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6.231 Dynamic Programming and Stochastic Control
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6.231 DYNAMIC PROGRAMMING

LECTURE 23

LECTURE OUTLINE

- Review of indirect policy evaluation methods
- Multistep methods, LSPE(λ)
- LSTD(λ)
- Q -learning
- Q -learning with linear function approximation
- Q -learning for optimal stopping problems

REVIEW: PROJECTED BELLMAN EQUATION

- For a fixed policy μ to be evaluated, consider the corresponding mapping T :

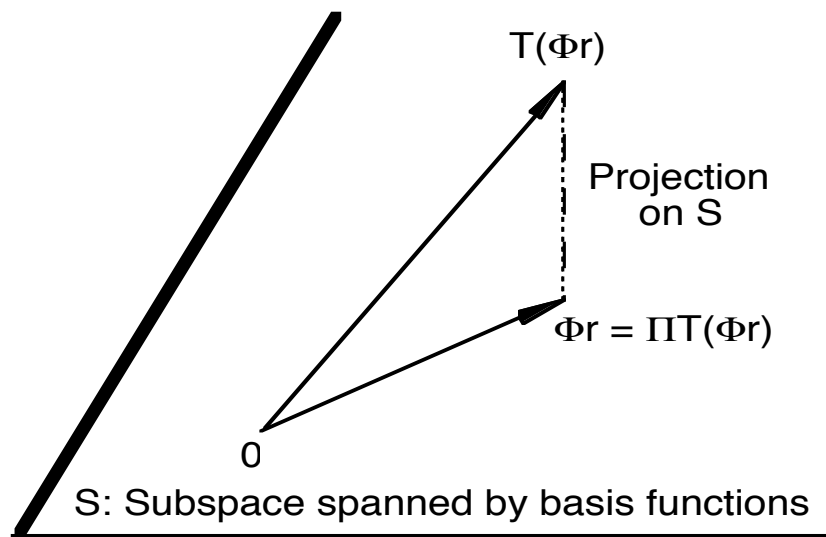
$$(TJ)(i) = \sum_{j=1}^n p_{ij} (g(i, j) + \alpha J(j)), \quad i = 1, \dots, n,$$

or more compactly,

$$TJ = g + \alpha PJ$$

- The solution J_μ of Bellman's equation $J = TJ$ is approximated by the solution of

$$\Phi r = \Pi T(\Phi r)$$



Indirect method: Solving a projected form of Bellman's equation

PVI/LSPE

- **Key Result:** ΠT is contraction of modulus α with respect to the weighted Euclidean norm $\|\cdot\|_\xi$, where $\xi = (\xi_1, \dots, \xi_n)$ is the steady-state probability vector. The unique fixed point Φr^* of ΠT satisfies

$$\|J_\mu - \Phi r^*\|_\xi \leq \frac{1}{\sqrt{1 - \alpha^2}} \|J_\mu - \Pi J_\mu\|_\xi$$

- **Projected Value Iteration (PVI):** $\Phi r_{k+1} = \Pi T(\Phi r_k)$, which can be written as

$$r_{k+1} = \arg \min_{r \in \mathcal{R}^s} \|\Phi r - T(\Phi r_k)\|_\xi^2$$

or equivalently

$$r_{k+1} = \arg \min_{r \in \mathcal{R}^s} \sum_{i=1}^n \xi_i \left(\phi(i)'r - \sum_{j=1}^n p_{ij} (g(i, j) + \alpha \phi(j)'r_k) \right)^2$$

- **LSPE (simulation-based approximation):** We generate an infinite trajectory (i_0, i_1, \dots) and update r_k after transition (i_k, i_{k+1})

$$r_{k+1} = \arg \min_{r \in \mathcal{R}^s} \sum_{t=0}^k (\phi(i_t)'r - g(i_t, i_{t+1}) - \alpha \phi(i_{t+1})'r_k)^2$$

JUSTIFICATION OF PVI/LSPE CONNECTION

- By writing the necessary optimality conditions for the least squares minimization, PVI can be written as

$$\left(\sum_{i=1}^n \xi_i \phi(i) \phi(i)' \right) r_{k+1} = \left(\sum_{i=1}^n \xi_i \phi(i) \sum_{j=1}^n p_{ij} (g(i, j) + \alpha \phi(j)' r_k) \right)$$

- Similarly, by writing the necessary optimality conditions for the least squares minimization, LSPE can be written as

$$\left(\sum_{t=0}^k \phi(i_t) \phi(i_t)' \right) r_{k+1} = \left(\sum_{t=0}^k \phi(i_t) (g(i_t, i_{t+1}) + \alpha \phi(i_{t+1})' r_k) \right)$$

- So LSPE is just PVI with the two expected values approximated by simulation-based averages.
- Convergence follows by the law of large numbers.
- The bottleneck in rate of convergence is the law of large of numbers/simulation error (PVI is a contraction with modulus α , and converges fast relative to simulation).

LEAST SQUARES TEMP. DIFFERENCES (LSTD)

- Taking the limit in PVI, we see that the projected equation, $\Phi r^* = \Pi T(\Phi r^*)$, can be written as $Ar^* + b = 0$, where

$$A = \sum_{i=1}^n \xi_i \phi(i) \left(\alpha \sum_{j=1}^n p_{ij} \phi(j) - \phi(i) \right)'$$
$$b = \sum_{i=1}^n \xi_i \phi(i) \sum_{j=1}^n p_{ij} g(i, j)$$

- A, b are expected values that can be approximated by simulation: $A_k \approx A, b_k \approx b$, where

$$A_k = \frac{1}{k+1} \sum_{t=0}^k \phi(i_t) \left(\alpha \phi(i_{t+1}) - \phi(i_t) \right)'$$
$$b_k = \frac{1}{k+1} \sum_{t=0}^k \phi(i_t) g(i_t, i_{t+1})$$

- **LSTD method:** Approximates r^* as

$$r^* \approx \hat{r}_k = -A_k^{-1} b_k$$

- Conceptually very simple ... but less suitable for optimistic policy iteration (hard to transfer info from one policy evaluation to the next).
- Can be shown that convergence rate is the same for LSPE/LSTD (for large k , $\|r_k - \hat{r}_k\| \ll \|r_k - r^*\|$).

MULTISTEP METHODS

- Introduce a multistep version of Bellman's equation $J = T^{(\lambda)} J$, where for $\lambda \in [0, 1)$,

$$T^{(\lambda)} = (1 - \lambda) \sum_{t=0}^{\infty} \lambda^t T^{t+1}$$

- Note that T^t is a contraction with modulus α^t , with respect to the weighted Euclidean norm $\|\cdot\|_{\xi}$, where ξ is the steady-state probability vector of the Markov chain.
- From this it follows that $T^{(\lambda)}$ is a contraction with modulus

$$\alpha_{\lambda} = (1 - \lambda) \sum_{t=0}^{\infty} \alpha^{t+1} \lambda^t = \frac{\alpha(1 - \lambda)}{1 - \alpha\lambda}$$

- T^t and $T^{(\lambda)}$ have the same fixed point J_{μ} and

$$\|J_{\mu} - \Phi r_{\lambda}^*\|_{\xi} \leq \frac{1}{\sqrt{1 - \alpha_{\lambda}^2}} \|J_{\mu} - \Pi J_{\mu}\|_{\xi}$$

where Φr_{λ}^* is the fixed point of $\Pi T^{(\lambda)}$.

- The fixed point Φr_{λ}^* depends on λ .
- Note that $\alpha_{\lambda} \downarrow 0$ as $\lambda \uparrow 1$, so error bound improves as $\lambda \uparrow 1$.

PVI(λ)

$$\Phi r_{k+1} = \Pi T^{(\lambda)}(\Phi r_k) = \Pi \left((1 - \lambda) \sum_{t=0}^{\infty} \lambda^t T^{t+1}(\Phi r_k) \right)$$

or

$$r_{k+1} = \arg \min_{r \in \mathfrak{R}^s} \left\| \Phi r - T^{(\lambda)}(\Phi r_k) \right\|_{\xi}^2$$

- Using algebra and the relation

$$(T^{t+1} J)(i) = E \left\{ \alpha^{t+1} J(i_{t+1}) + \sum_{k=0}^t \alpha^k g(i_k, i_{k+1}) \mid i_0 = i \right\}$$

we can write PVI(λ) as

$$r_{k+1} = \arg \min_{r \in \mathfrak{R}^s} \sum_{i=1}^n \xi_i \left(\phi(i)' r - \phi(i)' r_k - \sum_{t=0}^{\infty} (\alpha \lambda)^t E \{ d_k(i_t, i_{t+1}) \mid i_0 = i \} \right)^2$$

where

$$d_k(i_t, i_{t+1}) = g(i_t, i_{t+1}) + \alpha \phi(i_{t+1})' r_k - \phi(i_t)' r_k,$$

are the, so called, temporal differences (TD) - they are the errors in satisfying Bellman's equation.

LSPE(λ)

- Replacing the expected values defining PVI(λ) by simulation-based estimates we obtain LSPE(λ).
- It has the form

$$r_{k+1} = \arg \min_{r \in \mathbb{R}^s} \sum_{t=0}^k \left(\phi(i_t)' r - \phi(i_t)' r_k - \sum_{m=t}^k (\alpha \lambda)^{m-t} d_k(i_m, i_{m+1}) \right)^2$$

where (i_0, i_1, \dots) is an infinitely long trajectory generated by simulation.

- Can be implemented with convenient incremental update formulas (see the text).
- Note the λ -tradeoff:
 - As $\lambda \uparrow 1$, the accuracy of the solution Φr_λ^* improves - the error bound to $\|J_\mu - \Phi r_\lambda^*\|_\xi$ improves.
 - As $\lambda \uparrow 1$, the “simulation noise” in the LSPE(λ) iteration (2nd summation term) increases, so longer simulation trajectories are needed for LSPE(λ) to approximate well PVI(λ).

Q-LEARNING I

- Q-learning has two motivations:
 - Dealing with multiple policies simultaneously
 - Using a model-free approach [no need to know $p_{ij}(u)$ explicitly, only to simulate them]
- The Q-factors are defined by

$$Q^*(i, u) = \sum_{j=1}^n p_{ij}(u) \left(g(i, u, j) + \alpha J^*(j) \right), \quad \forall (i, u)$$

- In view of $J^* = TJ^*$, we have $J^*(i) = \min_{u \in U(i)} Q^*(i, u)$ so the Q factors solve the equation

$$Q^*(i, u) = \sum_{j=1}^n p_{ij}(u) \left(g(i, u, j) + \alpha \min_{u' \in U(j)} Q^*(j, u') \right), \quad \forall (i, u)$$

- $Q(i, u)$ can be shown to be the unique solution of this equation. **Reason:** This is Bellman's equation for a system whose states are the original states $1, \dots, n$, together with all the pairs (i, u) .
- Value iteration:

$$Q(i, u) := \sum_{j=1}^n p_{ij}(u) \left(g(i, u, j) + \alpha \min_{u' \in U(j)} Q(j, u') \right), \quad \forall (i, u)$$

Q-LEARNING II

- Use any probabilistic mechanism to select sequence of pairs (i_k, u_k) [all pairs (i, u) are chosen infinitely often], and for each k , select j_k according to $p_{i_k j}(u_k)$.
- At each k , Q-learning algorithm updates $Q(i_k, u_k)$ according to

$$Q(i_k, u_k) := (1 - \gamma_k(i_k, u_k)) Q(i_k, u_k) + \gamma_k(i_k, u_k) \left(g(i_k, u_k, j_k) + \alpha \min_{u' \in U(j_k)} Q(j_k, u') \right)$$

- Step size $\gamma_k(i_k, u_k)$ must converge to 0 at proper rate (e.g., like $1/k$).
- **Important mathematical point:** In the Q-factor version of Bellman's equation the order of expectation and minimization is reversed relative to the ordinary cost version of Bellman's equation:

$$J^*(i) = \min_{u \in U(i)} \sum_{j=1}^n p_{ij}(u) (g(i, u, j) + \alpha J^*(j))$$

- Q-learning can be shown to converge to true/exact Q-factors (a sophisticated proof).
- **Major drawback:** The large number of pairs (i, u) - no function approximation is used.

Q-FACTOR APPROXIMATIONS

- Introduce basis function approximation for Q -factors:

$$\tilde{Q}(i, u, r) = \phi(i, u)'r$$

- We cannot use LSPE/LSTD because the Q -factor Bellman equation involves minimization/multiple controls.

- An optimistic version of LSPE(0) is possible:

- Generate an infinitely long sequence $\{(i_k, u_k) \mid k = 0, 1, \dots\}$.

- At iteration k , given r_k and state/control (i_k, u_k) :

(1) Simulate next transition (i_k, i_{k+1}) using the transition probabilities $p_{i_k j}(u_k)$.

(2) Generate control u_{k+1} from the minimization

$$u_{k+1} = \arg \min_{u \in U(i_{k+1})} \tilde{Q}(i_{k+1}, u, r_k)$$

(3) Update the parameter vector via

$$r_{k+1} = \arg \min_{r \in \mathfrak{R}^s} \sum_{t=0}^k \left(\phi(i_t, u_t)'r - g(i_t, u_t, i_{t+1}) - \alpha \phi(i_{t+1}, u_{t+1})'r_k \right)^2$$

Q-LEARNING FOR OPTIMAL STOPPING

- Not much is known about convergence of optimistic LSPE(0).
- Major difficulty is that the projected Bellman equation for Q -factors may not be a contraction, and may have multiple solutions or no solution.
- There is one important case, **optimal stopping**, where this difficulty does not occur.
- Given a Markov chain with states $\{1, \dots, n\}$, and transition probabilities p_{ij} . We assume that the states form a single recurrent class, with steady-state distribution vector $\xi = (\xi_1, \dots, \xi_n)$.
- At the current state i , we have two options:
 - Stop and incur a cost $c(i)$, or
 - Continue and incur a cost $g(i, j)$, where j is the next state.
- Q -factor for the continue action:

$$Q(i) = \sum_{j=1}^n p_{ij} \left(g(i, j) + \alpha \min \{ c(j), Q(j) \} \right) \triangleq (FQ)(i)$$

- **Major fact:** F is a contraction of modulus α with respect to norm $\| \cdot \|_{\xi}$.

LSPE FOR OPTIMAL STOPPING

- Introduce Q -factor approximation

$$\tilde{Q}(i, r) = \phi(i)'r$$

- PVI for Q -factors:

$$\Phi r_{k+1} = \Pi F(\Phi r_k)$$

- LSPE

$$r_{k+1} = \left(\sum_{t=0}^k \phi(i_t)\phi(i_t)' \right)^{-1} \sum_{t=0}^k \phi(i_t) \left(g(i_t, i_{t+1}) + \alpha \min \{ c(i_{t+1}), \phi(i_{t+1})'r_k \} \right)$$

- Simpler version: Replace the term $\phi(i_{t+1})'r_k$ by $\phi(i_{t+1})'r_t$. The algorithm still converges to the unique fixed point of ΠF (see H. Yu and D. P. Bertsekas, “A Least Squares Q-Learning Algorithm for Optimal Stopping Problems”).