From Policy Gradient to Actor-Critic methods The policy gradient derivation (2/3)

Olivier Sigaud

Sorbonne Université http://people.isir.upmc.fr/sigaud



Limits of Algorithm 1

- Needs a large batch of trajectories or suffers from large variance
- ▶ The sum of rewards is not much informative
- Computing R from complete trajectories is not the best we can do

$$\begin{split} \nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}) &\sim & \frac{1}{m} \sum_{i=1}^{m} \sum_{t=1}^{H} \nabla_{\boldsymbol{\theta}} \mathrm{log} \pi_{\boldsymbol{\theta}}(\mathbf{a}_{t}^{(i)} | \mathbf{s}_{t}^{(i)}) R(\boldsymbol{\tau}^{(i)}) \\ &\sim & \frac{1}{m} \sum_{i=1}^{m} \sum_{t=1}^{H} \nabla_{\boldsymbol{\theta}} \mathrm{log} \pi_{\boldsymbol{\theta}}(\mathbf{a}_{t}^{(i)} | \mathbf{s}_{t}^{(i)}) [\sum_{k=1}^{H} r(\mathbf{s}_{k}^{(i)}, \mathbf{a}_{k}^{(i)})] \\ &* \text{split into two parts} \end{split}$$

$$\sim \frac{1}{m} \sum_{i=1}^{m} \sum_{t=1}^{H} \nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_{t}^{(i)} | \mathbf{s}_{t}^{(i)}) [\sum_{k=1}^{t-1} r(\mathbf{s}_{k}^{(i)}, \mathbf{a}_{k}^{(i)}) + \sum_{k=t}^{H} r(\mathbf{s}_{k}^{(i)}, \mathbf{a}_{k}^{(i)})]$$

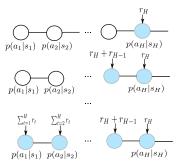
* past rewards do not depend on the current action

$$\sim \frac{1}{m} \sum_{i=1}^{m} \sum_{t=1}^{H} \nabla_{\boldsymbol{\theta}} \log \pi_{\boldsymbol{\theta}}(\mathbf{a}_{t}^{(i)}|\mathbf{s}_{t}^{(i)}) [\sum_{k=t}^{H} r(\mathbf{s}_{k}^{(i)}, \mathbf{a}_{k}^{(i)})]$$

https://www.youtube.com/watch?v=S_gwYj1Q-44 (28')



Algorithm 2



- ► Same as Algorithm 1
- ▶ But the sum is incomplete, and computed backwards
- ▶ Slightly less variance, because it ignores irrelevant rewards



Discounting rewards

$$\nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}) \sim \frac{1}{m} \sum_{i=1}^{m} \sum_{t=1}^{H} \nabla_{\boldsymbol{\theta}} \log \pi_{\boldsymbol{\theta}}(\mathbf{a}_{t}^{(i)} | \mathbf{s}_{t}^{(i)}) [\sum_{k=t}^{H} r(\mathbf{s}_{k}^{(i)}, \mathbf{a}_{k}^{(i)})]$$
* reduce the variance by discounting the rewards along the trajectory
$$1 \sum_{k=t}^{m} \sum_{t=1}^{H} \nabla_{\boldsymbol{\theta}} \mathbf{a}_{t}^{(i)} \mathbf{a}_{t}^{(i)} \mathbf{b}_{t}^{(i)} \sum_{k=t}^{H} c_{t}^{(i)} \mathbf{a}_{t}^{(i)} \mathbf{b}_{t}^{(i)}$$

$$\sim \frac{1}{m} \sum_{i=1}^{m} \sum_{t=1}^{H} \nabla_{\boldsymbol{\theta}} \log \pi_{\boldsymbol{\theta}}(\mathbf{a}_{t}^{(i)}|\mathbf{s}_{t}^{(i)}) [\sum_{k=t}^{H} \boldsymbol{\gamma}^{k-t} r(\mathbf{s}_{k}^{(i)}, \mathbf{a}_{k}^{(i)})]$$

$$\downarrow_{p(a_{1}|s_{1}) \ p(a_{2}|s_{2})} \cdots \underbrace{\downarrow_{p(a_{H}|s_{H})}}_{r_{H-1}+\gamma r_{H}} \underbrace{\downarrow_{r_{H}}}_{r_{H}}$$

$$\vdots \cdots \underbrace{\downarrow_{p(a_{H}|s_{H})}}_{p(a_{1}|s_{1}) \ p(a_{2}|s_{2})} \cdots \underbrace{\downarrow_{p(a_{H}|s_{H})}}_{r_{H-1}+\gamma r_{H}} \underbrace{\downarrow_{r_{H}}}_{r_{H}}$$



Introducing the action-value function

- lacksquare $\sum_{k=t}^{H} \gamma^{k-t} r(\mathbf{s}_k^{(i)}, \mathbf{a}_k^{(i)})$ can be rewritten $Q_{(i)}^{\pi_{m{ heta}}}(\mathbf{s}_t^{(i)}, \mathbf{a}_t^{(i)})$
 - $\nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}) \sim \frac{1}{m} \sum_{i=1}^{m} \sum_{t=1}^{H} \nabla_{\boldsymbol{\theta}} \mathsf{log} \pi_{\boldsymbol{\theta}}(\mathbf{a}_{t}^{(i)} | \mathbf{s}_{t}^{(i)}) Q_{(i)}^{\pi_{\boldsymbol{\theta}}}(\mathbf{s}_{t}^{(i)}, \mathbf{a}_{t}^{(i)})$
- It is just rewriting, not a new algorithm
- ▶ But suggests that the gradient could be just a function of the local step if we could estimate $Q_{(i)}^{\pi\theta}(\mathbf{s}_t, \mathbf{a}_t)$ in one step

Estimating $Q^{\pi_{\theta}}(s, a)$

- Instead of estimating $Q^{\pi_{\theta}}(s,a) = \mathbb{E}_{(i)}[Q^{\pi_{\theta}}_{(i)}(s,a)]$ from Monte Carlo,
- **b** Build a model $\hat{Q}_{\phi}^{\pi_{\theta}}$ of $Q^{\pi_{\theta}}$ through function approximation
- Two approaches:
 - Monte Carlo estimate: Regression against empirical return

$$\phi_{j+1} \to arg \min_{\phi_j} \frac{1}{m} \sum_{i=1}^{m} \sum_{t=1}^{H} (\sum_{k=t}^{H} \gamma^{k-t} r(\mathbf{s}_k^{(i)}, \mathbf{a}_k^{(i)}) - \hat{Q}_{\phi_j}^{\pi_{\theta}} (\mathbf{s}_t^{(i)}, \mathbf{a}_t^{(i)}))^2$$

 \blacktriangleright Temporal Difference estimate: init $\hat{Q}^{\pi_{\pmb{\theta}}}_{\pmb{\phi}_0}$ and fit using $(\mathbf{s},\mathbf{a},r,\mathbf{s}')$ data

$$\phi_{j+1} \rightarrow \min_{\phi_j} \sum_{(\mathbf{s}, \mathbf{a}, r, \mathbf{s}')} ||r + \gamma f(\hat{Q}^{\pi_{\theta}}_{\phi_j}(\mathbf{s}', .)) - \hat{Q}^{\pi_{\theta}}_{\phi_j}(\mathbf{s}, \mathbf{a})||^2$$

- $\qquad \qquad f(\hat{Q}_{\phi_j}^{\pi_{\theta}}(\mathbf{s}',.)) = \max_{\mathbf{a}} \hat{Q}_{\phi_j}^{\pi_{\theta}}(\mathbf{s}',\mathbf{a}) \text{ (Q-learning)}, \\ = \hat{Q}_{\phi_j}^{\pi_{\theta}}(\mathbf{s}',\pi_{\theta}(\mathbf{s}')) \text{ (AC)}...$
- May need some regularization to prevent large steps in φ https://www.youtube.com/watch?v=S_gwYj1Q-44 (36')

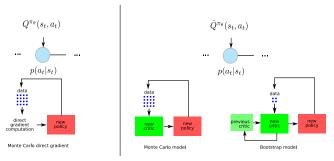


Martin Riedmiller. Neural fitted Q iteration-first experiences with a data efficient neural reinforcement learning method. In European Conference on Machine Learning, pp. 317–328. Springer, 2005



András Antos, Csaba Szepesvári, and Rémi Munos. Fitted Q-iteration in continuous action-space MDPs. In Advances in neural information processing systems, pp.9–16, 2008.

Monte Carlo versus Bootstrap approaches



- Three options:
 - MC direct gradient: Compute the true $Q^{\pi_{\theta}}$ over each trajectory
 - MC model: Compute a model $\hat{Q}_{\phi}^{\pi\theta}$ over rollouts using MC regression, throw it away after each policy gradient step
 - ▶ Bootstrap: Update a model $\hat{Q}^{\pi\theta}_{\phi}$ over samples using TD methods, keep it over policy gradient steps
- ▶ With bootstrap, update everything from the current state, see next lessons
- Next lesson: adding a baseline

Any question?



Send mail to: Olivier.Sigaud@upmc.fr





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