

Fitted Dynamic Programming

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**CS/Stat 184(0): Introduction to Reinforcement Learning
Fall 2024**

Today

- Feedback from last lecture
- Recap
- Neural networks
- Fitted value iteration
- Fitted policy iteration

Feedback from feedback forms

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Linear regression parameterizes $f(x)$ as $x^\top \theta$ and can work well when $\mathbb{E}[y | x]$ very smooth, high-dimensional (penalties like ridge/lasso help here), and/or there is a good featurization $\phi(x)$

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Parameter vector θ concatenates all W 's and b 's; $\dim(\theta)$ scales as $\text{width}^2 \times \text{depth}$

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Unfortunately, $L(\theta)$ is **non-convex**, i.e., it will in general have many local optima

We hope that SGD finds a good one... in practice there are optimization tricks that are like SGD but perform better, e.g., one very popular one is called **Adam**

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Practical Neural Networks are very far from “just” ERM

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Recall that Bellman equations state that the optimal value function $V^*(s)$ satisfies:

$$V^*(s) = \max_a \left\{ r(s, a) + \gamma \mathbb{E}_{s' \sim P(\cdot | s, a)} [V^*(s')] \right\}, \quad \forall s$$

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And the VI algorithm is a fixed-point algorithm to find V^* :

1. Initialization: $V^0(s) = 0, \forall s$

2. For $t = 0, \dots, T - 1$

$$V^{t+1}(s) = \max_a \left\{ r(s, a) + \gamma \sum_{s' \in \mathcal{S}} P(s' | s, a) V^t(s') \right\}, \quad \forall s$$

3. Return: $V^T(s)$

$$\pi(s) = \arg \max_a \left\{ r(s, a) + \gamma \mathbb{E}_{s' \sim P(\cdot | s, a)} V^T(s') \right\}$$

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For $h = H - 1, \dots, 0$, set:

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The above DP algorithm can just be seen as solving SH (Bellman) equations for the SH different values of $V(s, h)$, but doing so in an exact, efficient way via DP

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Recall from HW1 the Bellman equations for Q^* :

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Analogous Q-value DP, with same notational change as previous slide: h as argument

1. Initialization: $Q(s, a, H) = 0 \quad \forall s, a$

2. Solve (via dynamic programming):

$$Q(s, a, h) = r(s, a) + \mathbb{E}_{s' \sim P(s, a)} \left[\max_{a' \in A} Q(s', a', h + 1) \right] \quad \forall s, a, h$$

3. Return:

$$\pi_h(s) = \arg \max_a \left\{ Q(s, a, h) \right\}$$

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- We have N trajectories $\tau_1, \dots, \tau_N \sim \rho_{\pi_{data}}$

Each trajectory is of the form $\tau_i = \{s_0^i, a_0^i, \dots, s_{H-1}^i, a_{H-1}^i, s_H^i\}$

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Then we'd be happy if we found a

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Setting that aside for the moment, to fit supervised learning, we'd minimize a least-

squares objective function: $\hat{f}(x) = \arg \min_{f \in \mathcal{F}} \sum_{i=1}^{NH} (y_i - f(x_i))^2$

Connection to Supervised Learning (cont'd)

We can convert our data $\tau_1, \dots, \tau_N \sim \rho_{\pi_{data}}$, into (y, x) pairs; how many? NH

BUT, to compute each y , we need to already know Q !

Setting that aside for the moment, to fit supervised learning, we'd minimize a least-

squares objective function: $\hat{f}(x) = \arg \min_{f \in \mathcal{F}} \sum_{i=1}^{NH} (y_i - f(x_i))^2$

Then if we have enough data, choose a good \mathcal{F} , and optimize well,

$$Q(s_h, a_h, h) := \hat{f}(x) \approx \mathbb{E}[y | x] = \mathbb{E} \left[r(s_h, a_h) + \max_{a'} Q(s_{h+1}, a', h + 1) \mid s_h, a_h, h \right]$$

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Input: **offline dataset** $\tau_1, \dots, \tau_N \sim \rho_{\pi_{data}}$

1. Initialize fitted Q function at f_0

2. For $k = 1, \dots, K$:

$$f_k = \arg \min_{f \in \mathcal{F}} \sum_{i=1}^N \sum_{h=1}^{H-1} \left(f(s_h^i, a_h^i, h) - \left(r(s_h^i, a_h^i) + \max_a f_{k-1}(s_{h+1}^i, a, h+1) \right) \right)^2$$

3. With f_K as an estimate of Q^* , return $\pi_h(s) = \arg \max_a \left\{ f^K(s, a, h) \right\}$

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Q-Learning is an online version, i.e., draw new trajectories at each k based on f_k as Q -function

Today

- ✓ • Feedback from last lecture
- ✓ • Recap
- ✓ • Neural networks
- ✓ • Fitted value iteration
 - Fitted policy iteration

Recall: Policy Iteration (PI)

- Initialization: choose a policy $\pi^0 : S \mapsto A$

- For $k = 0, 1, \dots$

1. **Policy Evaluation:** Solve (via dynamic programming):

$$Q^{\pi^k}(s, a, h) = r(s, a) + \mathbb{E}_{s' \sim P(\cdot | s, a)} \left[Q^{\pi^k}(s', \pi^k(s), h + 1) \right] \quad \forall s, a, h$$

2. **Policy Improvement:** set $\pi_h^{k+1}(s) := \arg \max_a Q^{\pi^k}(s, a, h)$

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Yes! RHS can be written as $\mathbb{E} \left[r(s_h, a_h) + Q^{\pi^k}(s_{h+1}, \pi^k(s_h), h + 1) \mid s_h, a_h, h \right]$

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Spot the difference!

Fitted Policy Evaluation

Use exact same strategy as before: **fixed point iteration**

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Input: policy π , dataset $\tau_1, \dots, \tau_N \sim \rho_\pi$

1. Initialize fitted Q^π function at f_0

2. For $k = 1, \dots, K$:

$$f_k = \arg \min_{f \in \mathcal{F}} \sum_{i=1}^N \sum_{h=1}^{H-1} \left(f(s_h^i, a_h^i, h) - \left(r(s_h^i, a_h^i) + f_{k-1}(s_{h+1}^i, \pi(s_h^i), h + 1) \right) \right)^2$$

3. Return the function f_K as an estimate of Q^π

Fitted Policy Iteration:

- Initialization: choose a policy $\pi^0 : S \mapsto A$ and a sample size N
- For $k = 0, 1, \dots$
 1. **Fitted Policy Evaluation:** Using N sampled trajectories $\tau_1, \dots, \tau_N \sim \rho_{\pi^k}$, obtain approximation $\hat{Q}^{\pi^k} \approx Q^{\pi^k}$
 2. **Policy Improvement:** set $\pi_h^{k+1}(s) := \arg \max_a \hat{Q}^{\pi^k}(s, a, h)$

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Using the definition of the Q function, can do a **non-iterative** fitted policy evaluation

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Input: policy π , dataset $\tau_1, \dots, \tau_N \sim \rho_\pi$

Return:

$$\hat{Q}^\pi = \arg \min_{f \in \mathcal{F}} \sum_{i=1}^N \sum_{h=1}^{H-1} \left(f(s_h^i, a_h^i, h) - \sum_{t=h}^{H-1} r(s_t^i, a_t^i) \right)^2$$

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Summary:

- Neural Networks work well for complex function approximation with big data
- Incorporating supervised learning into PI and VI makes them RL algorithms!

Attendance:

bit.ly/3RcTC9T



Feedback:

bit.ly/3RHtlxy

