

Extensions to Policy Gradient

Reinforcement Learning School of Data Science University of Virginia

Last updated: March 25, 2025

<u>Agenda</u>

- > Recap of REINFORCE
- > Actor-Critic Methods
- > Trust Region Methods
- > Deep Deterministic Policy Gradient (DDPG)

Recap of REINFORCE

REINFORCE

This was our first policy gradient (PG) method

All PG methods are on-policy algorithms

REINFORCE is a Monte Carlo Gradient algorithm (uses full trajectories)

It works by:

- 1) simulating paths
- 2) calculating returns G from each path (evaluate)
- 3) taking update steps based on return and gradient of policy (improve)

REINFORCE, A Monte-Carlo Policy-Gradient Method (episodic)

Input: a differentiable policy parameterization $\pi(a|s, \theta), \forall a \in A, s \in S, \theta \in \mathbb{R}^n$

Initialize policy weights θ

Repeat forever:

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

For each step of the episode t = 0, ..., T-1:

$$G_t \leftarrow \text{return from step } t$$

$$\boldsymbol{\theta} \leftarrow \boldsymbol{\theta} + \alpha \gamma^t G_t \nabla_{\boldsymbol{\theta}} \log \pi(A_t | S_t, \boldsymbol{\theta})$$

REINFORCE Intuition

As we generate sample paths, we calculate return G over the paths.

The update step will increase parameter vector in direction of:

- greater return
- greater increase of probability repeating action A_t on future visits to state S_t

REINFORCE, A Monte-Carlo Policy-Gradient Method (episodic)

Input: a differentiable policy parameterization $\pi(a|s, \theta), \forall a \in A, s \in S, \theta \in \mathbb{R}^n$

Initialize policy weights θ

Repeat forever:

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

For each step of the episode t = 0, ..., T-1:

 $G_t \leftarrow \text{return from step } t$

$$\boldsymbol{\theta} \leftarrow \boldsymbol{\theta} + \alpha \gamma^t G_t \nabla_{\boldsymbol{\theta}} \log \pi(A_t | S_t, \boldsymbol{\theta})$$

REINFORCE Limitations

REINFORCE can have high variance and converge slowly

From any sampled (s,a) there may be many very different return estimates

We are estimating the gradient of the performance measure $\nabla \hat{J(\theta_t)}$

Can be noisy due to factors including:

- > Noise in environment
- > Sparse rewards
- > Length of sample trajectories can vary

REINFORCE with Baseline

REINFORCE can have high variance and converge slowly

We want to take positive gradient steps in direction of parameters that lead to high-reward trajectories We want to take negative steps for low-reward trajectories

Rather than working w rewards, we can work with relative rewards



REINFORCE with Baseline, contd.

Positive gradient steps in direction of parameters that lead to high *relative* reward trajectories Negative steps for low *relative* reward trajectories

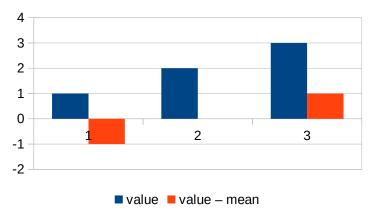
Q: What makes sense for relative reward?

One solution: subtract the average trajectory reward

This is actually the state value function V(s)

By subtracting off a baseline, it reduces variance

We will still have an unbiased estimate for gradient

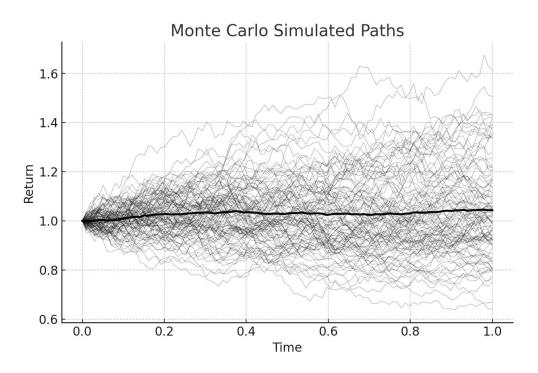


REINFORCE with Baseline, contd.

Plot showing many simulated paths. A gain can be computed for each path.

The dark curve shows average trajectory values. The gain can be computed and it represents V(s).

This can be treated as baseline.



REINFORCE with Baseline Algorithm

REINFORCE can have high variance and converge slowly

One solution: subtract baseline, which can be state value function

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 $\boldsymbol{\theta} \in \mathbb{R}^{d'}$ and state-value weights $\mathbf{w} \in \mathbb{R}^d$ (e.g., to 0)

Loop forever (for each episode):

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

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

$$G \leftarrow \sum_{k=t+1}^{T} R_k$$
 $\delta \leftarrow G - \hat{v}(S_t, \mathbf{w})$ Value function as baseline (G_t)

$$\mathbf{w} \leftarrow \mathbf{w} + \alpha^{\mathbf{w}} \gamma^t \delta \nabla \hat{v}(S_t, \mathbf{w})$$

$$\boldsymbol{\theta} \leftarrow \boldsymbol{\theta} + \alpha^{\boldsymbol{\theta}} \gamma^t \delta \nabla \ln \pi(A_t | S_t, \boldsymbol{\theta})$$

Actor-Critic Methods

Introducing the Critic

The state value function estimate is based on first state of each transition (time *t*)

$$\delta \leftarrow G - \hat{v}(S_t, \mathbf{w})$$

Another option is to use state value estimates from multiple time steps

Given two time steps, we can estimate a one-step return: $\gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})$

State value one step into future,
discounted to present value

Introducing the Critic

The state value function estimate is based on first state of each transition (time *t*)

$$\delta \leftarrow G - \hat{v}(S_t, \mathbf{w})$$

Another option is to use state value estimates from multiple time steps

Given two time steps, we can estimate a one-step return: $\gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})$

We can collect many trajectories, calculate the returns, and train a neural network

By using this model, it reduces the variance from the samples

It can be used to evaluate the policy, and is called the critic

The Actor

The policy network is called the *actor*

It guides the behavior of the agent

The actor uses a separate set of parameters

Actor-Critic

We now have two different parametrized functions:

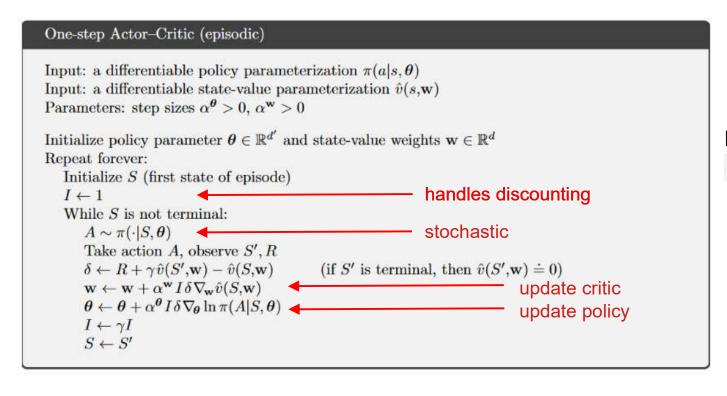
- > Policy function (actor)
- > State-value function (*critic*). Evaluates the policy.

For each function the params are updated on each pass

Can use neural networks for the functions, given sufficient training data

Actor-Critic Algorithm

There are two parametrized functions: one for policy (*actor*), one for state value (*critic*) For each function the params are updated on each pass



Estimate of one-step return $\gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})$

Actor-Critic vs REINFORCE w Baseline

To summarize the difference:

Actor critic:
$$\delta \leftarrow R + \gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})$$

REINFORCE with baseline:
$$\delta \leftarrow G - \hat{v}(S, \mathbf{w})$$

- > The purpose of both is to reduce variance
- > Actor critic is a refinement that uses two time points

Reappearance of the Advantage Function

Recall the advantage function which measures the benefit of a particular action:

$$A^{\pi}(s, a) = Q^{\pi}(s, a) - V^{\pi}(s)$$

This appears in the critic term (line 4 circled at right)

The Q-value estimate can be calculated from single step transitions. We have done this with TD learning.

```
While S is not terminal:

A \sim \pi(\cdot|S, \theta)

Take action A, observe S', R

\delta \leftarrow R + \gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})

\mathbf{w} \leftarrow \mathbf{w} + \alpha^{\mathbf{w}} I \delta \nabla_{\mathbf{w}} \hat{v}(S, \mathbf{w})

\theta \leftarrow \theta + \alpha^{\theta} I \delta \nabla_{\theta} \ln \pi(A|S, \theta)

I \leftarrow \gamma I

S \leftarrow S'
```

When estimating the advantage this way, the approach is sometimes called Advantage Actor Critic or A2C

Actor-Critic Summary

Roles:

- Actor learns the policy which drives actions
- Critic estimates value function V(s) or action-value function Q(s,a)

Updating:

- Actor updates policy based on critic feedback
- Critic updates value estimates based on TD error

Advantages:

- Stability and Efficiency: Critic reduces variance in actor policy updates
- Continuous Action Spaces: AC methods handle continuous action spaces

```
While S is not terminal:

A \sim \pi(\cdot|S, \theta)

Take action A, observe S', R

\delta \leftarrow R + \gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})

\mathbf{w} \leftarrow \mathbf{w} + \alpha^{\mathbf{w}} I \delta \nabla_{\mathbf{w}} \hat{v}(S, \mathbf{w})

\theta \leftarrow \theta + \alpha^{\theta} I \delta \nabla_{\theta} \ln \pi(A|S, \theta)

I \leftarrow \gamma I

S \leftarrow S'
```

Trust Region Methods

Trust Region Policy Optimization (TRPO)

In policy gradient, a single bad update to policy parameters can have large differences in performance

TRPO methods control step size to avoid bad updates

Source: OpenAl Spinning Up

21

Trust Region Policy Optimization (TRPO)

In policy gradient, a single bad update to policy parameters can have large differences in performance

TRPO methods control step size to avoid bad updates

TRPO methods have shown a large improvement over Actor Critic methods

TRPO updates policies by taking largest possible step to improve performance However, there is a constraint: new policy must not differ "too much" from old policy

Constraint is formalized using KL-divergence: distance between prob distributions

Source: OpenAl <u>Spinning Up</u>

K-L Divergence Definition

Given probability distributions P and Q defined on sample space \mathcal{X} ,

KL divergence measures log-difference between *P* and *Q*, where expectation is measured relative to *P*

Discrete case:
$$D_{ ext{KL}}(P \parallel Q) = \sum_{x \in \mathcal{X}} P(x) \log \left(rac{P(x)}{Q(x)}
ight)$$
 ,

Continuous case:
$$D_{ ext{KL}}(P \parallel Q) = \int_{-\infty}^{\infty} p(x) \log \left(rac{p(x)}{q(x)}
ight) dx,$$

See here for details and a nice example

K-L Divergence Facts

 $D_{\mathrm{KL}}(P \parallel Q) = \sum_{x \in \mathcal{X}} P(x) \log igg(rac{P(x)}{Q(x)}igg),$

KL divergence also called *relative entropy*

Interpreted as average difference of number of bits required for encoding samples of P using a code optimized for Q

Value is zero when distributions P and Q are identical

When P and Q are not identical, KL divergence is not symmetric: $D_{KL}(P \parallel Q)
eq D_{KL}(Q \parallel P)$

The triangle inequality also does not hold

Thus, KL divergence is not a metric

TRPO in Practice

TRPO is relatively hard to implement in practice.

Proximal Policy Optimization (PPO) is easier to implement

Thus, we see a quick overview of TRPO

Relative Policy Performance

Define surrogate advantage which measures relative performance of policies

 $\pi_{ heta}$ New policy

 $\pi_{ heta_k}$ Old policy

 $\mathcal{L}(heta_k, heta)$ Surrogate advantage of new policy over old policy

$$\mathcal{L}(\theta_k, \theta) = \mathop{\mathbf{E}}_{s, a \sim \pi_{\theta_k}} \left[\frac{\pi_{\theta}(a|s)}{\pi_{\theta_k}(a|s)} A^{\pi_{\theta_k}}(s, a) \right]$$

Expectation is taken under new policy
As usual, we estimate expectation by sampling

Relative Policy Performance: Constraint

Now we introduce KL divergence:

$$\bar{D}_{KL}(\theta||\theta_k) = \mathop{\mathbf{E}}_{s \sim \pi_{\theta_k}} \left[D_{KL} \left(\pi_{\theta}(\cdot|s) || \pi_{\theta_k}(\cdot|s) \right) \right]$$

And we constrain it:

$$\bar{D}_{KL}(\theta||\theta_k) \le \delta$$

Relative Policy Performance: Key Equations

An update step is subject to these key equations:

$$heta_{k+1} = \arg\max_{ heta} \mathcal{L}(heta_k, heta)$$
 (objective) s.t. $ar{D}_{KL}(heta|| heta_k) \leq \delta$ (constraint)

Relative Policy Performance: Key Equations

An update step is subject to these key equations:

$$heta_{k+1} = \arg\max_{ heta} \mathcal{L}(heta_k, heta)$$
 (objective) s.t. $\bar{D}_{KL}(heta|| heta_k) \leq \delta$ (constraint)

Objective: maximize surrogate advantage

Constraint: limit KL divergence

Taylor Approximation

$$\mathcal{L}(\theta_k, \theta) = \mathop{\mathbf{E}}_{s, a \sim \pi_{\theta_k}} \left[\frac{\pi_{\theta}(a|s)}{\pi_{\theta_k}(a|s)} A^{\pi_{\theta_k}}(s, a) \right]$$

TRPO uses first-order Taylor expansion for easier solution

Recall:

For function f(x) differentiable at point x = a, can consider approximations:

$$P_1(x) = f(a) + f'(a)(x-a)$$
 (First order)

$$P_2(x)=f(a)+f'(a)(x-a)+rac{f''(a)}{2}(x-a)^2.$$
 (Second order)

30

Taylor Approximation, contd.

$$\mathcal{L}(\theta_k, \theta) = \mathop{\mathbf{E}}_{s, a \sim \pi_{\theta_k}} \left[\frac{\pi_{\theta}(a|s)}{\pi_{\theta_k}(a|s)} A^{\pi_{\theta_k}}(s, a) \right]$$

TRPO uses Taylor expansions of objective, constraint for easier solution

Approximation has this form:

$$\mathcal{L}(\theta_k, \theta) \approx g^T(\theta - \theta_k)$$
$$\bar{D}_{KL}(\theta||\theta_k) \approx \frac{1}{2}(\theta - \theta_k)^T H(\theta - \theta_k)$$

where g is gradient and H is Hessian (2nd order derivative) of surrogate This can be useful where parameters are sufficiently close

Optimization Problem

Framing as an optimization problem we have:

$$\theta_{k+1} = \arg \max_{\theta} g^T(\theta - \theta_k)$$

s.t. $\frac{1}{2}(\theta - \theta_k)^T H(\theta - \theta_k) \le \delta$

Optimization Problem

Framing as an optimization problem we have:

$$\theta_{k+1} = \arg \max_{\theta} g^T(\theta - \theta_k)$$

s.t. $\frac{1}{2}(\theta - \theta_k)^T H(\theta - \theta_k) \le \delta$

The objective of TRPO is to maximize the gradient of surrogate advantage function

This is equivalent to policy gradient method, which moves in direction of gradient

Recall our objective to maximize this quantity from PG:

$$abla v_{\pi}(s) =
abla \left[\sum_{a} \pi(a|s) q_{\pi}(s,a)
ight]$$

Solving the Optimization Problem

Framing as an optimization problem we have:

$$\theta_{k+1} = \arg \max_{\theta} g^T(\theta - \theta_k)$$

s.t. $\frac{1}{2}(\theta - \theta_k)^T H(\theta - \theta_k) \le \delta$

Can be solved with this update rule:

$$\theta_{k+1} = \theta_k + \alpha^j \sqrt{\frac{2\delta}{g^T H^{-1} g}} H^{-1} g,$$

Hessian H needs to be inverted

TRPO Challenges

Several challenges make this hard to implement (see Bilgin p249 for list)

Includes:

- > Taylor approximation may be violated
- > For policy network with massive number of parameters,
 - Inverting Hessian H may be hard
 - May be hard to store H-1

Proximal Policy Optimization (PPO)

Same approach as TRPO:

- > Take policy improvement steps
- > Limit distance between old and new policy to retain good performance

TRPO used second-order Taylor expansion

PPO uses first-order methods and tricks
Easier to implement, competitive performance

PPO Variants

PPO-Penality: penalize KL-divergence in objective function

PPO-Clip: No KL-divergence term. Specialized clipping of objective fcn.

OpenAl uses PPO-Clip

Objective function from earlier:

$$\mathcal{L}(\theta_k, \theta) = \mathop{\mathbf{E}}_{s, a \sim \pi_{\theta_k}} \left[\frac{\pi_{\theta}(a|s)}{\pi_{\theta_k}(a|s)} A^{\pi_{\theta_k}}(s, a) \right]$$

We will limit policy changes based on ratio

PPO-Clip

Clipped form of objective function *L* has form below:

$$L(s, a, \theta_k, \theta) = \min \left(\frac{\pi_{\theta}(a|s)}{\pi_{\theta_k}(a|s)} A^{\pi_{\theta_k}}(s, a), \quad g(\epsilon, A^{\pi_{\theta_k}}(s, a)) \right),$$

This can be computed based on sample trajectories

The min() operator will serve to limit changes based on ratio of new and old policies

PPO-Clip: Example of Positive Case

If advantage for a state-action pair is positive, the contribution to objective reduces to

$$L(s, a, \theta_k, \theta) = \min\left(\frac{\pi_{\theta}(a|s)}{\pi_{\theta_k}(a|s)}, (1+\epsilon)\right) A^{\pi_{\theta_k}}(s, a).$$

Any benefit of changing the policy beyond clipped limit is removed

This acts as a regularizer

PPO-Clip Algorithm

- 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:

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

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

Deep Deterministic Policy Gradient (DDPG)

DPG - Motivation

We studied a lot of approaches using Q-function

For discrete / low cardinality action space, this makes sense: $a^*(s) = \arg\max_a Q^*(s,a)$

$$a^*(s) = \arg\max_{a} Q^*(s, a)$$

For continuous / high cardinality action space, the *max* is too costly

Instead, a function approximator is used in DDPG

This is the same idea that we used for state spaces with DQN

DDPG - Parametrization

The idea is to use: $\max_a Q(s,a) \approx Q(s,\mu(s))$

for policy $\mu(s)$

This assumes the policy function is differentiable wrt action

The policy is now deterministic

Uses the same features as DQN:

- Replay buffer for stored experiences
- Target network with soft updates (Polyak averaging):

$$\phi_{\text{targ}} \leftarrow \rho \phi_{\text{targ}} + (1 - \rho) \phi$$

DDPG - Training

The policy is deterministic

If the agent explores on-policy, it may not try enough actions

The trick for better exploration is to add noise to the actions during training

DDPG - Algorithm

Add noise for more exploration

Algorithm 1 Deep Deterministic Policy Gradient

- 1: Input: initial policy parameters θ , Q-function parameters ϕ , empty replay buffer \mathcal{D}
- 2: Set target parameters equal to main parameters $\theta_{\text{targ}} \leftarrow \theta$, $\phi_{\text{targ}} \leftarrow \phi$
- 3: repea
 - 4: Observe state s and select action $a = \text{clip}(\mu_{\theta}(s) + \epsilon, a_{Low}, a_{High})$, where $\epsilon \sim \mathcal{N}$
- 5: Execute a in the environment
- 6: Observe next state s', reward r, and done signal d to indicate whether s' is terminal
- 7: Store (s, a, r, s', d) in replay buffer \mathcal{D}
- 8: If s' is terminal, reset environment state.
- 9: **if** it's time to update **then**
- 10: **for** however many updates **do**
- 11: Randomly sample a batch of transitions, $B = \{(s, a, r, s', d)\}$ from \mathcal{D}
- 12: Compute targets

$$y(r, s', d) = r + \gamma (1 - d) Q_{\phi_{\text{targ}}}(s', \mu_{\theta_{\text{targ}}}(s'))$$

13: Update Q-function by one step of gradient descent using

$$\nabla_{\phi} \frac{1}{|B|} \sum_{(s,a,r,s',d) \in B} (Q_{\phi}(s,a) - y(r,s',d))^2$$

14: Update policy by one step of gradient ascent using

$$\nabla_{\theta} \frac{1}{|B|} \sum_{s \in B} Q_{\phi}(s, \mu_{\theta}(s))$$

15: Update target networks with

Soft updates

$$\phi_{\text{targ}} \leftarrow \rho \phi_{\text{targ}} + (1 - \rho) \phi
\theta_{\text{targ}} \leftarrow \rho \theta_{\text{targ}} + (1 - \rho) \theta$$

- 16: end for
- 17: end if
- 18: until convergence

Source: https://spinningup.openai.com/en/latest/algorithms/ddpg.html