Learning for policy improvement

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joint work w/ Wen Sun

Learning from ≪*109 steps*

- •Many successes of RL: need billions of training examples
- What if experience is expensive?
	- ‣ e.g., robot in real world
	- ▶ e.g., expert supervision
	- ‣ e.g., personalization
	- ‣ e.g., safety
- Idea: replace "SGD-like" algorithms (e.g., policy gradient) with algorithms based on *policy iteration*
- Different tradeoff between computation and data
	- ‣ compute better update directions, take bigger steps
	- "think before you act"

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Motivation

- Consider TD(*λ*) (and later algorithms based on it, like DQN)
- Each iteration: one new experience, one step of SGD
	- ‣ couples optimization efficiency and sample efficiency
	- ‣ if optimizer is slow, uses more data
- What if we did more computation on a minibatch of samples to determine a better update direction?
	- ‣ if done well: bigger updates, fewer total samples
	- ‣ if done poorly: don't update policy often enough, collected data is less relevant

(Exact) policy iteration

- Do at least once:
	- ‣ for all states s, actions a
		- \triangleright calculate current total exp. cost $Q^{\pi}(s, a)$, value $\forall \pi(s) = E_{a \sim \pi(s)}[Q\pi(s, a)]$, and (dis)advantage $A^{\pi}(s, a) = Q^{\pi}(s, a) - V^{\pi}(s)$
	- choose $\pi^{new}(s) = \argmin_{a} A^{\pi}(s, a)$

// evaluate

// improve

- Doesn't work in a real-size problem:
	- ‣ must sample (s, a) rather than iterating over all
	- \triangleright can't calculate A^{π} exactly, must estimate somehow
	- ‣ can't choose new policy freely, must work in some hypothesis class

proximate policy improvement

(meta-algorithm)

- Do at least once:
	- \triangleright estimate $A^{\pi}(s, a)$
	- \triangleright train π' to achieve low E[A π (s, a)]
	- adjust π toward π ' to get π ^{new}
- *// evaluate*
- *// improve*

- To instantiate: way to estimate $A^{\pi}(s, a)$, train π' , update π^{new}
	- \triangleright also starting π , stopping criterion

Simple analysis of approx PI

- Guarantee: cost of π^{new} is $\sqrt{\pi(s_0)} + T$ $E_{new}[A^{\pi}(s, a)]$
	- ‣ performance difference lemma
		- ighthrouporting sum: if we follow π we expect $\sqrt{\pi(s_0)}$; each time we instead take an action a $\neq \pi(s)$ we gain/lose $A^{\pi}(s, a)$
	- \triangleright improvement when $E_{\text{new}}[A^{\pi}(s, a)] < 0$ (i.e., π improvable, hypothesis class rich enough, and training succeeds)
- Difficulty: expectation is under distribution of (s, a) from π^{new} (not the distribution we used to collect data)
- Can we develop algorithms that guarantee improvement (w/
assumptions) despite this difficulty? Yes... assumptions) despite this difficulty?

Two routes to improvement

- •Small updates: if new policy is close enough to old, then difference between new/old gradients doesn't matter
	- ‣ "SGD-like" analysis
	- ‣ pro: frequent updates mean we can get to relevant policies faster
- Big updates: if we can guarantee $E_{\text{new}}[A^{\pi}(s, a)] < 0$ whp, doesn't matter how close new/old policies are
	- ‣ pro: might be able to do better optimizing each minibatch
- Range of algorithms, from aggressive to deliberate updates

Ex: natural policy gradient

- Exponential family policy $\pi(s, a) \sim \exp(\phi(s, a) \cdot w)$
	- ighthoroup Fig. is features $\phi(s, a)$, weights w, probs. normalized to sum to 1 at each s
- Evaluation step: fit $A^{\pi}(s, a) \approx \phi(s, a) \cdot v$ by regression
	- \triangleright compatible function approximation (same ϕ)
	- ▶ data: run π, collect $\phi(s, a) \rightarrow$ [total cost after taking $(s, a)] \hat{V}^{\pi}(s)$
- Improvement/update step: boosting
	- \rightarrow w \leftarrow w $-\eta v$ (step size η)

ok to use any estimate, even a biased one

try all acts, or try

a random one

and use IPS

- i.e., scale $\pi(s, a)$ by exp($-\eta \phi(s, a) \cdot v$), then renormalize
- •Analysis: v estimates *natural gradient* of total cost wrt w
	- improvement for small enough η , vs. noise level (batch size) and smoothness

Variation: policy boosting

- Functional gradient version of same algorithm
- Instead of linear model $\phi(s, a)$ · v, fit $A^{\pi}(s, a)$ from any function class (decision trees, deep nets, …) — say, at iteration t, $A_t(s, a)$
- Policy proportional to $exp(-\eta_1A_1 \eta_2A_2 ...)$

Ex: conservative policy iteration

- Policy: mixture of classifiers from hypothesis space *^H*
- Evaluation/improvement step: cost-sensitive classification
	- train π ' from *H* to minimize E[A π (s, a)] with s $\sim \pi$, a $\sim \pi$ ' '
	- sample s $\sim \pi$, a \sim arbitrary, record advantage $A^{\pi}(s, a)$ and score 1 / P(a | s)
	- e.g., $a \sim \pi$ and argmax regression: same setup as for natural gradient
- Update step: mixture
	- $\blacktriangleright \pi(s, a) \leftarrow (1-\eta) \pi(s, a) + \eta \pi'(s, a)$
- Analysis: π ' minimizes dot product w/ functional policy gradient
	- ‣ so, update step is Frank-Wolfe
	- nonconvex objective, but convergence guaranteed with, e.g., $\eta = 1$ /iter

Connection: imitation learning

- \bullet Let π be the expert policy, run one iteration of approx. policy improvement to find π ^{new} that does at least as well
- Note: if expert is suboptimal, learner might do strictly better
	- in fact, a single step of policy iteration is often very powerful
	- ‣ e.g., suggests that a *random* expert can be good enough for successful IL, as long as it reaches goals occasionally

Ex: AggreVaTe

- •Policy: mixture of classifiers from hypothesis space *^H*
- •Evaluation/improvement step: cost-sensitive *no-regret* classification
	- \triangleright train π ' from $\Delta(H)$ to minimize E_{π} '[A π (s, a)]: i.e., on its *own* distribution
- Update step:
	- ‣ slowly mix from expert to learned policy (original paper)
	- or wait and set $\pi \leftarrow \pi'$ when confident of improvement (justified by perf. diff. lemma)

Update step

- If we mix over to new policy slowly enough, improvement is guaranteed (analysis in original AggreVaTe paper)
- If we switch all at once, can be confident of improvement when observed advantage is sufficiently negative
	- ‣ compare no-regret guarantee to performance difference lemma $V_{\pi}(s_0) - V_{\pi}(s_0) = T E_{\pi}[A^{\pi}(s, a)]$
- Either way, don't have to use mixture policy (random selection from iterations of no-regret)
	- ‣ best component of the mixture is at least as good as the mixture itself

Variation: AggreVaTeD

- Pick online gradient descent as no-regret learner
- Pick a deep net as function representation (even though noregret property is not guaranteed in this case)
- Ignore the expert after initialization
	- ▶ learn initial policy by behavioral cloning
	- ‣ then aggressively update expert to be current policy (small minibatches)

AggreVaTeD on AI gym 2d walker

- Single PI step, starting from policy from TRPO (overnight)
- AggreVaTeD (regular and natural gradient versions) improves much faster than from-scratch RL, beats original expert

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Aggrevated On PUNDP *AggreVaTeD on POMDP*

- •Fun use of IL: let the expert cheat
- POMDP version of acrobot (can see only joint angles not angular velocities)
- Expert sees everything

Summary

- •Sample-efficient RL via approximate policy improvement
- •Family of RL algorithms (different ways to estimate advantages, represent step directions, improve policies; aggressive to deliberate updates)
	- ‣ gave several examples
	- ‣ many more seem possible
- Future work
	- ‣ more thorough exploration and analysis of possibilities
	- ‣ combine these algorithms with exploration mechanisms