In the literature, these are commonly represented with the letter u instead of a, but here we'll stick to a in order to illustrate the similarities with the discrete case of RL.

Definition 3.1.1: Optimal control problem

$$\min_{\pi_0, \dots, \pi_{T-1}: \mathcal{S} \to \mathcal{A}} \mathbb{E} \left[\sum_{t=0}^{T-1} c(s_t, a_t) \mid s_{t+1} = f(s_t, a_t), s_0 \sim \mu_0 \right]$$
(3.2)

The function $f: \mathcal{S} \times \mathcal{A} \to \mathcal{S}$ is the analogue of the state transitions. However, this might not be totally deterministic; there might be some underlying noise, or there might also be some measurement error in the state.

Our goal is to find a control policy π_t , which minimizes the *total cost* (undiscounted) over some finite number of steps T.

Note that this is a pretty hard problem the way it's written right now! We have some pretty strict constraints in the form of the state transitions.

How does this relate to the finite horizon case? If s_t and a_t were discrete, then we'd be able to work backwards using the dynamic programming algorithm we saw before.

As a matter of fact, let's consider what happens if we discretize the problem. Recall that $s \in \mathbb{R}^d$ and $a \in \mathbb{R}^k$, which are continuous. To make them discrete, let's round them to the nearest multiple of ϵ (for some choice of ϵ), so now the set of possible s and a is discrete. (Formally, we call this new set of rounded values an ϵ -grid over the

original continuous space. For example, if $\epsilon = 0.01$, then we're just rounding to two decimal spaces.)

If both these state and action spaces can be bounded, then the resulting sets are actually finite, so now we can use our previous tools for MDPs! But is this actually a feasible solution? Even if our S and A are finite, it might still be unfeasible to run the existing algorithms.

Indeed, this happens to be the case. Suppose our state and action spaces are bounded by some constants $\max_{s \in \mathcal{S}} ||s|| \leq B_s$ and $\max_{a \in \mathcal{A}} ||a|| \leq B_a$. Then to form our ε -net, we must divide each dimension into intervals of length ε , resulting in $(B_s/\varepsilon)^d$ and $(B_a/\varepsilon)^k$ points!

To get a sense of how quickly this grows, let's consider $\varepsilon = 0.01, d = k = 10$. Then the number of elements in our transition matrix $|\mathcal{S}|^2|\mathcal{A}|$, is on the order of $(100^{10})^2(100^{10}) = 10^{60}!$ Try finding a computer that'll fit that in memory!

So as we've seen, discretizing the problem isn't a feasible solution as soon as our action and state spaces are even moderately high-dimensional. How can we do better? Well, when doing the discretization, we implicitly relied on the assumption that rounding our value by some tiny amount ε wouldn't change the behavior much; namely, that the functions involved are relatively *continuous*. Can we use this structure in other ways? This brings us to the next topic, where we make some simplifying assumptions:

3.2 The Linear Quadratic Regulator Problem

Definition 3.2.1: Linear quadratic regulator problem

Linear dynamics: assume that the state transitions are linear with respect to the inputs:

$$s_{t+1} = f(s_t, a_t, w_t) = As_t + Ba_t + w_t$$

Note that we've also assumed that f is time-homogeneous: our state transitions behave the same way across all time steps.

Quadratic cost function:

$$c(s_t, a_t) = \begin{cases} s_t^\top Q s_t + a_t^\top R a_t & t < T \\ s_T^\top Q s_T & t = T \end{cases}$$

We want c to be a convex function so that there actually exists a minimum. This means that Q and R must both be positive definite.

In other words, we assume that we only take T actions total, and just ignore a_T .

Gaussian noise: $w_t \sim \mathcal{N}(0, \Sigma)$

Putting everything together, the optimization problem we want to solve is:

$$\min_{\substack{\pi_0, \dots, \pi_{T-1}: \mathcal{S} \to \mathcal{A}}} \quad \mathbb{E}\left[\left(\sum_{t=0}^{T-1} s_t^\top Q s_t + a_t^\top R a_t\right) + s_T^\top Q s_T\right]$$
s.t.
$$s_{t+1} = A s_t + B a_t + w_t$$

$$a_t = \pi_t(s_t)$$

$$w_t \sim \mathcal{N}(0, \Sigma)$$

$$s_0 \sim \mu_0$$

It might seem like we're oversimplifying, but in fact, like we mentioned above, one way to think of these simplifications is that we're just taking the best linear approximation to some general f and c. This is part of the reason why LQR is so well-studied. In fact, humans might even design systems to be linear in order to use results from LQR!

Of course, this only works if the state transitions and cost functions we're approximating are *roughly* smooth or quadratic respectively. For more complex, nonlinear systems, this basic design will break down. But later on, we'll see that we can generalize these ideas to get surprisingly good solutions.

Example 3.2.1: Driving down a road

Suppose we're driving down a road. At each time step, we can choose an action a_t : either we accelerate and apply a force forward $(a_t > 0)$, or reverse and apply a force backward $(a_t < 0)$. Suppose we can choose an action every δ seconds, and that our car has mass m.

Recall that Newtonian mechanics says that force = mass \times acceleration. We can write the acceleration as the change in velocity over time, and write the velocity as the change in position over time:

acceleration_t =
$$\frac{v_t - v_{t-1}}{\delta}$$

 $v_t = \frac{p_t - p_{t-1}}{\delta}$

How should we construct our state? We want to express everything in terms of these linear dynamics, and we also want our state to be Markov, so that we can apply dynamic programming like before. Then if we write our state as consisting of the position and velocity, then we can write

$$p_{t+1} = p_t + \delta v_t$$
$$v_{t+1} = v_t + \frac{\delta}{m} a_t$$

Writing everything out in matrix notation, we get:

$$s_{t+1} = \begin{bmatrix} 1 & \delta \\ 0 & 1 \end{bmatrix} \begin{bmatrix} p_t \\ v_t \end{bmatrix} + \begin{bmatrix} 0 \\ \frac{\delta}{m} \end{bmatrix} a_t$$

Let's derive a more compressed form for the state at time t as a summation over past time steps. Note that

$$s_{t} = As_{t-1} + Ba_{t-1} + w_{t-1}$$

$$= A(As_{t-2} + Ba_{t-2} + w_{t-2}) + Ba_{t-1} + w_{t-1}$$

$$= \cdots$$

$$= A^{t}s_{0} + \sum_{i=0}^{t-1} A^{i}(Ba_{t-i-1} + w_{t-i-1})$$

Let's consider the expected value of the state at this time. Since we assume that $\mathbb{E} w_t = 0$ (this is the zero vector in d dimensions), by linearity of expectation, the w_t term vanishes, and so we're left with

$$\mathbb{E}[s_t \mid s_{0:t-1}, a_{0:t-1}] = A^t s_0 + \sum_{i=0}^{t-1} A^i B a_{t-i-1}.$$

So now we have a good overview of the LQR setting. How can we now define an optimal policy in this setting? Recall that we allow our policies to be *time-dependent*.

It turns out that the optimal policy is one that is deterministic and *linear* at each time step! That is,

$$\pi_t^{\star}(s_t) = -K_t s_t.$$

We'll prove this more formally in Theorem 3.2.2. This should remind you somewhat of the way in which the optimal policy in the previous MDP setting was stationary and deterministic. In both cases, it turns out that the optimal policy has special structure!

Note that the average state at time t for the optimal policy is then

$$\mathbb{E}[s_t \mid s_0, a_t = -K_t s_t] = \left(\prod_{i=0}^{t-1} (A - BK_i)\right) s_0.$$

This introdces the quantity $A - BK_i$, which will show up frequently in discussions of LQR! For example, one important question is: will s_t remain bounded, or will it go to infinity as time goes on? We can answer this by analyzing this quantity $A - BK_i$, in particular its largest eigenvalue. Intuitively, if we imagine that these K_i s are equal (call this matrix K), then this expression looks like $(A - BK)^t s_0$. Now consider the maximum eigenvalue of A - BK, which we denote as λ_{max} . If $\lambda_{\text{max}} > 1$, then there's some initial state s_0^* for which

$$(A - BK)^t s_0^{\star} = \lambda_{\max}^t s_0^{\star} \xrightarrow{t \to \infty} \infty.$$

Definition 3.2.2: Value functions for LQR

Given a policy $\pi = (\pi_0, \dots, \pi_{t-1})$, we can define the value function $V_t^{\pi} : \mathcal{S} \to \mathbb{R}$ as

$$V_t^{\pi}(s) = \mathbb{E}\left[\sum_{i=t}^T c(s_i, a_i)\right]$$

$$= \mathbb{E}\left[\left(\sum_{i=t}^{T-1} s_i^{\top} Q s_i + a_i^{\top} R a_i\right) + s_T^{\top} Q s_T\right]$$
where $s_t = s$

$$a_i = \pi_i(s_i) \quad \forall i > t.$$

We call this expression inside the equation the **cost-to-go**, since it's just the total cost starting from timestep t.

Similarly, the Q function just additionally conditions on the first action we take:

$$Q_t^{\pi}(s, a) = \mathbb{E}\left[\sum_{i=t}^{T} c(s_i, a_i)\right]$$

$$= \mathbb{E}\left[\left(\sum_{i=t}^{T-1} s_i^{\top} Q s_i + a_i^{\top} R a_i\right) + s_T^{\top} Q s_T\right]$$
where $(s_t, a_t) = (s, a)$

$$a_i = \pi_i(s_i) \quad \forall i > t$$

As it turns out, we can now solve for the optimal policy π via dynamic programming in terms of these value and action-value functions.

3.2.1 Optimality for LQR

Definition 3.2.3: Optimal value functions for LQR

The optimal value function is the one that, in all states and across all timesteps, achieves *lowest cost* across all policies:

$$V_t^{\star}(s) = \min_{\pi} V_t^{\pi}(s)$$

$$= \min_{\pi_{t:T-1}} \mathbb{E}\left[\left(\sum_{i=t}^{T-1} s_t^{\top} Q s_t + a_t^{\top} R a_t\right) + s_T^{\top} Q s_T\right]$$
where $a_i = \pi_i(s_i) \quad \forall i \ge t$

$$s_t = s$$

Additionally, we'll show theorems 3.2.1 and 3.2.2 below, showing that the V_t^* is quadratic and that π_t^* is linear. Then, we'll show how to calculate the actual coefficients that specify these functions.

Theorem 3.2.1: V_t^{\star} in LQR is a quadratic function

Formally, we claim that

$$V_t^{\star}(s) = s^{\top} P_t s + p_t$$

for some $P_t \in \mathbb{R}^{d \times d}$ and $p_t \in \mathbb{R}^d$ where P_t is positive-definite. Note that this doesn't have a linear term, just a quadratic term plus a constant.

Theorem 3.2.2: Optimal policy in LQR is linear

That is,

$$\pi_t^{\star}(s) = -K_t s$$

for some $K_t \in \mathbb{R}^{k \times d}$. (The negative is just there by convention.)

We'll derive these theorems by induction, starting from the last timestep and working backwards in time. Note that induction has a very fundamental connection with dynamic programming: our inductive proof will naturally lend itself to a DP algorithm that allows us to calculate the optimal value and policy!

Base case: $V_T^{\star}(s)$ is quadratic.

Inductive hypothesis: Show that if $V_{t+1}^{\star}(s)$ is quadratic, then:

- 1. $Q_t^{\star}(s, a)$ is quadratic (in both s and a)
- 2. Derive the optimal policy $\pi_t^{\star}(s) = \arg\min_a Q_t^{\star}(s, a)$, and show that it's linear.
- 3. Show $V_t^{\star}(s)$ is quadratic.

Finally, this will have shown that $V_t^{\star}(s)$ is quadratic and $\pi_t^{\star}(s)$ is linear.

This is essentially the same proof as a finite-horizon MDP, except that now the state and action are *continuous* instead of finite.

Base case. Let's start by considering the final timestep V_T^{π} , for some policy π . Then the only expression is

$$V_T^{\star}(s) = s^{\top} Q s,$$

which is quadratic, as we desired. Pattern-matching to the expression from earlier, we see that $P_T = Q$ and $p_t = 0$.

Inductive step. Assume $V_{t+1}^{\star}(s) = s^{\top} P_{t+1} s + p_{t+1}$ for all states s. We'll start off by demonstrating that $Q_t^{\star}(s)$ is quadratic. Recall that the definition of $Q_t^{\star}: \mathcal{S} \times \mathcal{A} \to \mathbb{R}$ is

$$Q_t^{\star}(s, a) = c(s, a) + \mathop{\mathbb{E}}_{s' \sim f(s, a, w_{t+1})} V_{t+1}^{\star}(s').$$

We know $c(s, a) := s^{\top}Qs + a^{\top}Ra$. Let's consider the average value over the next timestep. The only randomness in the dynamics comes from the noise w_{t+1} , so we can write out this expected value as:

$$\mathbb{E}_{s' \sim f(s, a, w_{t+1})} V_{t+1}^{\star}(s') = \mathbb{E}_{w_{t+1} \sim \mathcal{N}(0, \sigma^2 I)} V_{t+1}^{\star}(As + Ba + w_{t+1})$$

$$= \mathbb{E}_{w_{t+1}} [(As + Ba + w_{t+1})^{\top} P_{t+1}(As + Ba + w_{t+1}) + p_{t+1}].$$

Summing these two expressions and combining like terms, we get

$$Q_{t}^{\star}(s,a) = s^{\top}Qs + a^{\top}Ra + \underset{w_{t+1}}{\mathbb{E}} [(As + Ba + w_{t+1})^{\top}P_{t+1}(As + Ba + w_{t+1}) + p_{t+1}]$$

$$= s^{\top}(Q + A^{\top}P_{t+1}A)s + a^{\top}(R + B^{\top}P_{t+1}B)a + 2s^{\top}A^{\top}P_{t+1}Ba + p_{t+1}$$

$$\underset{w_{t+1}}{\mathbb{E}} w_{t+1}^{\top}P_{t+1}w_{t+1}.$$

Now consider this last term. By writing out the product and using linearity of expectation, we can write this out as

$$\mathbb{E}_{w_{t+1}} w_{t+1}^{\top} P_{t+1} w_{t+1} = \sum_{i=1}^{d} \sum_{j=1}^{d} (P_{t+1})_{i,j} \mathbb{E}_{w_{t+1}} [(w_{t+1})_i (w_{t+1})_j].$$

When dealing with these quadratic forms, it's often helpful to consider the terms on the diagonal separately from those off the diagonal. On the diagonal, the expectation becomes $\mathbb{E}(w_{t+1})_i^2 = \text{Var}\left((w_{t+1})_i\right) = \sigma^2$. Off the diagonal, since the elements of w_{t+1} are independent, the expectation factors into $\mathbb{E}(w_{t+1})_i \mathbb{E}(w_{t+1})_j = 0$. Thus, the only terms left are the ones on the diagonal, so the sum of these can be expressed as the trace of $\sigma^2 P_{t+1}$:

$$\mathbb{E}_{w_{t+1}} w_{t+1}^{\top} P_{t+1} w_{t+1} = \text{Tr}(\sigma^2 P_{t+1}).$$

Substituting this back into the expression for Q_t^{\star} , we have:

Theorem 3.2.3: Optimal Q-Function in LQR

$$Q_t^{\star}(s,a) = s^{\top}(Q + A^{\top}P_{t+1}A)s + a^{\top}(R + B^{\top}P_{t+1}B)a + 2s^{\top}A^{\top}P_{t+1}Ba + \operatorname{Tr}(\sigma^2P_{t+1}) + p_{t+1}.$$

As we'd hoped, this expression is quadratic in s and a! (Phew!)

Now let's move on to the next part of the next part of proving the inductive hypothesis: showing that $\pi_t^*(s) = \arg\min_a Q_t^*(s, a)$ is linear. This becomes easy if Q_t^* is convex w.r.t. $a \dots$ Which it is!

Theorem 3.2.4: Q_t^{\star} is convex in a

Consider the part of Theorem 3.2.3 that is quadratic in a, namely $a^{\top}(R+B^{\top}P_{t+1}B)a$. Then Q_t^{\star} is convex w.r.t. a if $R+B^{\top}P_{t+1}B$ is positive definite.

To show this, recall that in our definition of LQR, we assumed that R is positive definite (see Definition 3.2.1). Also note that $B^{\top}P_{t+1}B$ is symmetric, and therefore positive definite. Since the sum of two positive-definite matrices is also positive-definite, we have that $R + B^{\top}P_{t+1}B$ is positive-definite, and so Q_t^{\star} is convex w.r.t. a.

This means that finding the minimum is easy: we can just take the gradient w.r.t. a and set it to zero! First, we calculate the gradient:

$$\nabla_a Q_t^{\star}(s, a) = \nabla_a [a^{\top} (R + B^{\top} P_{t+1} B) a + 2s^{\top} A^{\top} P_{t+1} B a]$$

= $2(R + B^{\top} P_{t+1} B) a + (2s^{\top} A^{\top} P_{t+1} B)^{\top}$

Setting this to zero, we get

$$0 = (R + B^{\top} P_{t+1} B) a + B^{\top} P_{t+1} A s$$

$$\pi_t^{\star}(s) := a = -(R + B^{\top} P_{t+1} B)^{-1} B^{\top} P_{t+1} A s$$

$$= -K_t s,$$
(3.3)

where
$$K_t = (R + B^{\top} P_{t+1} B)^{-1} B^{\top} P_{t+1} A$$
.

We're now almost there! To complete our inductive proof, we must show that the inductive hypothesis is true at time t; that is, we must prove that $V_t^{\star}(s)$ is quadratic. Using the identity $V_t^{\star}(s) = Q_t^{\star}(s, \pi^{\star}(s))$, we have:

$$V_t^*(s) = Q_t^*(s, \pi^*(s))$$

$$= s^{\top} (Q + A^{\top} P_{t+1} A) s + (-K_t s)^{\top} (R + B^{\top} P_{t+1} B) (-K_t s) + 2s^{\top} A^{\top} P_{t+1} B (-K_t s)$$

$$+ \operatorname{Tr}(\sigma^2 P_{t+1}) + p_{t+1}$$

Note that w.r.t. s, this is the sum of a quadratic term and a constant, which is exactly what we were aiming for! The constant term is clearly $p_t = \text{Tr}(\sigma^2 P_{t+1}) + p_{t+1}$. We can simplify the quadratic term by substituting in K_t . Notice that when we do this, the $(R + B^{T} P_{t+1} B)$ term in the expression is cancelled out by its inverse, and the remaining terms combine to give what is known as the *Ricatti equation*:

Theorem 3.2.5: Ricatti equation

$$P_t = Q + A^{\top} P_{t+1} A - A^{\top} P_{t+1} B (R + B^{\top} P_{t+1} B)^{-1} B^{\top} P_{t+1} A.$$

There are several nice things to note about this expression:

- 1. It's defined recursively; Given P_T , A, B, and the state coefficients Q, we can recursively calculate all values of P_t across timesteps.
- 2. It appears frequently in calculations surrounding optimality, such as in V^* and Q^* .
- 3. Together with A, B, and the action coefficients R, it fully defines the optimal policy.

The optimal policy also has some interesting properties: in addition to being independent of the starting distribution μ_0 , which also happened for our finite-horizon MDP solution, it's fully deterministic and doesn't depend on any noise! (Compare this with the discrete MDP case, where calculating our optimal policy required taking an expectation over the state transitions.)