

Fraunhofer-Institut für Integrierte Schaltungen IIS

Reinforcement Learning

Lecture 7: Policy-based RL 2

Christopher Mutschler

But how do we maximize this?

→ Gradient Ascent! Suppose we know how to calculate the gradient w.r.t. the parameters:

 $\nabla_{\theta} \mathbb{E}_{\tau \sim \pi_{\theta}} G(\tau)$

• Then we can update our parameters θ with a learning rate α in the direction of the gradient:

θ

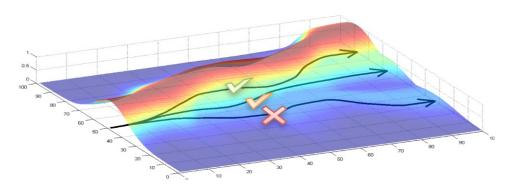
$$\leftarrow \theta + \alpha \nabla_{\theta} \mathbb{E}_{\tau \sim \pi_{\theta}} G(\tau)$$
Policy Gradient
often in literature referred to as $\nabla_{\theta} J(\pi_{\theta})$



Plugging grad-log-prob into the gradient update gives:

$$\nabla_{\theta} \mathbb{E}_{\tau \sim \pi_{\theta}} G(\tau) = \mathbb{E}_{\tau \sim \pi_{\theta}} (\nabla_{\theta} \log P(\tau|\theta) G(\tau))$$
$$= \mathbb{E}_{\tau \sim \pi_{\theta}} \left(\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) G(\tau) \right)$$

- But what is the intuition behind this gradient?
- The gradient tries to
 - Increase probability of paths with positive G
 - Decrease probability of paths with negative G
 In practice:
 - Increase probability of paths with small positive G a bit
 - Increase probability of paths with large positive G much more



Pieter Abbeel. DeepRL Bootcamp 4A Policy Gradients.



Plugging grad-log-prob into the gradient update gives:

$$\nabla_{\theta} \mathbb{E}_{\tau \sim \pi_{\theta}} G(\tau) = \mathbb{E}_{\tau \sim \pi_{\theta}} (\nabla_{\theta} \log P(\tau|\theta) G(\tau))$$
$$= \mathbb{E}_{\tau \sim \pi_{\theta}} \left(\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) G(\tau) \right)$$

Reduce variance: as this is an expectation, we can estimate it with a sample mean using Monte-Carlo sampling of *L* trajectories:
 each action in the episode in influenced by the reward of the whole episode???

$$\nabla_{\theta} \mathbb{E}_{\tau \sim \pi_{\theta}} G(\tau) \approx \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) G(\tau)$$
$$\approx \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \sum_{t'=t}^{T} \gamma^{t'-t} R(s_{t'}, a_{t'})$$



- Monte-Carle Policy Gradient Control
 - Update parameters by stochastic gradient ascent
 - Using policy gradient theorem
 - Using return-to-go $Q^{\pi_{\theta}}(s_t, a_t)$ as an unbiased sample of G:

 $\Delta \theta_t = \alpha \nabla_\theta \log \pi_\theta(a_t | s_t) \, \gamma G_t$

REINFORCE: Monte-Carlo Policy-Gradient Control (episodic) for π_*

```
Input: a differentiable policy parameterization \pi(a|s, \theta)

Algorithm parameter: step size \alpha > 0

Initialize policy parameter \theta \in \mathbb{R}^{d'} (e.g., to 0)

Loop forever (for each episode):

Generate an episode S_0, A_0, R_1, \dots, S_{T-1}, A_{T-1}, R_T, following \pi(\cdot|\cdot, \theta)

Loop for each step of the episode t = 0, 1, \dots, T - 1:

G \leftarrow \sum_{k=t+1}^{T} \gamma^{k-t-1} R_k

\theta \leftarrow \theta + \alpha \gamma^t G \nabla \ln \pi(A_t|S_t, \theta)
(G<sub>t</sub>)
```

Sutton, R. S., & Barto, A. G. (2018). Reinforcement learning: An introduction. MIT press.



Policy-based RL

Agenda

- Policy-based RL 1
 - Intro to Policy-based RL
 - Policy Gradients
- Policy-based RL 2
 - Variance & Baselines
 - Actor-Critics
 - Trust-Region Policy Optimization (TRPO)
 - Proximal Policy Optimization (PPO)
 - Deep Deterministic Policy Gradient (DDPG)



Policy Gradients: where to hide the variance?

- Expected Grad-Log-Prob Lemma*:
 - P_{θ} is a parameterized probability distribution over a random variable x, then:

$$\mathbb{E}_{x \sim P_{\theta}}[\nabla_{\theta} \log P_{\theta}(x)] = 0$$

• From this follows that for any function *b* that *only depends on states*:

 \mathbb{E}_a

$$\sum_{t \sim \pi_{\theta}} [\nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t})b(s_{t})] = 0,$$
 because:

$$= \mathbb{E} \left[\sum_{a} \pi(a|s_{t})b\nabla_{\theta} \log \pi(a|s_{t}) \right]$$

$$= \mathbb{E} \left[b\nabla_{\theta} \sum_{a} \pi(a|s_{t}) \right]$$

$$= \mathbb{E} [b\nabla_{\theta} 1]$$

$$= 0$$
 inverse score-function trick prob-distribution and we sum over all actions

* You can find one version of the proof here: https://spinningup.openai.com/en/latest/spinningup/rl_intro3.html#id1

DIGRESSION



Policy Gradients: where to hide the variance?

 This allows us to add or subtract any of such terms (i.e., baseline functions) to our policy gradient without changing its expectation:

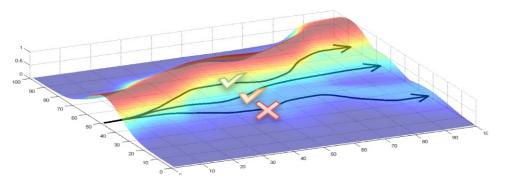
$$\nabla_{\theta} \mathbb{E}_{\tau \sim \pi_{\theta}} G(\tau) = \mathbb{E}_{\tau \sim \pi_{\theta}} \left(\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \left(G_t - b(s_t) \right) \right)$$

* You can find one version of the proof here: https://spinningup.openai.com/en/latest/spinningup/rl_intro3.html#id1



Advantage Functions: Intuition

- But what is the intuition behind this gradient?
- The gradient tries to
 - Increase probability of paths with positive G
 - Decrease probability of paths with negative G



Pieter Abbeel. DeepRL Bootcamp 4A Policy Gradients.

- You mostly keep on increasing everything (but some more than others)
 - And this requires a lot of rollouts to average the effect out!
 - Ideal: in-/decrease probs of paths that are better/worse than the average!



Advantage Functions: Introduce a Baseline

But how does this help?

- We can introduce a baseline to our targets!
- We can define *advantage functions* that reduce variance
- Intuition: if an agent sees what it expected, it feels *neutral* about it
- *Hint: we already did this when we introduced the reward-to-go variant!*
- For instance, we can use $b(s_t) = v^{\pi}(s_t)$ to reduce variance in the sample estimate of the policy gradient:

 $A^{\pi}(s_t, a_t) = Q^{\pi}(s_t, a_t) - V^{\pi}(s_t)$

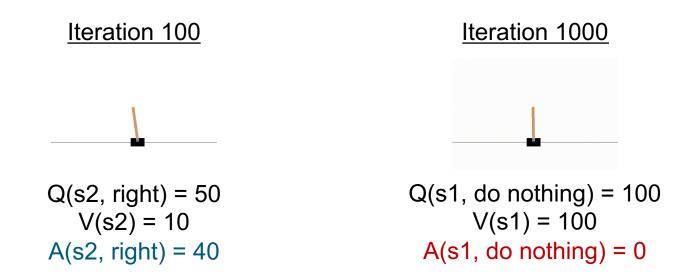
→ faster and more stable policy learning!



Advantage Functions: Intuition

Question: what helps our agent learn faster: information about

- 1. invariant actions on good states, or
- 2. information about good actions in bad/challenging states?





REINFORCE w/ Baselines

REINFORCE with Baseline (episodic), for estimating $\pi_{\theta} \approx \pi_*$

Input: a differentiable policy parameterization $\pi(a|s, \theta)$ Input: a differentiable state-value function parameterization $\hat{v}(s, \mathbf{w})$ Algorithm parameters: step sizes $\alpha^{\theta} > 0$, $\alpha^{\mathbf{w}} > 0$ Initialize policy parameter $\theta \in \mathbb{R}^{d'}$ and state-value weights $\mathbf{w} \in \mathbb{R}^{d}$ (e.g., to $\mathbf{0}$) Loop forever (for each episode): Generate an episode $S_0, A_0, R_1, \dots, S_{T-1}, A_{T-1}, R_T$, following $\pi(\cdot|\cdot, \theta)$ Loop for each step of the episode $t = 0, 1, \dots, T - 1$: $G \leftarrow \sum_{k=t+1}^{T} \gamma^{k-t-1} R_k$ (G_t) $\delta \leftarrow G - \hat{v}(S_t, \mathbf{w})$ \leftarrow advantage $\mathbf{w} \leftarrow \mathbf{w} + \alpha^{\mathbf{w}} \delta \nabla \hat{v}(S_t, \mathbf{w})$ $\theta \leftarrow \theta + \alpha^{\theta} \gamma^t \delta \nabla \ln \pi(A_t|S_t, \theta)$ \leftarrow policy gradient update

Sutton, R. S., & Barto, A. G. (2018). Reinforcement learning: An introduction. MIT press.



REINFORCE w/ Baselines

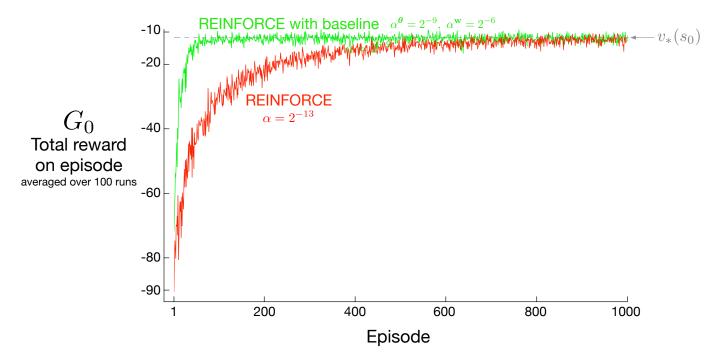


Figure 13.2: Adding a baseline to REINFORCE can make it learn much faster, as illustrated here on the short-corridor gridworld (Example 13.1). The step size used here for plain REINFORCE is that at which it performs best (to the nearest power of two; see Figure 13.1).

Sutton, R. S., & Barto, A. G. (2018). Reinforcement learning: An introduction. MIT press.



Vanilla Policy Gradient Algorithm

Pseudocode

Algorithm 1 Vanilla Policy Gradient Algorithm

1: Input: initial policy parameters θ_0 , initial value function parameters ϕ_0

2: for $k = 0, 1, 2, \dots$ do

- 3: Collect set of trajectories $\mathcal{D}_k = \{\tau_i\}$ by running policy $\pi_k = \pi(\theta_k)$ in the environment.
- 4: Compute rewards-to-go \hat{R}_t .
- 5: Compute advantage estimates, \hat{A}_t (using any method of advantage estimation) based on the current value function V_{ϕ_k} .
- 6: Estimate policy gradient as

$$\hat{g}_k = \frac{1}{|\mathcal{D}_k|} \sum_{\tau \in \mathcal{D}_k} \sum_{t=0}^T \nabla_\theta \log \pi_\theta(a_t|s_t)|_{\theta_k} \hat{A}_t.$$

7: Compute policy update, either using standard gradient ascent,

 $\theta_{k+1} = \theta_k + \alpha_k \hat{g}_k$

or via another gradient ascent algorithm like Adam

8: Fit value function by regression on mean-squared error:

$$\phi_{k+1} = \arg\min_{\phi} \frac{1}{|\mathcal{D}_k|T} \sum_{\tau \in \mathcal{D}_k} \sum_{t=0}^T \left(V_{\phi}(s_t) - \hat{R}_t \right)^2,$$

typically via some gradient descent algorithm.

9: end for

Problem:

- Learning rate is a heuristic that cannot be easily tuned
- If we take large steps (especially in the beginning of training) we might never recover

Estimate the Advantages

Calculate the gradient and take a gradient step

Update the Critic (used for the Advantage estimation)



Policy-based RL

Agenda

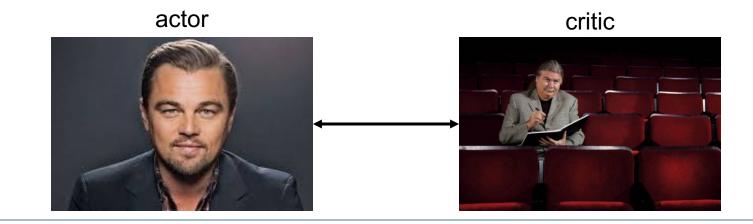
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Policy Gradients: Variance (revisited)

$$\nabla_{\theta} \mathbb{E}_{\pi_{\theta}} G(\tau) \approx \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \sum_{t'=t}^{T} \gamma^{t'-t} R(s_{t'}, a_{t'}) - b(s_t)$$
$$= Q^{\pi}(s_t, a_t)$$

- Monte-Carlo policy gradient is sampled and has high variance
- Idea: we can use a critic that estimates the Q





Policy Gradients: Variance (revisited)

$$\nabla_{\theta} \mathbb{E}_{\pi_{\theta}} G(\tau) \approx \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \sum_{\substack{t'=t \\ t'=t}}^{T} \gamma^{t'-t} R(s_{t'}, a_{t'}) - b(s_t)$$
$$= Q^{\pi}(s_t, a_t)$$

- Monte-Carlo policy gradient is sampled and has high variance
- Idea: we can use a critic that estimates the Q
- → In practice, (as we already know) $Q^{\pi}(s_t, a_t)$ cannot be "computed"
- \rightarrow we instead need to approximate it with a neural network with parameters ϕ and standard SGD over k epochs minimizing the MSE:

$$\phi_k = \arg\min_{\phi} \mathbb{E}_{s_t; a_t, \hat{R}_t \sim \pi_k} \left[\left(Q_{\phi}(s_t, a_t) - \hat{R}_t \right)^2 \right]$$



Actor-Critics Actor-Critic

$$\nabla_{\theta} \mathbb{E}_{\pi_{\theta}} G(\tau) \approx \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \sum_{t'=t}^{T} \gamma^{t'-t} R(s_{t'}, a_{t'}) - b(s_t) = Q^{\pi}(s_t, a_t)$$

- Monte-Carlo policy gradient is sampled and has high variance
- Idea: we can use a critic (e.g., a NN as in DQN) to estimate the Q and learn two sets of parameters separately
 - Actor: update θ by policy gradient ?
 - Critic: Update parameters ϕ of v_{ϕ}^{π} , e.g., by n-step TD
- We call such algorithms *actor-critic algorithms*



Actor-Critic

• Typically, we estimate $v^{\pi}(s_t; \phi)$ explicitly, and then sample

 $q^{\pi}(s_t, a_t) \approx G_t^{(n)}$

- For instance: $\hat{G}_t^{(1)} = R_t + \gamma v^{\pi}(s_{t+1}; \phi)$
- Then we arrive at:

$$\nabla_{\theta} \mathbb{E}_{\pi_{\theta}} \widehat{G}(\tau) = \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \left(\widehat{G}_t - b(s_t) \right)$$

 \rightarrow We use an *approximate* gradient in the direction suggested by the critic



Actor-Critics Actor-Critic: Advantage Functions

$$\nabla_{\theta} \mathbb{E}_{\pi_{\theta}} G(\tau) \approx \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \left(\hat{G}_t - b(s_t) \right)$$
$$:= A^{\pi}(s_t, a_t)$$

Calculating the advantage function is (embarrassingly) straight forward:

• As before, we can simply use the TD error:

$$A^{\pi}(s_t, a_t) = Q^{\pi}(s_t, a_t) - V^{\pi}(s_t) = r + \gamma \cdot v^{\pi}(s_{t+1}) - v^{\pi}(s_t)$$



Actor-Critics Actor-Critic: Advantage Functions

$$\nabla_{\theta} \mathbb{E}_{\pi_{\theta}} G(\tau) \approx \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \left(\hat{G}_t - b(s_t) \right)$$
$$:= A^{\pi}(s_t, a_t)$$

Calculating the advantage function is (embarrassingly) straight forward:

 We can also use Generalized Advantage Estimation (GAE) (multi-step TD-error like TD with n-step returns)

$$\begin{split} \hat{A}_{t}^{(1)} &\coloneqq \delta_{t}^{V} &= -V(s_{t}) + r_{t} + \gamma V(s_{t+1}) \\ \hat{A}_{t}^{(2)} &\coloneqq \delta_{t}^{V} + \gamma \delta_{t+1}^{V} &= -V(s_{t}) + r_{t} + \gamma V(s_{t+1}) + \gamma^{2} V(s_{t+2}) \\ \hat{A}_{t}^{(3)} &\coloneqq \delta_{t}^{V} + \gamma \delta_{t+1}^{V} + \gamma^{2} \delta_{t+2}^{V} &= -V(s_{t}) + r_{t} + \gamma V(s_{t+1}) + \gamma^{2} V(s_{t+2}) + \gamma^{3} V(s_{t+3}) \\ \hat{A}_{t}^{(k)} &\coloneqq \sum_{l=0}^{k-1} \gamma^{l} \delta_{t+l}^{V} &= -V(s_{t}) + r_{t} + \gamma r_{t+1} + \dots + \gamma^{k-1} r_{t+k-1} + \gamma^{k} V(s_{t+k}) \\ \hat{A}_{t}^{(\infty)} &= \sum_{l=0}^{\infty} \gamma^{l} \delta_{t+l}^{V} = -V(s_{t}) + \sum_{l=0}^{\infty} \gamma^{l} + r_{t+l} \end{split}$$



Actor-Critics Actor-Critic: Advantage Functions

$$\nabla_{\theta} \mathbb{E}_{\pi_{\theta}} G(\tau) \approx \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \left(\widehat{G}_t - b(s_t) \right)$$
$$:= A^{\pi}(s_t, a_t)$$

Calculating the advantage function is (embarrassingly) straight forward:

- Full returns: high variance
- One-step TD: high bias
- n-step TD: somewhere in the middle

But still: approximating the policy gradient introduces bias

- It is important to use on-policy targets (i.e., can be corrected using importance sampling)
- Alternative idea: bootstrap (with $\lambda = 0$) whenever policies differ



Continuous Actions: Gaussian Policy

- Because we directly update the policy parameters of the policy, we can easily deal with continuous action spaces
 - Most algorithms discussed can be used for both discrete and continuous actions
- In continuous action spaces, a Gaussian policy is common, e.g., mean is some function of state $\mu(s)$
- For simplicity, lets consider fixed variance of σ^2 (which can be parameterized as well, instead)
- Policy is Gaussian: $a \sim \mathcal{N}(\mu(s), \sigma^2)$
- The gradient of the log of the policy is then

$$\nabla_{\theta} \log \pi_{\theta}(s, a) = \frac{a - \mu(s)}{\sigma^2} \nabla \mu(s)$$

This can be used, for instance, in REINFORCE or advantage actor-critic



Continuous Actions: Exploration-Exploitation

• Exploration vs. Exploitation of Policy Gradient methods:

- PG trains a stochastic policy in an on-policy way
- Actions are sampled from the environment according to the latest version of its stochastic policy
- Randomness in selecting actions
 - (initially) depends on the initialization
 - Becomes less over the course of training
 - ...as the update rule encourages to exploit rewards that the policy already has found
- Policy-gradients only consider improvement under current data
 → easy to get stuck in local optima



One more thing: Exploration-Exploitation

- Could we use ϵ -greedy? Yes! but it is not ideal...
 - Wildly different actions cause breakage
 - Exploration is mostly uninformed about current best guess
- Alternative idea: make sure that the entropy of the policy is not too low:

$$-\sum_{s} \mu(s) \sum_{a} \pi(a|s) \log \pi(a|s) = -\mathbb{E}[\log \pi(a_t|s_t)]$$

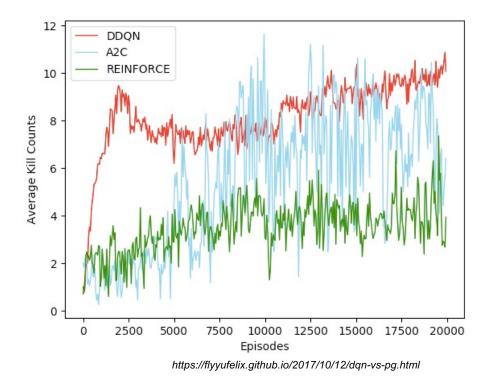
- ightarrow Add a regularization term that pushes entropy up slightly each step
- Encourages exploration and does not pick fully randomly
 - May increase variance in Gaussian policies
 - Makes softmax slightly more uniform
- Similar to (spoiler!) KL-regularization (in TRPO)
- Works well in practice



Conclusion

- It is substantially different from DQNs
 - no replay buffer, no stored experiences
 - learn directly, on-policy

- Once a batch has been used \rightarrow discard experience
 - less sample efficient



• Learning off-policy (which would allow to reuse experience) is not (or only with certain tricks) possible



Conclusion

• The policy gradient has many forms:

 $\begin{aligned} \nabla_{\theta} J(\theta) &= \mathbb{E}_{\pi_{\theta}} [\nabla_{\theta} \log \pi_{\theta}(s, a) G_{t}] \\ \nabla_{\theta} J(\theta) &= \mathbb{E}_{\pi_{\theta}} [\nabla_{\theta} \log \pi_{\theta}(s, a) (G_{t} - b(s_{t})] \\ \nabla_{\theta} J(\theta) &= \mathbb{E}_{\pi_{\theta}} \left[\nabla_{\theta} \log \pi_{\theta}(s, a) \hat{A}_{t}^{(\infty)} \right] \\ \nabla_{\theta} J(\theta) &= \nabla_{\theta} Q_{t}(s, \pi_{\theta}(s)) \end{aligned}$

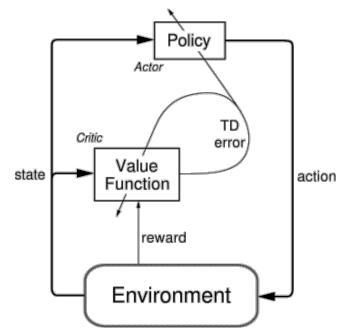
REINFORCE

REINFORCE w/ baseline

advantage actor-critic

deterministic policy gradient (see DDPG)

- Each leads to a stochastic gradient ascent algorithm
- Critic uses policy evaluation (e.g., MC or TD) to estimate $Q^{\pi}(s, a)$ or $V^{\pi}(s)$

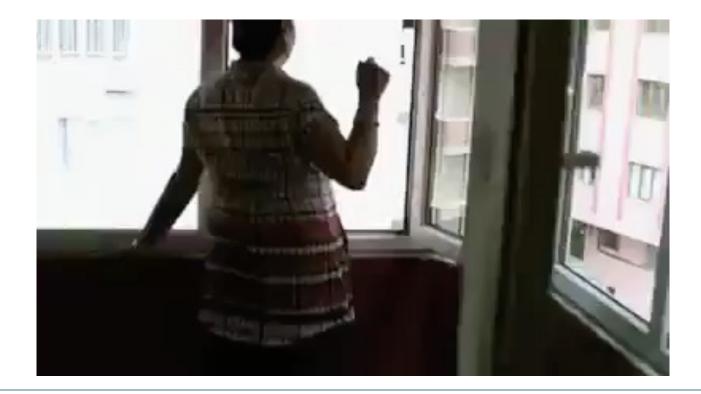


Sutton et a.: Reinforcement learning: An introduction. 2018.





"Always try to solve a problem the direct way – but be careful with the high variance."



https://www.youtube.com/watch?v=ek3hgVQ1RqM



Policy-based RL

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Policy Gradients so far

• We wanted to do the following:

$$\max_{\theta} J(\pi_{\theta}) = \mathbb{E}_{\tau \sim \pi_{\theta}} \left(\sum_{t=0}^{\infty} \gamma^{t} r_{t} \right)$$

• Then we defined the policy gradient:

$$g = \nabla_{\theta} J(\pi_{\theta}) = \nabla_{\theta} \mathbb{E}_{\pi_{\theta}} G(\tau) \approx \frac{1}{L} \sum_{\tau} \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \hat{A}_t$$

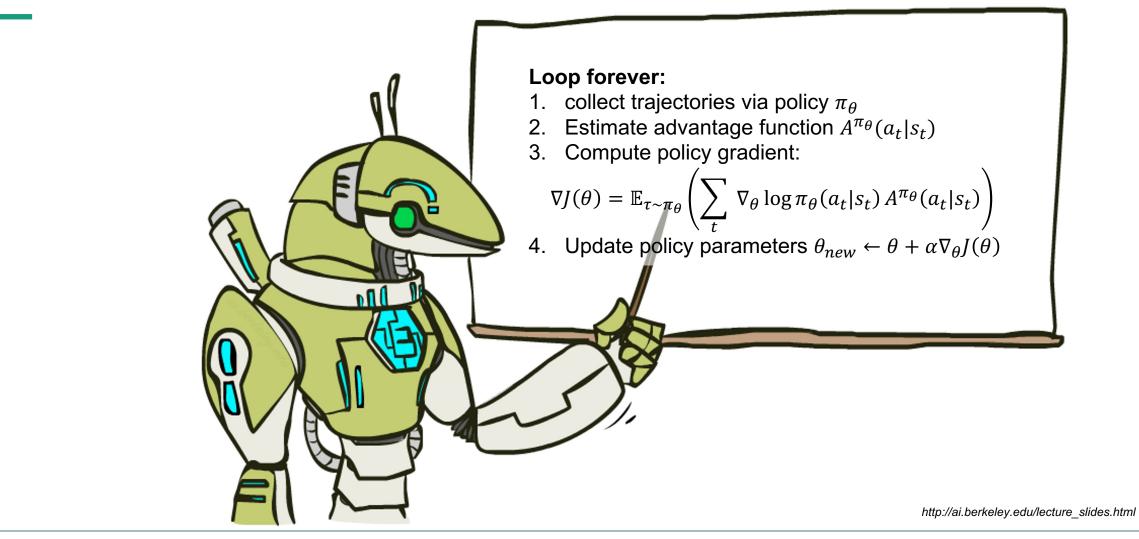
- If $\hat{A}_t > 0 \ (< 0)$:
 - Gradient becomes positive (negative)
 - The probability of taking a_t in s_t is increased (decreased)
- We update our policy parameters by

$$\theta_{k+1} = \theta_k + \alpha \cdot g$$

Take a gradient step in updating the policy (gradient ascent)



Policy Gradients so far





Policy Gradients so far

Problem

- Run gradient descent/ascent on one batch of collected experience
- Note: the advantage function (which is a noisy estimate) may not be accurate
 - Too large steps may lead to a disaster (even *if* the gradient is *correct*)
 - Too small steps are also bad
- Definition and scheduling of learning rates in RL is tricky as the underlying data distribution changes with updates to the policy
- Mathematical formulization:
 - First-order derivatives approximate the (parameter) surface to be flat
 - But if the surface exhibits high curvature it gets dangerous
 - Projection: small changes in parameter space might lead to large changes in policy space!
- Parameters θ get updated to areas too far out of the range from where previous data was collected (note: a bad policy leads to bad data – unlike in value-based RL!)

Images taken from https://medium.com/@jonathan_hui/rl-trust-region-policy-optimization-trpo-explained-a6ee04eeeee9 and http://www.taiwanoffthebeatentrack.com/2012/08/23/mount-hua-华山-the-most-dangerous-hike-in-the-world/



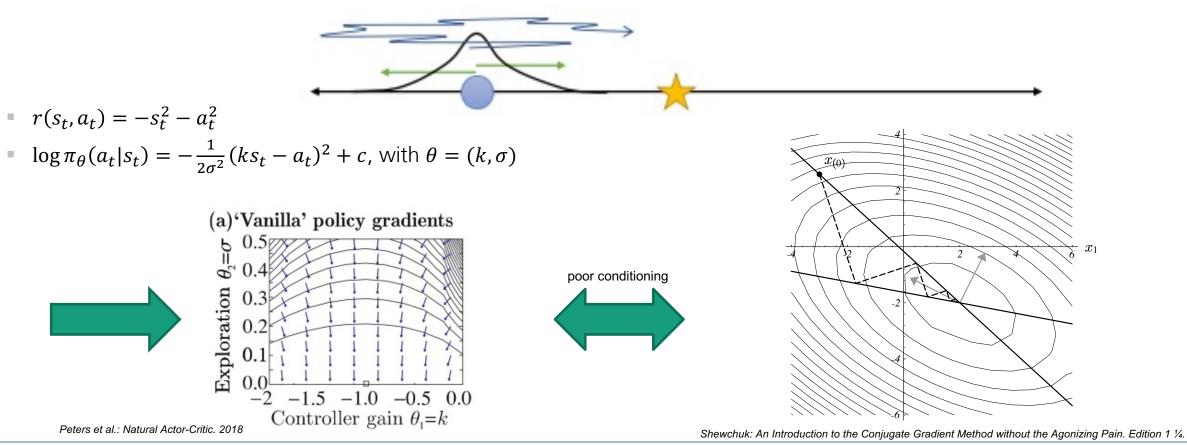






Primer: Natural Policy Gradient

There is something wrong with the policy gradient. Example:



* Example take from Sergey Levine's CS285: Advanced Policy Gradients



Primer: Natural Policy Gradient

• Given:

 $\theta \leftarrow \theta + \alpha \nabla_{\theta} J(\theta) \quad \text{and} \quad \pi_{\theta}(a_t | s_t)$

- Some parameters change probabilities a lot more than others! \rightarrow a single value for α is not sufficient
- What we essentially do (optimization perspective on 1st order gradient descent):

e θ' ← arg max_θ(θ' − θ)^T∇_θJ(θ), subject to
$$||θ' − θ||^2 ≤ ε$$
θ' ← arg max_θ(θ' − θ)^T∇_θJ(θ), subject to $D(π_{θ'}, π_{θ}) ≤ ε$
"trust region"
(b) Natural policy gradients
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"Simple" Idea

Regularize updates to the policy parameters, such that the policy does not change too much.



only for reference

Primer: Trust-Region Methods

Optimization in Machine Learning

- It is common to formulate ML problems as optimization problems:
 - Minimize the squared error
 - Minimize the cross entropy
 - Maximize the log-likelihood
 - Maximize the discounted sum of rewards
 - Minimize the sum of control cost



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Primer: Trust-Region Methods

Optimization in Machine Learning: two classes

- 1. Line Search, e.g., gradient descent
 - find a (some) direction of improvement
 - (cleverly) select a step length
- 2. Trust-Region Methods
 - select a trust region (analog to max step length)
 - find a point of improvement in that region







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Primer: Trust-Region Methods

- Idea:
 - Approximate the real objective f with something simpler, i.e., \tilde{f}
 - Solve $\tilde{x}^* = \arg \min_x \tilde{f}(x)$
- Problem:
 - The optimum \tilde{x}^* might be in a region where \tilde{f} poorly approximates f
 - \tilde{x}^* might be far from optimal
- Solution:
 - Restrict the search to a region tr where we trust \tilde{f} to approximate f well

• Solve $\tilde{x}^* = \arg\min_{x \in tr} \tilde{f}(x)$



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Primer: Trust-Region Methods

2nd order Taylor approximation

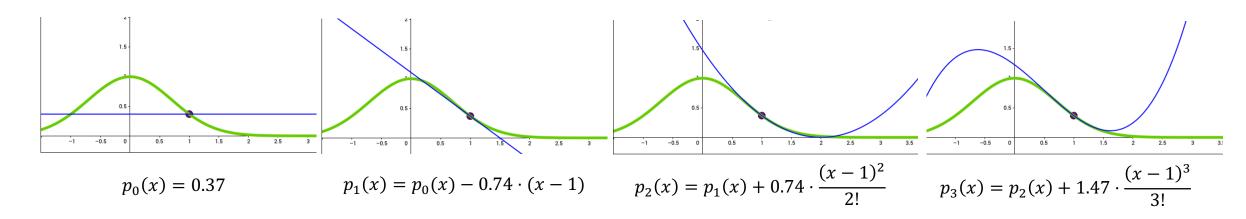
• Example: chose \tilde{f} to be quadratic approximation of f:

$$f(x) \approx \tilde{f}(x) = f(c) + \nabla f(c)^T (x - c) + \frac{1}{2!} (x - c)^T H(c) (x - c),$$

where ∇f is the gradient and H is the Hessian

Example: nth order Taylor approximation of $f(x) = e^{-x^2}$, c = 1

$$\Rightarrow p_n(x) = \sum_{j=0}^n \frac{1}{j!} f^{(j)}(a) (x-a)^j$$



https://mathinsight.org/applet/taylor_polynomial



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Primer: Trust-Region Methods

2nd order Taylor approximation

• Example: chose \tilde{f} to be quadratic approximation of f:

$$f(x) \approx \tilde{f}(x) = f(c) + \nabla f(c)^T (x - c) + \frac{1}{2!} (x - c)^T H(c) (x - c),$$

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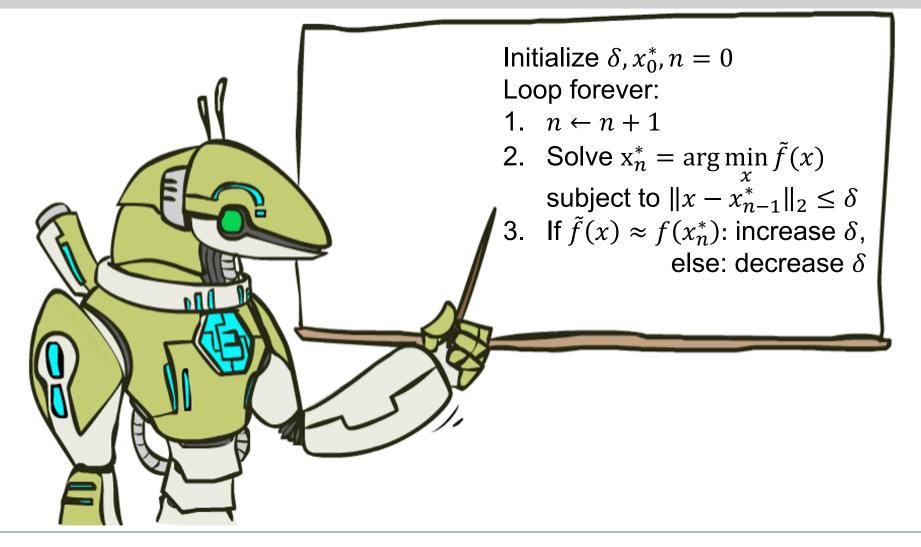
• The trust region is often chosen to be a hypersphere:

$$\|x - c\|_2 \le \delta$$



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Primer: Trust-Region Methods





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Primer: Trust-Region Methods

2nd order Taylor approximation

• \tilde{f} often chosen to be quadratic approximation of f:

$$\min_{x} f(c) + \nabla f(c)^{T} (x-c) + \frac{1}{2!} (x-c)^{T} H(c) (x-c),$$

subject to $||x - c||_2 \le \delta$.

- When *H* is positive semi-definite
 - Convex optimization
 - Simple and globally optimal solution
 - e.g.: conjugate gradient
- When *H* is not positive semi-definite
 - Non-convex optimization
 - Simple heuristics that guarantee improvement



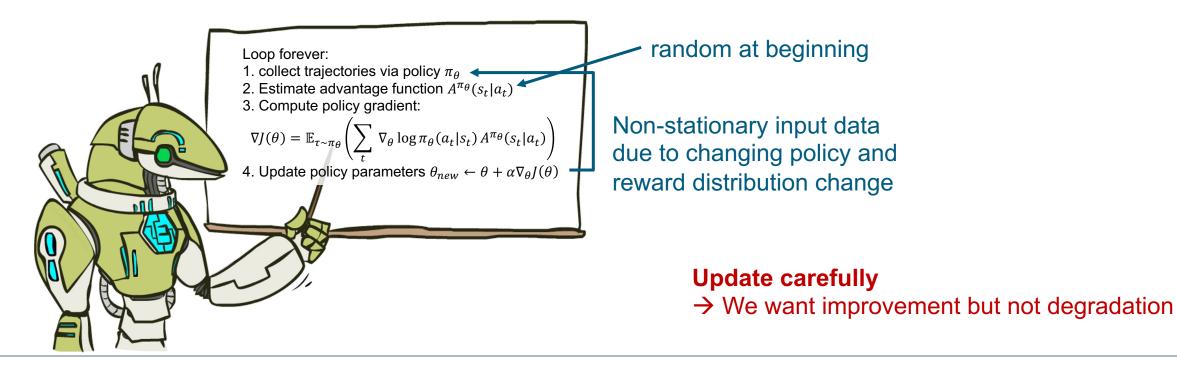


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The problem(s) of the Policy Gradient (PG) is that

Back to RL...

- PG keeps old and new policy close in parameter space (not in policy space!), while
- small changes can lead to large differences in performance, and
- "large" step-sizes hurt performance (whatever "large" means...)





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• We want to optimize $\eta(\pi)$, i.e., the expected return of policy π :

$$\eta(\pi) = \mathbb{E}_{s_0 \sim \rho_0, a^t \sim \pi_{old}(\cdot \mid s_t)} \left[\sum_{t=0}^{\infty} \gamma^t r_t \right]$$

- We collect data with π_{old} and optimize to get a new policy π_{new}
- In fact, so far...
 - we did not really optimize anything, as
 - we just favored those $(a_t|s_t)$ that had a higher advantage
- Question:
 - How can we write down an optimization problem that allows to do small updates on a policy π based on data sampled from π (on-policy data)?

¹ Kakade et al.: Approximately Optimal Approximate Reinforcement Learning. ICML 2002.

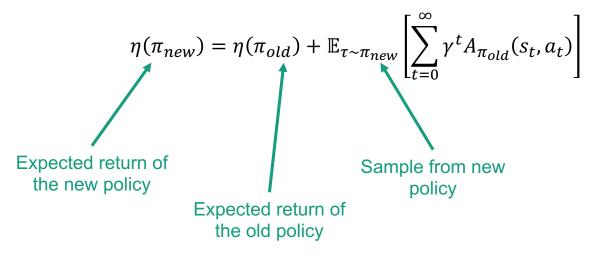


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• We want to optimize $\eta(\pi)$, i.e., the expected return of policy π :

$$\eta(\pi) = \mathbb{E}_{s_0 \sim \rho_0, a^t \sim \pi_{old}(\cdot | s_t)} \left[\sum_{t=0}^{\infty} \gamma^t r_t \right]$$

- We collect data with π_{old} and optimize to get a new policy π_{new}
- Let's express $\eta(\pi_{new})$ in terms of advantage over the original policy¹:



¹ Kakade et al.: Approximately Optimal Approximate Reinforcement Learning. ICML 2002.



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• We want to optimize $\eta(\pi)$, i.e., the expected return of policy π :

$$\eta(\pi) = \mathbb{E}_{s_0 \sim \rho_0, a^t \sim \pi_{old}(\cdot | s_t)} \left[\sum_{t=0}^{\infty} \gamma^t r_t \right]$$

- We collect data with π_{old} and optimize to get a new policy π_{new}
- Let's express $\eta(\pi_{new})$ in terms of advantage over the original policy¹:

$$\eta(\pi_{new}) = \eta(\pi_{old}) + \mathbb{E}_{\tau \sim \pi_{new}} \left[\sum_{t=0}^{\infty} \gamma^t A_{\pi_{old}}(s_t, a_t) \right]$$
$$= \eta(\pi_{old}) + \sum_s \rho_{\pi_{new}}(s) \sum_a \pi_{new}(a|s) A_{\pi_{old}}(s, a)$$

<u>Discounted</u> visitation frequency according to **new** policy: $\rho_{\pi_{new}}(s) = P(s_0 = s) + \gamma P(s_1 = s) + \gamma^2 P(s_2 = s) + \cdots$

¹ Schulman et al.: Trust-Region Policy Optimization. ICML 2015.



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• We want to optimize $\eta(\pi)$, i.e., the expected return of policy π :

$$\eta(\pi) = \mathbb{E}_{s_0 \sim \rho_0, a^t \sim \pi_{old}(\cdot | s_t)} \left[\sum_{t=0}^{\infty} \gamma^t r_t \right]$$

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$$\eta(\pi_{new}) = \eta(\pi_{old}) + \mathbb{E}_{\tau \sim \pi_{new}} \left[\sum_{t=0}^{\infty} \gamma^t A_{\pi_{old}}(s_t, a_t) \right]$$

$$= \eta(\pi_{old}) + \sum_{s} \rho_{\pi_{new}}(s) \sum_{a} \pi_{new}(a|s) A_{\pi_{old}}(s, a)$$

$$0$$
If we can guarantee this...

→ New objective guarantees improvement from $\pi_{old} \rightarrow \pi_{new}$



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• We want to optimize $\eta(\pi)$, i.e., the expected return of policy π :

$$\eta(\pi) = \mathbb{E}_{s_0 \sim \rho_0, a^t \sim \pi_{old}(\cdot | s_t)} \left[\sum_{t=0}^{\infty} \gamma^t r_t \right]$$

- We collect data with π_{old} and optimize to get a new policy π_{new}
- Let's express $\eta(\pi_{new})$ in terms of advantage over the original policy:

η

$$(\pi_{new}) = \eta(\pi_{old}) + \mathbb{E}_{\tau \sim \pi_{new}} \left[\sum_{t=0}^{\infty} \gamma^t A_{\pi_{old}}(s_t, a_t) \right]$$
$$= \eta(\pi_{old}) + \left[\sum_{s} \rho_{\pi_{new}}(s) \sum_{a} \pi_{new}(a|s) A_{\pi_{old}}(s, a) \right]$$

However, this cannot be easily estimated. The state visitations that we sampled so far are coming from the old policy!

 \rightarrow we cannot optimize this in the current form!

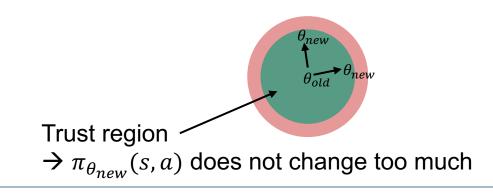


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approximate
locally
$$\eta(\pi_{new}) = \eta(\pi_{old}) + \sum_{s} \rho_{\pi_{new}}(s) \sum_{a} \pi_{new}(a|s) A_{\pi_{old}}(s, a)$$

$$\approx L(\pi_{new}) = \eta(\pi_{old}) + \sum_{s} \rho_{\pi_{old}}(s) \sum_{a} \pi_{new}(a|s) A_{\pi_{old}}(s, a)$$
This we already sampled
 \Rightarrow We already have this!

- The approximation is accurate within step size δ (trust region)
 - δ needs to be chosen based on a lower-bound approximation error
- Monotonic improvement guaranteed
 - (within the green region!)



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- If we want to optimize $L(\theta_{new})$ instead of $\eta(\theta_{new})$... with a guarantee of monotonic improvement on $\eta(\theta_{new})$, we need a bound on $L(\theta_{new})$.
- It can be proven that there exists the following bound^{1,2}:

$$\eta(\pi_{new}) \ge L(\pi_{new}) - C \cdot D_{KL}^{max}(\pi_{old}, \pi_{new})$$
, where $C = \frac{4\epsilon\gamma}{(1-\gamma)^2}$



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Primer: KL-Divergence

• A measure of distance between two probability distributions *P* and *Q*:

$$D_{KL}(P||Q) = \mathbb{E}_{x \sim P}\left[\log \frac{P(x)}{Q(x)}\right] = \mathbb{E}_{x \sim P}\left[\log P(x) - \log Q(x)\right]$$

- Intuition stems from theory of channel coding:
 - The amount of information I (in bits or nats) that is transmitted from a sender to a receiver depends on the actual probability of the actual incident/event

$$I(x) = \log_a \left(\frac{1}{p_x}\right) = \log_a 1 - \log_a p_x = -\log_a p(x)$$

- Example #1: two events A and B have a probability of 0.75 and 0.25; then the information that A has happened is $-log_2(0.75) = 0.41$ (and for B it is 2)
- Example #2: A and B are equally likely, i.e., 0.5 each; then the information about which of them has happened is $-log_2(0.5) = 1$
- If we use base 2 it tells us how much uncertainty is cut off by half!
- For simplification we use nats as $\log_2 x = \log x / \log 2$ (log 2 is constant)



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Primer: KL-Divergence

• The **entropy** (of a probability distribution) is given by:

$$H(p) = -\sum_{i} p_i \log(p_i)$$

- It measures the average amount of information from one sample drawn from the underlying probability distribution $p \rightarrow it$ measures how unpredictable p is
 - ightarrow defines the lower bound of the optimal bit-encoding
- As usual, we can also write H(p) in its expectation and draw x from p:

 $H(p) = \mathbb{E}_{x \sim p}[-\log p(x)]$



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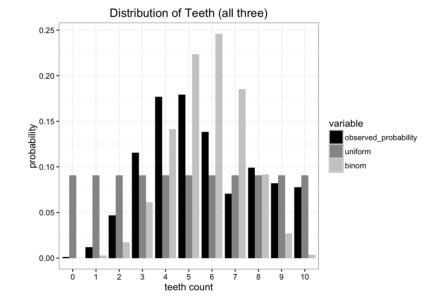
Primer: KL-Divergence

• The **cross-entropy** defines the average number of bits to identify a sampled event if it is encoded using *q* rather than *p*:

$$H(p,q) = -\sum_{i} p_i \log_2(q_i)$$

- tells us the difference about the true probability distribution p and the predicted probability distribution q (given the encoding)
- If p and q are equal, then the cross-entropy and the entropy are equal
- As before, we can write H(p,q) in its expectation:

 $H(p,q) = \mathbb{E}_{x \sim p}[-\log q(x)]$



from https://www.countbayesie.com/blog/2017/5/9/kullback-leibler-divergence-explained



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Primer: KL-Divergence

• **KL-Divergence** tells us how much far away the cross-entropy is from the entropy (note that $H(p,q) \ge H(p)$):

 $D_{KL}(p||q) = H(p,q) - H(p)$

• Measure of dissimilarity between two probability distributions *P* and *Q*:

$$D_{KL}(p||q) = \mathbb{E}_{x \sim p}[-\log q(x)] - \mathbb{E}_{x \sim p}[-\log p(x)]$$

= $\mathbb{E}_{x \sim p}[-\log q(x) - (-\log p(x))]$
= $\mathbb{E}_{x \sim p}[-\log q(x) + \log p(x)]$
= $\mathbb{E}_{x \sim p}[\log p(x) - \log q(x)] = \mathbb{E}_{x \sim p}\left[\log \frac{p(x)}{q(x)}\right]$

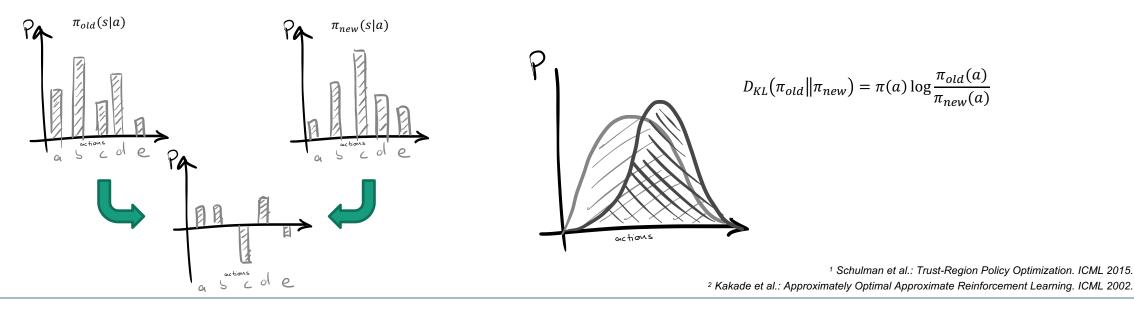
$$D_{KL}(p||q) = \sum_{i} p(i) \log \frac{p(i)}{q(i)} \quad \text{and} \quad D_{KL}(p||q) = \int P(x) \log \frac{p(x)}{q(x)} dx$$

Note: KL-Divergence is asymmetric and hence no distance metric!

Back to RL...

- If we want to optimize L(θ_{new}) instead of η(θ_{new}) ...
 with a guarantee of monotonic improvement on η(θ_{new}), ...
 ... we need a bound on L(θ_{new}).
- It can be proven that the following bound holds^{1,2}:

$$\eta(\pi_{new}) \ge L(\pi_{new}) - C \cdot D_{KL}^{max}(\pi_{old}, \pi_{new})$$
, where $C = \frac{4\varepsilon\gamma}{(1-\gamma)^2}$

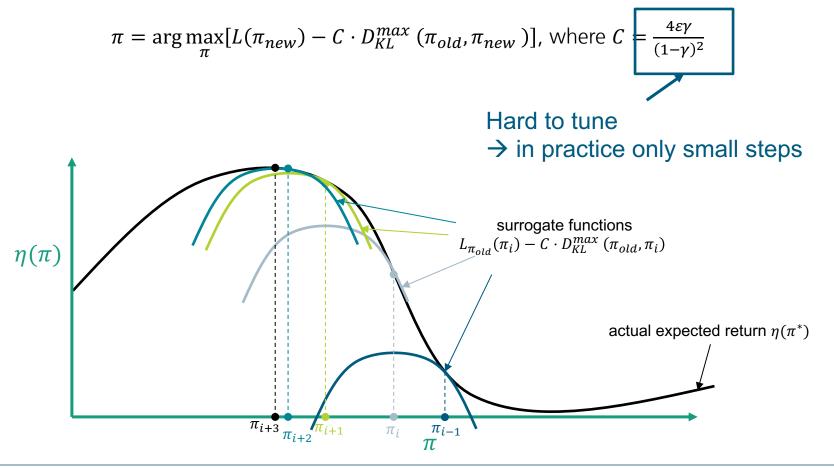




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Minorization-Maximization Algorithm

• A monotonically increasing policy can be defined by:





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Minorization-Maximization Algorithm

• A monotonically increasing policy can be defined by:

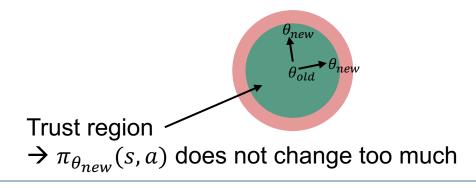
$$\pi = \arg \max_{\pi} [L(\pi_{new}) - C \cdot D_{KL}^{max}(\pi_{old}, \pi_{new})], \text{ where } C = \frac{4\varepsilon\gamma}{(1-\gamma)^2}$$

Side-note:

 A constraint on the KL-divergence between new and old policy (i.e., a trust region constraint) allows larger step sizes while being mathematically equivalent:

$$\pi = \arg \max_{\pi} L_{\pi_{old}}$$
, such that $D_{KL}^{max}(\pi_{old}, \pi) \leq \delta$

• Approximation with *L* is accurate within δ \rightarrow here, monotonic improvement guaranteed





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• Solving the KL-penalized problem (btw: directly over θ not π_{θ})

$$\arg\max_{\theta} \left[L(\tilde{\theta}) - C \cdot D_{KL}^{max}(\theta, \theta_{old}) \right]$$

• Use mean KL-divergence instead of max:

$$\arg \max_{\theta} \begin{bmatrix} L(\hat{\theta}) - C \cdot \overline{D_{KL}}(\theta, \theta_{old}) \end{bmatrix}$$

linear approximation of *L*

$$\arg \max_{\theta} \begin{bmatrix} \frac{\partial}{\partial \theta} L(\theta) \Big|_{\theta = \theta_{old}} \cdot (\theta - \theta_{old}) - \frac{C}{2} (\theta - \theta_{old})^T \cdot \frac{\partial^2}{\partial^2 \theta} \overline{D_{KL}}(\theta, \theta_{old}) \Big|_{\theta = \theta_{old}} \cdot (\theta - \theta_{old}) \end{bmatrix}$$

$$\arg \max_{\theta} \begin{bmatrix} g \cdot (\theta - \theta_{old}) - \frac{C}{2} (\theta - \theta_{old})^T F(\theta - \theta_{old}) \end{bmatrix}$$

$$g = \frac{\partial}{\partial \theta} L(\theta) \Big|_{\theta = \theta_{old}} \quad F = \frac{\partial^2}{\partial^2 \theta} \overline{D_{KL}}(\theta, \theta_{old}) \Big|_{\theta = \theta_{old}}$$



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• Derive with respect to θ and set to 0:

$$\begin{split} 0 &= \frac{\partial}{\partial \theta} \left[g \cdot (\theta - \theta_{old}) - \frac{c}{2} (\theta - \theta_{old})^T F(\theta - \theta_{old}) \right] \\ 0 &= 1 \cdot \left[g - \frac{c}{2} (\theta - \theta_{old})^T F \right] + (\theta - \theta_{old}) \left(-\frac{c}{2} F \right) \\ g - \frac{c}{2} (\theta - \theta_{old})^T F &= \frac{c}{2} (\theta - \theta_{old}) F \\ F^{-1} g - \frac{c}{2} (\theta - \theta_{old})^T &= \frac{c}{2} (\theta - \theta_{old}) \\ F^{-1} g &= c (\theta - \theta_{old}) \\ \frac{1}{c} F^{-1} g &= \theta - \theta_{old} \end{split}$$

• Solution: iterative optimization with Newton's method:

$$\theta_{new} = \theta_{old} + \frac{1}{c} H^{-1} g$$



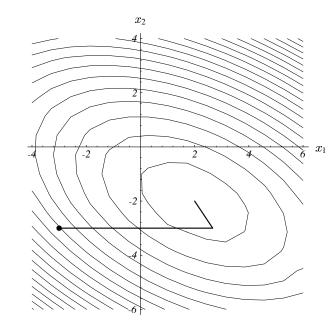
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• The update step

$$\theta_{new} = \theta_{old} + \frac{1}{c} H^{-1} g$$

involves computing the inverse of the Hessian \rightarrow too expensive for large $|\theta|$

- Conjugate Gradient (CG) computes $F^{-1}g$ approximately without forming F explicitly
 - CG solves $x = A^{-1}b$ without explicitly forming A
 - After k iterations, CG has minimized $\frac{1}{2}x^TAx bx$



Jonathan Richard Shewchuk: An Introduction to the Conjugate Gradient Method Without the Agonizing Pain. Edition 1 1/4. 1994.

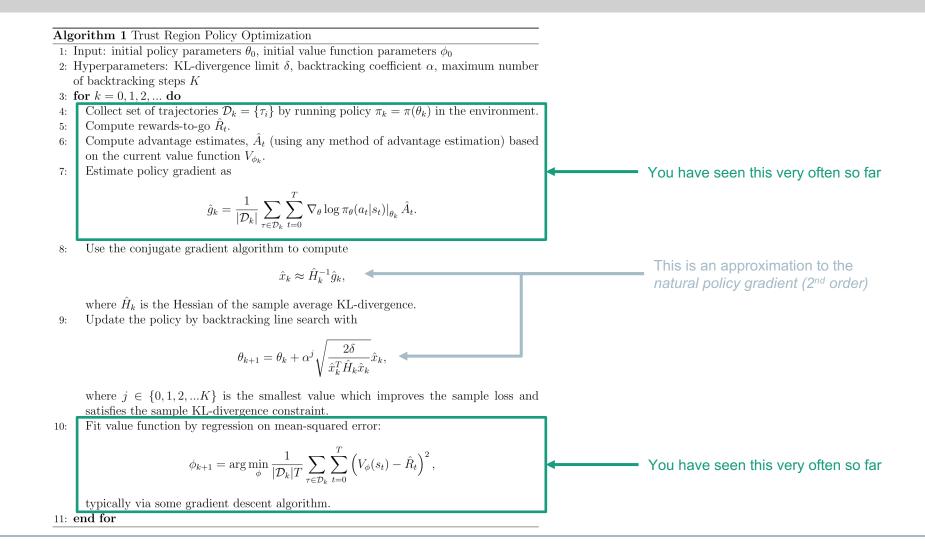


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- Unconstrained problem: $\arg \max_{\theta} L(\theta) C \cdot \overline{D_{KL}}(\theta_{old}, \theta)$
- Constrained problem: $\arg \max_{\alpha} L(\theta)$ subject to $C \cdot \overline{D_{KL}}(\theta, \theta_{old}) \le \delta$
 - δ is a hyper-parameter, remains fixed over the whole learning process
- Solve constrained quadratic problem: compute $F^{-1}g$ and then rescale step to get correct KL:
 - $\max_{\theta} g \cdot (\theta \theta_{old}) \text{ subject to } \frac{1}{2} (\theta \theta_{old})^T F)(\theta \theta_{old}) \leq \delta$
 - 2. Lagrangian: $\mathcal{L}(\theta, \lambda) = g \cdot (\theta \theta_{old}) \frac{c}{2} [(\theta \theta_{old})^T F(\theta \theta_{old}) \delta]$
 - 3. Differentiate with respect to θ and get $\theta \theta_{old} = \frac{1}{c}F^{-1}g$
 - 4. We want $\frac{1}{2}s^T F s = \delta$
 - 5. Given candidate step $s_{unscaled}$ rescale $s = \sqrt{\frac{2\delta}{s_{unscaled} \cdot F \cdot s_{unscaled}}} \cdot s_{unscaled}$



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Intuition & Summary

- We approximate the expected return function locally around the current policy.
 - The accuracy decreases when the new policy and the current policy diverge from each other.
- But we can establish an upper bound for the error.
 - Therefore, we can guarantee a policy improvement if we optimize the local approximation within a trusted region.
 - Outside this region, the bet is off.
- Even it may have a better-calculated value, its range of error fails the improvement guarantee. With such a guarantee inside the trust region, we can locate the optimal policy iteratively. So even it takes a while to prove it mathematically, the reasoning is simple.



Intuition & Summary

- Shortcomings:
 - TRPO minimizes a quadratic equation to approximate the inverse of the Fisher Information Matrix (FIM), i.e., the Hessian
 - To do this for every policy is expensive
 - It requires a large batch of rollouts to approximate it correctly
 - Less sample-efficient to other PG methods when trained with first-order methods such as Adam

 \rightarrow These become a significant issue for large/deep networks!

 And: Yes, TRPO is a complex algorithm that is hard to understand & implement



Policy-based RL

Agenda

- Policy-based RL 1
 - Intro to Policy-based RL
 - Policy Gradients
- Policy-based RL 2
 - Variance & Baselines
 - Actor-Critics
 - Trust-Region Policy Optimization (TRPO)
 - Proximal Policy Optimization (PPO)
 - Deep Deterministic Policy Gradient (DDPG)



- The main motivation behind PPO is the same as for TRPO:
 - Make the biggest possible improvement step
 - Do not step too far such that the performance accidentally collapses
- PPO addresses the shortcomings of TRPO:
 - PPO uses 1st order methods with a few tricks
 - Significantly simpler to implement
 - Shows similar performance to TRPO (empirically)
- There are two variants:
 - PPO-penalty: TRPO with KL-penalization instead of constraint (penalty coefficient is adjusted and scaled automatically over the course of training: Adaptive KL Penalty Coefficient)
 - PPO-clip: no constraints! Adds a clipping to the objective function to remove incentives to move too far

Spoiler: PPO is (1) much simpler to understand and to implement, and (2) much better (empirically)



Where are we so far?

• TRPO maximizes a "surrogate" objective subject to a constraint on the size of the policy update:

$$\max_{\theta} \widehat{\mathbb{E}}_t \left[\frac{\pi_{\theta}(a_t | s_t)}{\pi_{\theta_{old}}(a_t | s_t)} \hat{A}_t \right], \quad \text{subject to} \quad \widehat{\mathbb{E}}_t \left[KL \left[\pi_{\theta_{old}}(\cdot | s_t), \pi_{\theta}(\cdot | s_t) \right] \right] \le \delta$$

with θ_{old} being the policy parameters before the update

- We did not explicitly formulate it like this, but the intuition behind it is:
 - We want to measure how π_{θ} performs relative to $\pi_{\theta_{old}}$ (using data from the old policy)
 - The original objective (see TRPO slides) can be exactly reformulated to this one
- We can solve this with CG after making a linear approximation to the objective and a quadratic approximation to the KLconstraint



Where are we so far?

• TRPO maximizes a "surrogate" objective subject to a constraint on the size of the policy update:

$$\max_{\theta} \widehat{\mathbb{E}}_t \left[\frac{\pi_{\theta}(a_t | s_t)}{\pi_{\theta_{old}}(a_t | s_t)} \hat{A}_t \right], \quad \text{subject to} \quad \widehat{\mathbb{E}}_t \left[KL \left[\pi_{\theta_{old}}(\cdot | s_t), \pi_{\theta}(\cdot | s_t) \right] \right] \leq \delta$$

with θ_{old} being the policy parameters before the update

• Let us define the probability ratio $r_t(\theta)$:

$$r_t(\theta) = \frac{\pi_{\theta}(a_t|s_t)}{\pi_{\theta_{old}}(a_t|s_t)} \quad \text{, i.e., } r_t(\theta_{old}) = 1.$$

In other words, TRPO maximizes the following objective:

$$L^{CPI}(\theta) = \widehat{\mathbb{E}}_t \big[r_t(\theta) \hat{A}_t \big]$$

penalizing changes to the policy that move $r_t(\theta)$ (too far) away from 1.



• The PPO objective we want to maximize is given by

$$L(\theta) = \widehat{\mathbb{E}}_t \left[\min \left(r_t(\theta) \hat{A}_t, \operatorname{clip}(r_t(\theta), 1 - \epsilon, 1 + \epsilon) \hat{A}_t \right) \right]$$

where ϵ is a hyperparameter (i.e., 0.1 or 0.2) that defines how far π_{new} may go away from π_{old}

- First term inside the min is $L^{CPI}(\theta)$
- Second term inside the min clips the probability ratio
 \rightarrow removes the incentive for moving r_t outside of the interval [1 ϵ , 1 + ϵ]

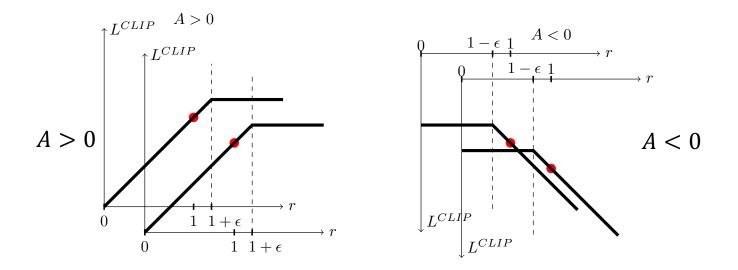
We take the minimum of the clipped and unclipped objective
 → the final objective is a lower bound (i.e., a pessimistic bound) on the unclipped objective

 $g(\epsilon, \hat{A}_t) = \begin{cases} (1+\epsilon)A, \text{ if } A \ge 0\\ (1-\epsilon)A, \text{ if } A < 0 \end{cases}$

 $=g\left(\epsilon,\hat{A}_{t}(s,a)\right)$



• The clipping operator is a *pessimistic bound* of the unclipped objective



- Plot show a single timestep of the surrogate function L^{CLIP} as a function of r
- The red circle shows the starting point for the optimization, i.e., r = 1



• Automated Tuning of the gradient step *without calculating the Hessian*

Algorithm 1 PPO-Clip

1: Input: initial policy parameters θ_0 , initial value function parameters ϕ_0

2: for k = 0, 1, 2, ... do

- 3: Collect set of trajectories $\mathcal{D}_k = \{\tau_i\}$ by running policy $\pi_k = \pi(\theta_k)$ in the environment.
- 4: Compute rewards-to-go \hat{R}_t .
- 5: Compute advantage estimates, \hat{A}_t (using any method of advantage estimation) based on the current value function V_{ϕ_k} .
- 6: Update the policy by maximizing the PPO-Clip objective:

$$\theta_{k+1} = \arg\max_{\theta} \frac{1}{|\mathcal{D}_k|T} \sum_{\tau \in \mathcal{D}_k} \sum_{t=0}^T \min\left(\frac{\pi_{\theta}(a_t|s_t)}{\pi_{\theta_k}(a_t|s_t)} A^{\pi_{\theta_k}}(s_t, a_t), \ g(\epsilon, A^{\pi_{\theta_k}}(s_t, a_t))\right),$$

Advantage Clipping for conservative policy updates

typically via stochastic gradient ascent with Adam.

7: Fit value function by regression on mean-squared error:

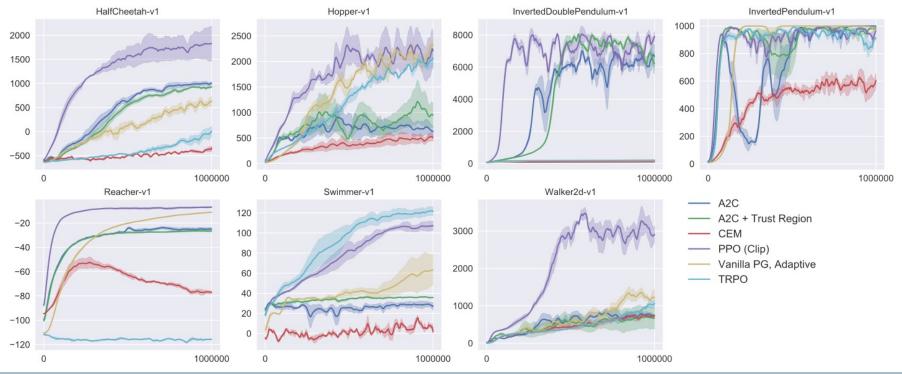
$$\phi_{k+1} = \arg\min_{\phi} \frac{1}{|\mathcal{D}_k|T} \sum_{\tau \in \mathcal{D}_k} \sum_{t=0}^T \left(V_{\phi}(s_t) - \hat{R}_t \right)^2,$$

typically via some gradient descent algorithm.

8: end for



- Results of PPO-clip:
 - Against well-known competitors
 - On well-known environments
- Those results are impressive!





Practical Considerations

 Algorithm 1 PPO, Actor-Critic Style

 for iteration=1, 2, ..., N do

 for actor=1, 2, ..., N do

 Run policy $\pi_{\theta_{old}}$ in environment for T timesteps

 Compute advantage estimates $\hat{A}_1, \ldots, \hat{A}_T$

 end for

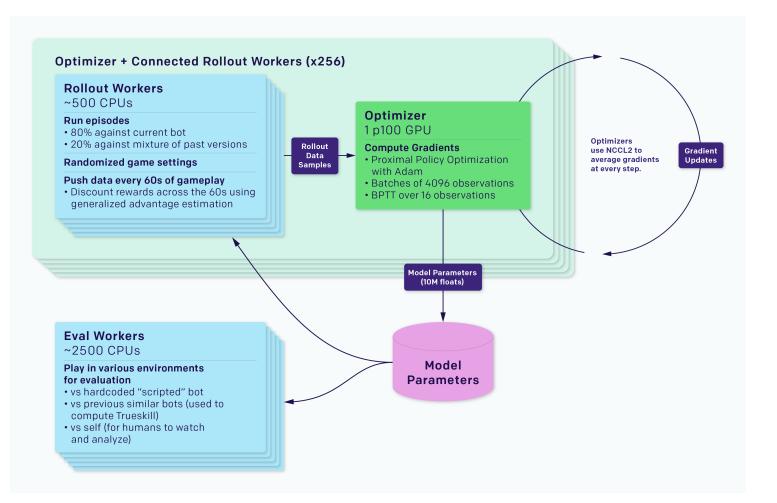
 Optimize surrogate L wrt θ , with K epochs and minibatch size $M \leq NT$
 $\theta_{old} \leftarrow \theta$

 end for

- There is two alternating threads in PPO:
 - 1. Policy interacts with the environment, collects data and computes advantage estimates (using fitted baselines estimates)
 - 2. 2nd thread collects all the experiences and runs SGD to optimize the policy using the clipped objective



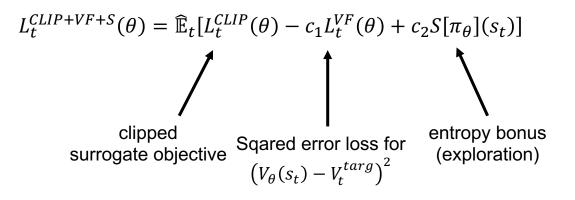
PPO in Action: OpenAl Five on DOTA II





Practical Considerations

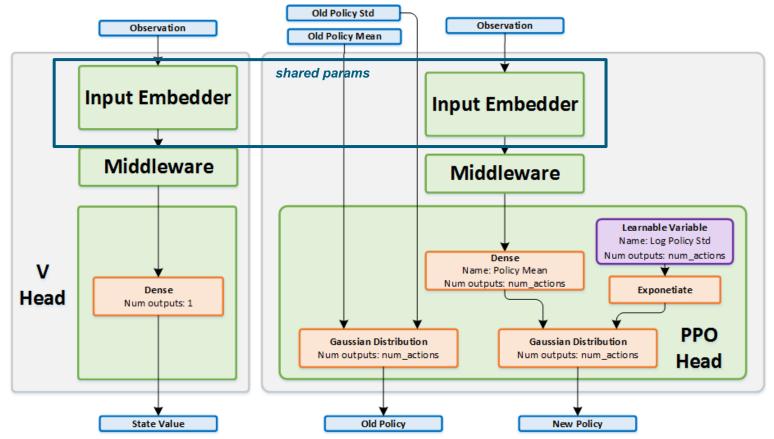
- One more thing....
- PPO combines a few more things in the final objective:



Note: you likely need similar features to represent the policy and the state-values
 → OpenAl Five shares parameters between the policy network and the value network
 → Both error terms are combined in a single loss function



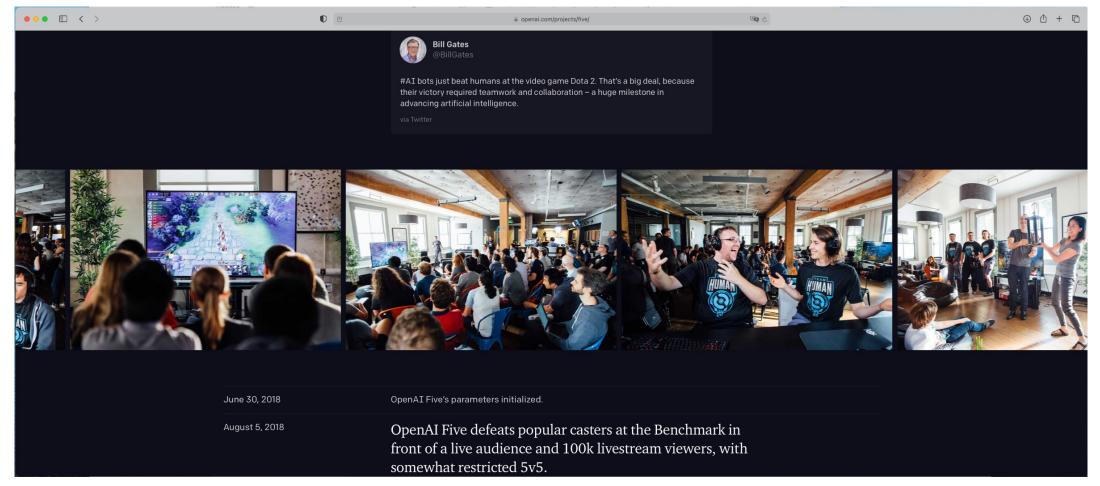
PPO in Action: OpenAl Five vs. DOTA II



https://intellabs.github.io/coach/components/agents/policy_optimization/ppo.html



PPO in Action: OpenAl Five vs. DOTA II



https://openai.com/projects/five/



PPO in Action: OpenAl Five vs. DOTA II

High-level info:

- 180 years of self-play per day and hero (~900 yrs/day), no human data
- Running on 256 P100 GPUs and 128,000 CPU cores

Technical stats:

- Observation size: ~36.8kB @ ~7Hz
- Batch size: 1,048,576 observations = 36 GB \odot
- Separate single-layer, 1024 unit LSTM per hero
- See <u>https://openai.com/blog/openai-five/</u> for an interactive demo!
- Reward: net worth, kills, deaths, assist, last hits, etc.
- "Team spirit" trade own rewards over team reward (heroes do not communicate)

Challenge: exploring combinatorial-vast space of combining actions w/ long planning horizons

- 80% of games against itself, 20% against past selves (avoid "strategy collapse")
- After several hours: concepts such as laning, farming or fighting emerged
- After several days: basic human strategies such as steal bounty runes from opponents, rotate heroes around the map to gain lane advantage etc.



Policy-based RL

Agenda

- Policy-based RL 1
 - Intro to Policy-based RL
 - Policy Gradients
- Policy-based RL 2
 - Variance & Baselines
 - Actor-Critics
 - Trust-Region Policy Optimization (TRPO)
 - Proximal Policy Optimization (PPO)
 - Deep Deterministic Policy Gradient (DDPG)



Question: can we also use ideas from value-based RL for continuous action spaces?

- Why can't we use Q-Learning and DQNs?
- \rightarrow Not so easily!

But what was the original problem with Q-Learning?

- Q-Learning and variants (including DQNs) do not work with continuous actions
- Why is that? Remember:
 - We calculated the targets $r_t + \gamma \cdot \max_{a' \in \mathcal{A}} \hat{Q}(s', a'; \theta_{i-1})$ by a single pass through the network
 - Our network was "static" and had $|\mathcal{A}|$ outputs
- Evaluating a continuous action space requires an exhaustive search over the available actions (and this becomes highly non-trivial!)¹

¹ However, see approaches such as Lim et al.: Actor-Expert: A Framework for using Q-learning in Continuous Action Spaces.



- DDPG learns a Q-function and a policy
 - Off-policy data + Bellman equation to learn Q-function
 - Make use of the Q-function to learn the policy
- The intuition relies on Q-learning:
 - If you know $Q^*(s, a)$ then in each state the optimal action $a^*(s)$ can be simply found by

 $a^*(s) = \arg\max_a Q^*(s, a)$

DDPG jointly learns approximations to $Q^*(s, a)$ and $a^*(s)$

...and specifically adapts this for continuous action spaces!



Main idea

- We assume the function $Q^*(s, a)$ to be differentiable with respect to a
- This allows for a gradient-based learning rule for a policy $\mu(s)$ that exploits this
- Instead of exhaustively looking for $\max_{a} Q(s, a)$ we approximate it:

 $\max_a Q(s,a) \approx Q(s,\mu(s))$

- We look at two sides of DDPG:
 - 1. Its Q-Learning side
 - 2. Its policy gradient side



The Q-Learning side of DDPG

• Recap the Bellman optimality equation:

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

• Then we can optimize the mean-squared Bellman error (MSBE) (neural network parameters ϕ , set of transitions $D, d \in \{0; 1\}$ indicates if s' is terminal):

$$L(\phi, D) = \mathbb{E}_{(s,a,r,s',d)) \sim D} \left[\left(Q_{\phi}(s,a) - \left(r + \gamma(1-d) \max_{a'} Q_{\phi}(s',a') \right) \right)^2 \right]$$

- For optimization using SGD we apply the well-known tricks:
 - Replay buffers (off-policy!)
 - Use target networks and update it with delay by $\phi_{targ} \leftarrow p\phi_{targ} + (1-p)\phi$, $p \in [0; 1]$



The Q-Learning side of DDPG

- Calculating the max over the actions in the target:
 - Target policy network μ_{θ} computes an action that approximately maximizes $Q_{\phi_{targ}}$
 - Target policy updates also computed using polyak averaging (see above)
 - Q-Learning in DDPG minimizes using SGD:

$$L(\phi, D) = \mathbb{E}_{(s,a,r,s',d)\sim D}\left[\left(Q_{\phi}(s,a) - \left(r + \gamma(1-d)Q_{\phi_{targ}}\left(s', \mu_{\theta_{targ}}(s')\right)\right)\right)^{2}\right]$$



The Policy Learning side of DDPG: simple

- Policy $\mu_{\theta}(s)$ is deterministic
- $\mu_{\theta}(s)$ should return the action that maximizes $Q_{\phi}(s, a)$
- Action space is continuous, and we assume Q to be differentiable with respect to actions a
- Hence, we can use gradient ascent (with respect to the policy parameters θ only) and solve:

 $\max_{\theta} \mathbb{E}_{s \sim D} \left[Q_{\phi}(s, \mu_{\theta}(s)) \right]$

(Q-function parameters ϕ are treated as constants here)



Exploration-Exploitation

- DDPG trains off-policy; policy is deterministic
- Hence an on-policy exploration is often not enough
- Solution: add noise to the actions at training time
 - Originally time-correlated Ornstein-Uhlenbeck (OU) process noise has been proposed
 - More recent work suggests to use zero-mean Gaussian noise as it is simpler and exhibits same performance
 - Over the course of training, we may reduce the scale of noise (but there is only a limited effect from this)
 - At test time we (of course) omit to add the noise



Pseudo Code

| Algorithm 1 DDPG algorithm | |
|--|---|
| Randomly initialize critic network $Q(s, a \theta^Q)$ and actor $\mu(s \theta^\mu)$ with weights θ^Q and θ^μ . Initialize target network Q' and μ' with weights $\theta^{Q'} \leftarrow \theta^Q$, $\theta^{\mu'} \leftarrow \theta^\mu$ Initialize replay buffer R | Use target networks (as in DQN) for both the actor and the critic |
| for episode = 1. M do | As we don't have a stochastic policy, |
| Initialize a random process \mathcal{N} for action exploration Receive initial observation state s_1 | we have to define a process for exploration |
| for $t = 1$, T do | • |
| Select action $a_t = \mu(s_t \theta^{\mu}) + \mathcal{N}_t$ according to the current policy and exploration noise Execute action a_t and observe reward r_t and observe new state s_{t+1} Store transition (s_t, a_t, r_t, s_{t+1}) in R Sample a random minibatch of N transitions (s_i, a_i, r_i, s_{i+1}) from R | Execute action, store transition in the replay buffer and sample at random some transitions (exactly as in DQN) |
| Set $y_i = r_i + \gamma Q'(s_{i+1}, \mu'(s_{i+1} \theta^{\mu'}) \theta^{Q'})$ Update critic by minimizing the loss: $L = \frac{1}{N} \sum_i (y_i - Q(s_i, a_i \theta^Q))^2$ | Update critic using the target networks for both the Actor and the Critic |
| Update the actor policy using the sampled policy gradient: | |
| $\nabla_{\theta^{\mu}} J \approx \frac{1}{N} \sum_{i} \nabla_{a} Q(s, a \theta^{Q}) _{s=s_{i}, a=\mu(s_{i})} \nabla_{\theta^{\mu}} \mu(s \theta^{\mu}) _{s_{i}}$ | Update the Actor using the Deterministic Policy Gradient Theorem (Silver et al. 2014) |
| Update the target networks: $\theta^{Q'} \leftarrow \tau \theta^Q + (1-\tau) \theta^{Q'}$ | Progressively update the target networks |
| $\begin{aligned} \theta^{Q'} &\leftarrow \tau \theta^Q + (1 - \tau) \theta^{Q'} \\ \theta^{\mu'} &\leftarrow \tau \theta^{\mu} + (1 - \tau) \theta^{\mu'} \end{aligned}$ | |

end for end for



Summary

- Special case in actor-critic-algorithms:
- Works only for continuous action spaces
- Is an off-policy algorithm utilizing the replay buffer trick from DQN
- Solves for deterministic policies instead of stochastic ones
- + impressive results both in simulation and in real world problems
- + one of the de-facto algorithms to use for continuous (or very large) action spaces
- very sensitive to the exploration process
- can be hard to tune sensitive to hyperparameters
- slower (wall clock time) compared to other actor-critic-algorithms



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