

Policy Search Methods

(CSC2547: Introduction to Reinforcement Learning)

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Introduction

- So far, we have described methods for computing the optimal policy based on the computation of the value function.
- Only the value function was **explicitly** represented.
 - The policy could be computed based on it.
- There are methods based on explicit representation of the policy and optimizing the performance of the agent by searching in the space of policies.
- We call them **policy search methods**.
- Hybrid methods: explicit representation of both value and policy.

Policy Parametrization

- Consider a stochastic policy $\pi_\theta : \mathcal{X} \rightarrow \mathcal{M}(\mathcal{A})$ that is parameterized by a $\theta \in \Theta$.
- The set Θ is the parameter space, e.g., a subset of \mathbb{R}^p .
- The space of all parameterized policies:

$$\Pi_\Theta = \{ \pi_\theta : \mathcal{X} \rightarrow \mathcal{M}(\mathcal{A}) : \theta \in \Theta \}. \quad (1)$$

- This space depends on the mapping π_θ and Θ .

Policy Parametrization: Examples

- Many choices for how we can parameterize a policy π_θ .
- A generic example is based on the Boltzmann (or softmax) distribution.
- Given a function $f_\theta : \mathcal{X} \times \mathcal{A} \rightarrow \mathbb{R}$ (e.g., a DNN or decision tree parameterized by θ), the density of choosing action a at state x is

$$\pi_\theta(a|x) = \frac{\exp(f_\theta(x, a))}{\int \exp(f_\theta(x, a')) da'}$$

Policy Parametrization: Examples

- A special case would be when $f_{\theta}(x, a) = \phi(x, a)^{\top} \theta$ for some features $\phi : \mathcal{X} \times \mathcal{A} \rightarrow \mathbb{R}^p$ and $\theta \in \mathbb{R}^p$:

$$\pi_{\theta}(a|x) = \frac{\exp(\phi(x, a)^{\top} \theta)}{\int \exp(\phi(x, a')^{\top} \theta) da'}.$$

- When the action space \mathcal{A} is discrete, $\pi_{\theta}(a|x)$ denotes the probability of choosing action a at state x (instead of its density):

$$\pi_{\theta}(a|x) = \frac{\exp(\phi(x, a)^{\top} \theta)}{\sum_{a' \in \mathcal{A}} \exp(\phi(x, a')^{\top} \theta)}.$$

Policy Parametrization: Examples

- Another example: $\pi_\theta(\cdot|x)$ defining a Normal distribution over action space with θ parameterization its mean and covariance:

$$\pi_\theta(\cdot|x) = \mathcal{N}(\mu_\theta(x), \Sigma_\theta(x)).$$

- If the action space is $d_{\mathcal{A}}$ -dimensional:
 - Mean: $\mu_\theta : \mathcal{X} \rightarrow \mathbb{R}^{d_{\mathcal{A}}}$
 - Covariance: $\Sigma_\theta : \mathcal{X} \rightarrow S_+^{d_{\mathcal{A}}}$. Here $S_+^{d_{\mathcal{A}}}$ refers to the set of $d_{\mathcal{A}} \times d_{\mathcal{A}}$ positive semi-definite matrices.

Ease of Work in Continuous Action Spaces

- Explicit parameterization of policy allows us to easily choose a continuous action
- For value-based methods, this can be challenging:
 - Even if we know Q^* , computing the optimal policy $\pi^*(x) = \pi_g(x; Q^*)$ requires an optimization problem in \mathcal{A} .
 - This is challenging if \mathcal{A} is a high-dimensional space.
 - VI and PI requires repeated calculation of the greedy policy.
- Sure, action selection might be easy!
- **Question:** How can we optimize the performance of a parametrized policy?

Performance Measure

- The performance can be measured in various ways.
- We focus on the expected return of following π_θ , the value function.
 - We can also incorporate the variance or some other risk measures, relatively easily.
- **Goal:** Find a policy that maximizes this performance measure.
- We are restricted to choosing policies within Π_Θ (1).

Performance Measure on a Single State

- Assume that we only care about the performance at state $x \in \mathcal{X}$.
- The goal of policy search:

$$\operatorname{argmax}_{\pi \in \Pi_{\Theta}} V^{\pi}(x) = \operatorname{argmax}_{\theta \in \Theta} V^{\pi_{\theta}}(x). \quad (2)$$

- **Interpretation:** We are interested in finding a policy that if the agent starts at this particular state x , its performance, measured according to its expected return, is maximized.

Performance Measure on a Single State

$$\operatorname{argmax}_{\pi \in \Pi_{\Theta}} V^{\pi}(x) = \operatorname{argmax}_{\theta \in \Theta} V^{\pi_{\theta}}(x).$$

- The optimal policy π^* not only maximizes the value function at this particular x , but also at any other $x' \in \mathcal{X}$.
- But the optimal policy may not be in Π_{Θ} .
- If $\pi^* \notin \Pi_{\Theta}$, we will not be able to find a policy that maximizes the value at all states.
 - In that case, we may want to find a policy that is only good at our starting state x , and ignore the performance at other states.
 - The obtained policy is going to be initial-state-dependent. If we change x to another state $x' \neq x$, the optimal policy within Π_{Θ} might change.

Performance Measure with $X_1 \sim \rho$

- Instead of the extreme case of considering a single initial state x , we can consider when the initial state is distributed according to some distribution $\rho \in \mathcal{M}(\mathcal{X})$.
- The performance measure would be the average of following π_θ with the initial state $X_1 \sim \rho$.

$$J(\pi_\theta) = J_\rho(\pi_\theta) \triangleq \int V^{\pi_\theta}(x) d\rho(x) = \mathbb{E}_{X \sim \rho} [V^{\pi_\theta}(X)]. \quad (3)$$

Performance Measure with $X_1 \sim \rho$

$$J(\pi_\theta) = J_\rho(\pi_\theta) \triangleq \int V^{\pi_\theta}(x) d\rho(x) = \mathbb{E}_{X \sim \rho} [V^{\pi_\theta}(X)].$$

- The optimal policy maximizes J_ρ .
- $J_\rho(\pi^*) \geq J_\rho(\pi_\theta)$ for any $\pi_\theta \in \Pi_\Theta$.
- If $\pi^* \notin \Pi_\Theta$, the inequality is strict.

Performance Measure with $X_1 \sim \rho$

$$J(\pi_\theta) = J_\rho(\pi_\theta) \triangleq \int V^{\pi_\theta}(x) d\rho(x) = \mathbb{E}_{X \sim \rho} [V^{\pi_\theta}(X)].$$

- In policy search methods, we aim to find the maximizer of the performance measure within Π_Θ .

$$\bar{\pi} \leftarrow \operatorname{argmax}_{\pi_\theta \in \Pi_\Theta} J_\rho(\pi_\theta). \quad (4)$$

- The corresponding policy is denoted by $\bar{\theta}$, i.e., $\bar{\pi} = \pi_{\bar{\theta}}$.
- For different ρ , we may get different optimizers.
- To emphasize the dependence of the maximizer on ρ , we may use $\bar{\pi}_\rho$.
- We may sometimes denote $J(\pi_\theta)$ or $J_\rho(\pi_\theta)$ simply by $J_\rho(\theta)$.

Policy Search as an Optimization Problem

- **Question:** How can we solve the optimization problem (4) to find π_θ that maximizes the performance measure J_ρ ?
- This is an optimization problem, so we can benefit from the arsenal of optimization algorithms.
- Being an RL problem, however, brings both challenges and opportunities.
 - **Challenge:** The value of J_ρ is not readily available, and has to be estimated through interaction with the environment.
 - **Opportunity:** The special structure of the RL problem, such as the value function satisfying the Bellman equation.

Policy Search as an Optimization Problem

- Optimization methods, broadly speaking, can be categorized based on the information they need about their objective.
- **Zero-order** methods only use the value of the objective at various query points.
 - They compute $J_\rho(\theta)$ at various θ s in order to guide the optimization process.
- **First-order** methods use the derivative of the objective instead of, or in addition to, the value of the objective.
 - They use $\nabla_\theta J_\rho(\theta)$ in order to guide the search.
 - The quintessential first-order optimization method is the gradient descent (and its stochastic variant).

Zero-Order Methods

- We first consider the case when the policy parameter space Θ is finite.
- This helps us understand some of the challenges.
- We then extend our discussion to case when Θ is continuous function space.

Zero-Order Methods: Finite Policy Parameter Space

- Assume that we are given a finite $\Theta = \{\theta_1, \dots, \theta_m\}$ policy parameters.
- This defines the finite policy space $\Pi_\Theta = \{\pi_\theta : \theta \in \Theta\}$.
- Find the policy $\pi_\theta \in \Pi_\Theta$ such that $J_\rho(\pi_\theta)$ is maximized (4).
- If we can easily compute $J_\rho(\pi_\theta)$ for each $\theta \in \Theta$, this is an easy problem, at least in principle.
- So the main issue is to compute $J_\rho(\pi_\theta)$.

Zero-Order Methods: Finite Policy Parameter Space

- The performance measure $J_\rho(\pi_\theta) = \mathbb{E}_{X \sim \rho} [V^{\pi_\theta}(X)]$, i.e., the expectation of $V^{\pi_\theta}(X)$ w.r.t. $X \sim \rho$.
- We can try to compute $V^{\pi_\theta}(x)$ for all $x \in \mathcal{X}$, using any of the PE methods that we have developed, and take the weighted average according to ρ .
- If \mathcal{X} is discrete, this would be

$$J_\rho(\pi_\theta) = \sum_{x \in \mathcal{X}} \rho(x) V^{\pi_\theta}(x).$$

- If \mathcal{X} is large:
 - Computing V^{π_θ} itself is not going to be easy.
 - Computing the integral $\int V^{\pi_\theta}(x) d\rho(x)$ is going to be challenging.

Zero-Order Methods: Finite Policy Parameter Space

- Alternative: Computing an unbiased estimate of $J_\rho(\pi_\theta)$ instead, using MC estimation.
- We derive this in two steps.
 - Assume that we know V^{π_θ} , estimate $J_\rho(\pi_\theta)$.
 - Replace $V^{\pi_\theta}(x)$ with the return $G^{\pi_\theta}(x)$.

Zero-Order Methods: Finite Policy Parameter Space

- We assume that we know V^{π_θ} , and we want to estimate $J_\rho(\pi_\theta)$.
- If we sample $X \sim \rho$, we have that $V^{\pi_\theta}(X)$ is an unbiased estimate of $J_\rho(\rho_\pi)$ as

$$\mathbb{E}[V^{\pi_\theta}] = \int V^{\pi_\theta}(x) d\rho(x) = J_\rho(\pi_\theta).$$

- If we draw n independent samples $X_1, \dots, X_n \sim \rho$, the estimator

$$\frac{1}{n} \sum_{i=1}^n V^{\pi_\theta}(X_i)$$

would be an unbiased as well.

Zero-Order Methods: Finite Policy Parameter Space

$$\frac{1}{n} \sum_{i=1}^n V^{\pi_{\theta}}(X_i)$$

- Variance:

$$\frac{\text{Var} [V^{\pi_{\theta}}(X)]}{n}.$$

- This variance goes to 0 as n increases.
- The variance $\text{Var} [V^{\pi_{\theta}}(X)]$ is a measure of dispersion of the value function across states samples according to ρ .
 - If the value function is constant, it will be zero.
 - If it is changing slowly as a function of the state, it would be small.
 - If the value function varies greatly, the variance is large.
- The variance is a function of the policy π_{θ} , so for each $\pi_{\theta} \in \Pi_{\theta}$, we get a different variance.

Zero-Order Methods: Finite Policy Parameter Space

- The second step is to replace $V^{\pi_\theta}(x)$ with the return $G^{\pi_\theta}(x)$.
- The return $G^{\pi_\theta}(x)$ is an unbiased estimate of $V^{\pi_\theta}(x)$.
- Computation of $G^{\pi_\theta}(x)$ requires starting the agent from state x and following π_θ (i.e., performing a rollout from x) until the end of episode for episodic tasks, or until infinity for continual tasks.
- If $X \sim \rho$, $G^{\pi_\theta}(X)$ is an unbiased estimate of $J_\rho(\pi_\theta)$ as

$$\begin{aligned}\mathbb{E}_{X \sim \rho} [G^{\pi_\theta}(X)] &= \mathbb{E}_{X \sim \rho} [\mathbb{E} [G^{\pi_\theta}(X) \mid X]] \\ &= \mathbb{E}_{X \sim \rho} [V^{\pi_\theta}(X)] = J_\rho(\pi_\theta).\end{aligned}$$

Zero-Order Methods: Finite Policy Parameter Space

- If we draw n independently selected $X_1, \dots, X_n \sim \rho$, we can form

$$\hat{J}_n(\pi_\theta) = \frac{1}{n} \sum_{i=1}^n G^{\pi_\theta}(X_i), \quad (5)$$

as an unbiased estimate of $J_\rho(\pi_\theta)$.

Zero-Order Methods: Finite Policy Parameter Space

Proposition

The estimator $\hat{J}_n(\pi_\theta)$ (5) is an unbiased estimator for $J_\rho(\pi_\theta)$ and has the variance of

$$\text{Var} \left[\hat{J}_n(\pi_\theta) \right] = \frac{1}{n} (\mathbb{E} [\text{Var} [G^{\pi_\theta}(X) \mid X]] + \text{Var} [V^{\pi_\theta}(X)]).$$

- If we have a finite number of parameters in Θ , we can estimate $J_\rho(\pi_{\theta_i}) \approx \hat{J}_n(\pi_{\theta_i}) \pm O_P(\frac{1}{\sqrt{n}})$ for each $\theta_i \in \Theta$.
- We can use these estimates to select the best among them:

$$\hat{\pi} = \pi_{\hat{\theta}} \leftarrow \underset{\theta \in \Theta}{\text{argmax}} \hat{J}_n(\pi_\theta). \quad (6)$$

Zero-Order Methods: Finite Policy Parameter Space

- As there is an $O_P(\frac{1}{\sqrt{n}})$ error in estimation of each $J_\rho(\pi_\theta)$, the selected policy $\hat{\pi}$ may not be the same as the maximizer $\bar{\pi}$ of (4).
- The error can happen if $\hat{J}_n(\hat{\pi}) > \hat{J}_n(\bar{\pi})$ (which leads to preferring $\hat{\pi}$ to $\bar{\pi}$ according to the empirical performance measure) even though $J_\rho(\hat{\pi}) < J_\rho(\bar{\pi})$.

Zero-Order Methods: Finite Policy Parameter Space

- Even if we make an error in selecting the best policy, the gap in their performance is within $O_P(\frac{1}{\sqrt{n}})$.
- As we increase n , the error in estimating $J_\rho(\pi_\theta)$ decreases and the probability of selecting an optimal policy increases.
- This increased accuracy, however, is at the cost of increased sample and computational complexity, which would be $n|\Theta|$ rollouts.

Zero-Order Methods: Finite Policy Parameter Space

Proposition

Consider $\hat{\pi} = \pi_{\hat{\theta}} \leftarrow \operatorname{argmax}_{\theta \in \Theta} \hat{J}_n(\pi_{\theta})$ (6). Assume that the returns $G^{\pi_{\theta}}(x)$ are all Q_{max} -bounded for any $\theta \in \Theta$ and $x \in \mathcal{X}$. Furthermore, suppose that $|\Theta| < \infty$. For any $\delta > 0$, we have that

$$J_{\rho}(\hat{\theta}) \geq \max_{\theta \in \Theta} J_{\rho}(\theta) - 2Q_{max} \sqrt{\frac{2 \ln \left(\frac{2|\Theta|}{\delta} \right)}{n}},$$

with probability at least $1 - \delta$.

Zero-Order Methods: Random Search

- If Θ is not finite, we cannot evaluate $\hat{J}_n(\pi_\theta)$ for all $\theta \in \Theta$.
- There are several generic methods for searching in a large parameter space:
 - Random Search (RS)
 - Simulated Annealing
 - Various evolutionary algorithms

Zero-Order Methods: Random Search

- We randomly pick m policy parameters $\theta_1, \dots, \theta_m \in \Theta$.
- Evaluate $\hat{J}_n(\pi_{\theta_i})$
- Pick the one with the highest value.
- Intuition of why this works:
 - With large enough m , one of θ_i might hit close to the optimal

$$\hat{\theta} \leftarrow \operatorname{argmax}_{\theta \in \Theta} \hat{J}_n(\pi_{\theta}).$$

- If n is large enough, the difference between $\hat{J}_n(\theta)$ and $J_{\rho}(\theta)$ would be small for all randomly selected θ .

Zero-Order Methods: Random Search

Require: Distribution $\nu \in \mathcal{M}(\Theta)$; Number of rollouts n ;

Maximum number of iterations K

- 1: Draw a parameter $\theta_1 = \theta'_1 \sim \nu$
- 2: Evaluate $\hat{J}_n(\pi_{\theta_1})$
- 3: **for** $k = 2, 3, \dots, K$ **do**
- 4: Draw a parameter $\theta'_k \sim \nu$
- 5: Evaluate $\hat{J}_n(\pi_{\theta'_k})$
- 6: **if** $\hat{J}_n(\pi_{\theta'_k}) > \hat{J}_n(\pi_{\theta_k})$ **then**
- 7: $\theta_k \leftarrow \theta'_k$
- 8: **else**
- 9: $\theta_k \leftarrow \theta_{k-1}$
- 10: **end if**
- 11: **end for**
- 12: **return** π_{θ_K}

Zero-Order Methods: Random Search

- We can provide guarantee that RS finds the optimal point, asymptotically.
- RS is not the most efficient way to search a parameter space.
- The way it is presented here does not benefit from all previous evaluation of the function when suggesting a new θ'_k .
- That knowledge can be useful by helping us focus on more promising regions of the search space, instead of blindly sampling from the same distribution ν .
- This can be achieved by adaptively changing the distribution ν_k to be a function of previous evaluations.

Evolutionary Algorithms

- A large class of optimization methods are inspired by the process of evolution.
- Heritable characteristics of individuals in a population change over generations due to processes such as natural selection.
- The evolution leads to the adaptation of individuals, which means that they become better to live in their habitat.

Evolutionary Process for Optimization

- Identifying a solution to an optimization problem as an individual in a population
- The value of the function to be optimized for a particular solution as the fitness of that individual
- Emulate the evolution:
 - Mutation
 - Reproduction (cross-over)
 - Selection
- There are many variations in how we can do this:
 - Genetic Algorithms
 - Genetic Programming
 - Evolutionary Strategy

Evolutionary Strategy (ES) (1 + 1)

Require: Initial point $\theta_0 \in \Theta$; Rollouts n ; Iterations K

Require: Initial standard deviation of mutation operator: $\sigma_1 > 0$

Require: Adaptation parameters: $c_+ > 0$ and $c_- < 0$.

- 1: Evaluate $\hat{J}_n(\pi_{\theta_0})$
- 2: **for** $k = 1, 2, \dots, K$ **do**
- 3: Draw a perturbation $\eta \sim \mathcal{N}(0, \mathbf{I})$
- 4: $\theta'_k \leftarrow \theta_k + \sigma_k \eta$ ▷ Mutation
- 5: Evaluate $\hat{J}_n(\pi_{\theta'_k})$
- 6: **if** $\hat{J}_n(\pi_{\theta'_k}) > \hat{J}_n(\pi_{\theta_k})$ **then** ▷ Selection
- 7: $\theta_{k+1} \leftarrow \theta'_k$
- 8: $\sigma_{k+1} \leftarrow \sigma_k e^{c_+}$
- 9: **else**
- 10: $\theta_{k+1} \leftarrow \theta_{k-1}$
- 11: $\sigma_{k+1} \leftarrow \sigma_k e^{c_-}$
- 12: **end if**
- 13: **end for**
- 14: **return** π_{θ_K}

Evolutionary Strategy (ES) (1 + 1) and Beyond

- Evolutionary Strategy (ES) (1 + 1) is one of the simplest evolutionary algorithms
 - Similar to RS, but guided choice of randomness
 - Has some theoretical analysis
- A modification of this algorithm is called ES(1, λ) with $\lambda > 1$ being an integer number.
 - The parent θ_k generates λ offsprings:

$$\theta'_{k,j} = \theta_k + \sigma_k \eta_j, \quad j = 1, \dots, \lambda.$$

- The competition would only be between the offsprings $\{\theta'_{k,j}\}_{j=1}^{\lambda}$, and not with the parent. Only one of the λ offsprings gets to the next generation.

Beyond Evolutionary Strategy

- ES does not have any sexual reproduction.
- There are other evolutionary algorithms that have the reproduction component too, e.g., Genetic Algorithm (GA).
- Evolutionary algorithms can be quite complicated algorithms.
 - Many heuristics, inspired by nature.
 - Their performance is often evaluated only empirically.
 - Analyzing them theoretically can be quite complicated.
 - Current available results are limited to simple algorithms, such as $ES(1 + 1)$, which may not be the best performing ones in practice.

Evolutionary Algorithms and RL

- Studying evolutionary algorithms to solve RL problems is a niche area in the RL community.
- Sometimes (often?), they are not considered as a part of the RL research proper.
- Knowing about them is useful!
- Both evolution and learning have been crucial adaptation mechanisms to get to the point where we have relatively smart species.
- Building AI agents with comparable capabilities to animals may require borrowing ideas from both learning and evolution.
 - Learning: Within the lifespan of the agent
 - Evolution: Across generations of the agents

First-Order Methods and the Policy Gradient

- The gradient of $J_\rho(\pi_\theta)$ w.r.t. θ allows us to design first-order optimization methods.
- Potentially more efficient in finding an optimum of the performance compared to zero-order methods.
- Not obvious how to compute the gradient:
 - The performance $J_\rho(\pi_\theta)$ depends on V^{π_θ} .
 - Not a simple function of π_θ .
 - The value function is a complicated function of the policy, reward distribution \mathcal{R} and the transition dynamics \mathcal{P} .

Finite Difference Approximation of the Policy Gradient

- Use **Finite Difference (FD)** approximation of the policy gradient.
- Can be computed using the value of the performance objective itself.
- Recall that given a function $f : \mathbb{R} \rightarrow \mathbb{R}$, the FD approximation of the derivative $f'(x) = \frac{df}{dx}(x)$ is

$$f'_{\text{FD}}(x) = \frac{f(x + \Delta x) - f(x)}{\Delta x}, \quad (7)$$

where Δx is a small number. This is called the **forward difference approximation**.

Finite Difference Approximation of the Policy Gradient

- By the Taylor's theorem, assuming twice differentiability, we have

$$f(x + \Delta x) = f(x) + f'(x)\Delta x + f''(z)|_{x < z < x + \Delta x} \frac{(\Delta x)^2}{2!}.$$

Therefore,

$$f'(x) = \frac{f(x + \Delta x) - f(x)}{\Delta x} - f''(z)|_{x < z < x + \Delta x} \frac{(\Delta x)^2}{2!}.$$

- The error between the FD approximation (7) and $f'(x)$ is

$$\left| f''(z) \Big|_{x < z < x + \Delta x} \frac{(\Delta x)^2}{2!} \right|,$$

that is, $O((\Delta x)^2)$.

Finite Difference Approximation of the Policy Gradient

- Central difference approximation:

$$f'_{\text{FD}}(x) = \frac{f(x + \Delta x) - f(x - \Delta x)}{2\Delta x}.$$

- Error is $O((\Delta x)^3)$.

Finite Difference Approximation of the Policy Gradient

To compute the gradient of $J_\rho(\pi_\theta)$ w.r.t. $\theta \in \mathbb{R}^p$, we need to compute $2p$ evaluations of J_ρ :

$$\nabla_\theta J_\rho(\pi_\theta) \approx \nabla_\theta^{(\text{FD})} J_\rho(\pi_\theta) = \begin{bmatrix} \frac{J_\rho(\theta + \varepsilon e_1) - J_\rho(\theta - \varepsilon e_1)}{2\varepsilon} \\ \vdots \\ \frac{J_\rho(\theta + \varepsilon e_i) - J_\rho(\theta - \varepsilon e_i)}{2\varepsilon} \\ \vdots \\ \frac{J_\rho(\theta + \varepsilon e_p) - J_\rho(\theta - \varepsilon e_p)}{2\varepsilon} \end{bmatrix},$$

where e_i is a unit vector along dimension i of \mathbb{R}^p .

Finite Difference Approximation of the Policy Gradient

- We cannot directly compute $J_\rho(\pi_\theta)$.
- We can only compute $\hat{J}_n(\pi_\theta)$ using rollouts.
- Replace each J_ρ above with their corresponding \hat{J}_n .
- This requires $2pn$ rollouts in total.
- Given the approximated gradient, which has error caused by both the FD approximation and using \hat{J}_n instead of J_ρ , we may use the **gradient ascent** to move towards higher value of $J_\rho(\pi_\theta)$:

$$\theta_{k+1} \leftarrow \theta_k + \alpha_k \nabla_\theta^{(\text{FD})} \hat{J}_n(\pi_{\theta_k}). \quad (8)$$

- Even though this is a feasible approach, we can compute the gradient more elegantly.

Policy Gradient for the Immediate Reward Problem

- Suppose that we want to find a policy $\pi_\theta : \mathcal{X} \rightarrow \mathcal{M}(\mathcal{A})$ with $\theta \in \mathbb{R}^p$ that maximizes the performance for the immediate reward problem.
- Recall that the interaction protocol is
 - At episode t , $X_t \sim \rho \sim \mathcal{M}(\mathcal{X})$
 - The agent chooses action $A_t \sim \pi_\theta(\cdot | X_t)$
 - The agent receives reward $R_t \sim \mathcal{R}(\cdot | X_t, A_t)$.
 - The agent starts the new (independent) episode $t + 1$.
- This is an **RL setting** as ρ and \mathcal{R} are not directly available to the agent, but only through samples.

Performance Measure

- The performance measure is

$$\begin{aligned} J_\rho(\pi_\theta) &= \int V^{\pi_\theta}(x) d\rho(x) = \int r^{\pi_\theta}(x) d\rho(x) \\ &= \int r(x, a) \pi_\theta(a|x) d\rho(x) da, \end{aligned}$$

as the value function V^{π_θ} for the immediate reward problem is the same as r^{π_θ} .

- Considered the action space to be continuous and we assume that $\pi_\theta(\cdot|x)$ provides a density over the state space.
- If we had a discrete action space, we would have

$$\int_{\mathcal{X}} d\rho(x) \sum_{a \in \mathcal{A}} r(x, a) \pi_\theta(a|x).$$

- We may switch back and forth between continuous and discrete action spaces.

Policy Gradient

- The gradient of $J_\rho(\pi_\theta)$ w.r.t. θ :

$$\begin{aligned}\nabla_\theta J_\rho(\pi_\theta) &= \int r(x, a) \nabla_\theta \pi_\theta(a|x) d\rho(x) da \\ &= \int d\rho(x) \int r(x, a) \nabla_\theta \pi_\theta(a|x) da \\ &= \mathbb{E}_{X \sim \rho} \left[\int r(X, a) \nabla_\theta \pi_\theta(a|X) da \right].\end{aligned}\quad (9)$$

- For discrete action spaces, the inner integral becomes $\sum_{a \in \mathcal{A}} r(x, a) \nabla_\theta \pi_\theta(a|x)$.
- We call $\nabla_\theta J_\rho(\pi_\theta)$ the **Policy Gradient** (PG).

Improving Performance Measure using Policy Gradient

If we can compute PG, we can update the policy parameters, using a gradient ascent method:

$$\theta_{k+1} \leftarrow \theta_k + \alpha_k \nabla_{\theta} J_{\rho}(\pi_{\theta_k}), \quad (10)$$

similar to what we have done using the FD approximation (8).

Computing the Policy Gradient

$$\nabla_{\theta} J_{\rho}(\pi_{\theta}) = \int r(x, a) \nabla_{\theta} \pi_{\theta}(a|x) d\rho(x) da$$

- How can we compute this gradient?
- We build this gradually in several steps.
- At each step, we relax some assumptions until we get to a procedure that can use the data available by the interaction protocol above.

Computing the Policy Gradient – Known ρ and r

$$\nabla_{\theta} J_{\rho}(\pi_{\theta}) = \int r(x, a) \nabla_{\theta} \pi_{\theta}(a|x) d\rho(x) da$$

- If we know ρ and r , we have all the necessary information for computing the gradient.
 - For each $x \in \mathcal{X}$, we compute the summation (or integral) over all $a \in \mathcal{A}$ of $r(x, a) \nabla_{\theta} \pi_{\theta}(a|x)$.
 - We weigh that term proportional to $\rho(x)$.
 - Take average over all x .
- But this is not the RL setting described as the interaction protocol at the beginning of the section.

Computing the Policy Gradient – Known r , unknown ρ

- Assume that r is known, but ρ can only be sampled.
- Approximately solve this problem by sampling $X_i \sim \rho$ ($i = 1, \dots, n$) and computing

$$\frac{1}{n} \sum_{i=1}^n \sum_{a \in \mathcal{A}} r(X_i, a) \nabla_{\theta} \pi_{\theta}(a|X_i). \quad (11)$$

or

$$\frac{1}{n} \sum_{i=1}^n \int r(X_i, a) \nabla_{\theta} \pi_{\theta}(a|X_i) da.$$

As $X_i \sim \rho$, this is an unbiased estimate of

$$\nabla_{\theta} J_{\rho}(\pi_{\theta}) = \mathbb{E}_{X \sim \rho} \left[\sum_{a \in \mathcal{A}} r(X, a) \nabla_{\theta} \pi_{\theta}(a|X) \right]$$

or $\mathbb{E}_{X \sim \rho} \left[\int r(x, a) \nabla_{\theta} \pi_{\theta}(a|x) da \right]$ (continuous).

Computing the Policy Gradient – Known r , unknown ρ

- As $X_i \sim \rho$, this is an unbiased estimate of

$$\nabla_{\theta} J_{\rho}(\pi_{\theta}) = \mathbb{E}_{X \sim \rho} \left[\sum_{a \in \mathcal{A}} r(X, a) \nabla_{\theta} \pi_{\theta}(a|X) \right]$$

or

$$\mathbb{E}_{X \sim \rho} \left[\int r(x, a) \nabla_{\theta} \pi_{\theta}(a|x) da \right].$$

- This is still not feasible if r is unknown in our interaction protocol:
 - the agent is initialized at state x
 - it has to choose its action according to $A \sim \pi_{\theta}(\cdot|x)$.

Computing the Policy Gradient – Unknown r and ρ

- The term

$$\sum_{a \in \mathcal{A}} r(x, a) \nabla_{\theta} \pi_{\theta}(a|x)$$

can be interpreted as the expectation of

$$r(x, A) \nabla_{\theta} \pi_{\theta}(A|x)$$

when A is coming from a uniform distribution with $q(a) = \frac{1}{|\mathcal{A}|}$ (for $a \in \mathcal{A}$).

- We have

$$\sum_{a \in \mathcal{A}} r(x, a) \nabla_{\theta} \pi_{\theta}(a|x) = |\mathcal{A}| \sum_{a \in \mathcal{A}} q(a) r(x, a) \nabla_{\theta} \pi_{\theta}(a|x). \quad (12)$$

- Similar for the continuous case.

Computing the Policy Gradient – Unknown r and ρ

- If the actions were coming from a uniform distribution, we could easily form an empirical estimate of these terms.
- But the actions in the interaction protocol comes from distribution $\pi_{\theta}(\cdot|x)$, which in general is different distribution than a uniform one.
- We have some form of off-policy sampling scenario in the distribution of actions.
- Some approaches to deal with it:
 - Estimate $\hat{r} \approx r$ using data (model-based approach).
 - Modify $r(x, A)\nabla_{\theta}\pi_{\theta}(A|x)$ to a quantity that can be estimated from data.

Computing the Policy Gradient – Unknown r and ρ

- Observation: for a function $f : \mathbb{R} \rightarrow \mathbb{R}$, we have

$$\frac{d \log f(x)}{dx} = \frac{\frac{df}{dx}(x)}{f(x)},$$

or more generally, for a function $f : \mathbb{R}^p \rightarrow \mathbb{R}$,

$$\nabla_x \log f(x) = \frac{\nabla_x f(x)}{f(x)}.$$

Computing the Policy Gradient – Unknown r and ρ

- Using this observation, we get

$$\begin{aligned}\int r(x, a) \nabla_{\theta} \pi_{\theta}(a|x) da &= \int r(x, a) \pi_{\theta}(a|x) \nabla_{\theta} \log \pi_{\theta}(a|x) da \\ &= \mathbb{E}_{A \sim \pi_{\theta}(\cdot|x)} [r(x, A) \nabla_{\theta} \log \pi_{\theta}(A|x)].\end{aligned}$$

- The desired quantity can be written as the expectation of

$$r(x, A) \nabla_{\theta} \log \pi_{\theta}(A|x)$$

when $A \sim \pi_{\theta}(\cdot|x)$.

- The sampling distribution is the same as the one agent uses to choose its actions.
- We are in the on-policy sampling scenario over the choice of actions.

Computing the Policy Gradient – Unknown r and ρ

- If $X \sim \rho$ and $A \sim \pi_\theta(\cdot|X)$, the random variable

$$r(X, A)\nabla_\theta \log \pi_\theta(A|X) \quad (13)$$

is an unbiased estimate of $\nabla_\theta J_\rho(\pi_\theta)$, i.e.,

$$\begin{aligned} \nabla_\theta J_\rho(\pi_\theta) &= \mathbb{E} [r(X, A)\nabla_\theta \log \pi_\theta(A|X)] \\ &= \mathbb{E}_{X \sim \rho} [\mathbb{E}_{A \sim \pi_\theta(\cdot|X)} [r(X, A)\nabla_\theta \log \pi_\theta(A|X) | X]]. \end{aligned} \quad (14)$$

- We can estimate the gradient of the performance w.r.t. the parameters of the policy using data available through the interaction of the agent with its environment.
- We may use this estimate in (10) to update the policy parameters using unbiased but noisy estimate of the gradient.
- This makes it a stochastic gradient ascent.

Two Sources of Variance

$$r(X, A) \nabla_{\theta} \log \pi_{\theta}(A|X)$$

- Unbiased estimate of the gradient
- But it has variance due to two sources of randomness:
 - Variance of estimating

$$g(x; \theta) \triangleq \mathbb{E}_{A \sim \pi_{\theta}(\cdot|X)} [r(X, A) \nabla_{\theta} \log \pi_{\theta}(A|X) \mid X = x]$$

with a single sample $r(X, A) \nabla_{\theta} \log \pi_{\theta}(A|X)$.

- Variance of estimating $\mathbb{E}_{X \sim \rho} [g(X; \theta)]$ using a single sample.

Two Sources of Variance

$$r(X, A) \nabla_{\theta} \log \pi_{\theta}(A|X)$$

- One can show that the variance along the i -th dimension of this r.v. is

$$\begin{aligned} \text{Var} \left[r(X, A) \frac{\partial \log \pi_{\theta}(A|X)}{\partial \theta_i} \right] = \\ \mathbb{E}_{X \sim \rho} \left[\text{Var} \left[r(X, A) \frac{\partial \log \pi_{\theta}(A|X)}{\partial \theta_i} \mid X \right] \right] + \text{Var}_{X \sim \rho} [g_i(X; \theta)]. \end{aligned} \tag{15}$$

Two Sources of Variance

- Let us define

$$g(x; \theta) = \mathbb{E}_{A \sim \pi_\theta(\cdot|x)} [r(x, A) \nabla_\theta \log \pi_\theta(A|x)]. \quad (16)$$

- The function $g : \mathcal{X} \times \Theta \rightarrow \mathbb{R}^p$ is the gradient of r^{π_θ} w.r.t. θ at state x , and is a p -dimensional vector.
- If we knew $r(x, a)$ and we could compute $g(x; \theta)$, we wouldn't have the first source of variance, but we still would have the second one.
- The variance would be

$$\text{Var}_{X \sim \rho} [g_i(X; \theta)].$$

- These two sources of variance make our estimate of the gradient inaccurate.
- There are ways to reduce them.

Variance Reduction – Randomness of States

- Suppose we can compute $g(x; \theta)$ exactly for any given $x \in \mathcal{X}$.
- Instead of a single sample $g(X_1; \theta)$, we use multiple independent samples X_1, \dots, X_n , all distributed according to ρ , to estimate the PG:

$$\begin{aligned}\nabla_{\theta} J_{\rho}(\pi_{\theta}) &\approx \frac{1}{n} \sum_{i=1}^n g(X_i; \theta) \\ &= \frac{1}{n} \sum_{i=1}^n \mathbb{E}_{A \sim \pi_{\theta}(\cdot | X_i)} [r(X_i, A) \nabla_{\theta} \log \pi_{\theta}(A | X_i)].\end{aligned}$$

- The variance of this estimator, along dimension i , is

$$\frac{1}{n} \text{Var}_{X \sim \rho} [g_i(X; \theta)].$$

As $n \rightarrow \infty$, the variance goes to zero. This leads to more accurate estimate of the PG, hence more accurate update of the policy.

Variance Reduction – Randomness of Actions

Consider the variance of estimating $g(x; \theta)$ (16) using a single sample $r(x, A) \nabla_{\theta} \log \pi_{\theta}(A|x)$ with $A \sim \pi_{\theta}(\cdot|x)$.

For each dimension i , we have

$$\begin{aligned}\mathbb{E} \left[\frac{\partial \log \pi_{\theta}(A|x)}{\partial \theta_i} b(x) \mid x \right] &= \int \pi_{\theta}(a|x) \frac{\partial \log \pi_{\theta}(a|x)}{\partial \theta_i} b(x) da \\ &= \int \frac{\partial \pi_{\theta}(a|x)}{\partial \theta_i} b(x) da \\ &= b(x) \int \frac{\partial \pi_{\theta}(a|x)}{\partial \theta_i} da \\ &= b(x) \frac{\partial}{\partial \theta_i} \underbrace{\int \pi_{\theta}(a|x) da}_{=1} = 0.\end{aligned}$$

Variance Reduction – Randomness of Actions

- This shows that

$$\mathbb{E} \left[\frac{\partial \log \pi_{\theta}(A|x)}{\partial \theta_i} r(x, A) \mid x \right] = \mathbb{E} \left[\frac{\partial \log \pi_{\theta}(A|x)}{\partial \theta_i} (r(x, A) + b(x)) \mid x \right]. \quad (17)$$

- Adding a state-dependent function $b : \mathcal{X} \rightarrow \mathbb{R}$ to $r(x, a)$ does not change the expectation.
- But it may change the variance!

Variance Reduction – Randomness of Actions

- For each dimension i of the PG, we can use a different state-dependent function.
- For any state-dependent function $b : \mathcal{X} \rightarrow \mathbb{R}^p$, the PG (14) is

$$\begin{aligned}\nabla_{\theta} J_{\rho}(\pi_{\theta}) &= \mathbb{E} [r(X, A) \nabla_{\theta} \log \pi_{\theta}(A|X)] = \\ &\quad \mathbb{E} [(r(X, A)\mathbf{1} + b(X)) \odot \nabla_{\theta} \log \pi_{\theta}(A|X)],\end{aligned}$$

where $\mathbf{1}$ is a p -dimensional vector with all components equal to 1, and \odot is a pointwise (Hadamard) product of two vectors, i.e., for $u, v \in \mathbb{R}^p$, $[u \odot v]_i = u_i v_i$.

- If we simply choose a scalar function b , which is often the case in practice, we have

$$\nabla_{\theta} J_{\rho}(\pi_{\theta}) = \mathbb{E} [(r(X, A) + b(X)) \nabla_{\theta} \log \pi_{\theta}(A|X)].$$

- The function b is called the **baseline**.

Variance Reduction – Randomness of Actions – Baseline

- The baseline can be used in order to minimize the variation of p -dimensional random vector.
- We use the variance for this purpose.
- We would like to find a function $b : \mathcal{X} \rightarrow \mathbb{R}^p$ such that for all $x \in \mathcal{X}$,

$$\min_b \sum_{i=1}^p \text{Var} \left[(r(x, A) + b_i(x)) \frac{\partial \log \pi_\theta(A|x)}{\partial \theta_i} \mid x \right] =$$
$$\text{Tr Cov} ((r(X, A)\mathbf{1} + b(x)) \odot \nabla_\theta \log \pi_\theta(A|x) \mid x)$$

Variance Reduction – Randomness of Actions – Baseline

$$b_i(x) = \frac{-\mathbb{E} \left[r(x, A) \left(\frac{\partial \log \pi_\theta(A|x)}{\partial \theta_i} \right)^2 \mid x \right]}{\mathbb{E} \left[\left(\frac{\partial \log \pi_\theta(A|x)}{\partial \theta_i} \right)^2 \mid x \right]}. \quad (18)$$

We could choose a single scalar function $b : \mathcal{X} \rightarrow \mathbb{R}$ instead. In that case, the solution would be

$$b(x) = \frac{-\mathbb{E} \left[r(x, A) \|\nabla_\theta \log \pi_\theta(A|x)\|_2^2 \mid x \right]}{\mathbb{E} \left[\|\nabla_\theta \log \pi_\theta(A|x)\|_2^2 \mid x \right]}.$$

Policy Gradient for Continuing Tasks

- We derive the PG for continuing tasks.
- The difference with the immediate reward case is that the performance $J_\rho(\pi_\theta)$ depends on the dynamics \mathcal{P}^{π_θ} too.
- As we change θ , the dynamics \mathcal{P}^{π_θ} changes as well.
- This seems to complicate the gradient computation.
- It turns out that despite this challenge, the PG can be written in an elegant, and relatively easy to compute, form.

Discounted Future-State Distribution

- New notations to present the results more compactly.
- Recall that $\mathcal{P}^\pi(\cdot|x; k) = \mathcal{P}^\pi(\cdot|x)^k$ is the probability distribution of following policy π for $k \geq 0$ steps.
- We introduce **discounted future-state distribution** of starting from $x \in \mathcal{X}$ and following π as

$$\rho_\gamma^\pi(\cdot|x) = \rho_\gamma(\cdot|x; \mathcal{P}^\pi) \triangleq (1 - \gamma) \sum_{k \geq 0} \gamma^k \mathcal{P}^\pi(\cdot|x; k). \quad (19)$$

- It is easy to verify that $\rho_\gamma^\pi(\cdot|x)$ is a valid probability distribution, e.g., $\rho_\gamma^\pi(\mathcal{X}|x) = 1$.

Discounted Future-State Distribution

- The relevance of this distribution becomes more clear if we note that

$$\begin{aligned} V^\pi(x) &= \mathbb{E} \left[\sum_{t \geq 0} \gamma^t R_t \mid X_0 = x \right] \\ &= \sum_{t \geq 0} \gamma^t \mathbb{E} [R_t \mid X_0 = x] \\ &= \sum_{t \geq 0} \gamma^t \int \mathcal{P}^\pi(dx' \mid x; t) r(x') \\ &= \frac{1}{1 - \gamma} \int \rho_\gamma^\pi(dx' \mid x) r(x') = \frac{1}{1 - \gamma} \mathbb{E}_{X' \sim \rho_\gamma^\pi(\cdot \mid x)} [r(X')] . \end{aligned}$$

- The value function at a state x is the expected reward when X' is distributed according to $\rho_\gamma^\pi(\cdot \mid x)$.

Discounted Future-State Distribution

- Interpretation: The agent starts from state x and at each time step, it decides to follow π with probability γ or terminates the episode with probability $1 - \gamma$.
- We can also define discounted future-state distribution of starting from ρ and following π as

$$\rho_\gamma^\pi(\cdot) = \rho_\gamma(\cdot | \mathcal{P}^\pi) \triangleq \int \rho_\gamma(\cdot | x; \mathcal{P}^\pi) d\rho(x).$$

- The performance measure $J(\pi_\theta)$ (3) is

$$J(\pi_\theta) = \mathbb{E}_{X \sim \rho} [V^{\pi_\theta}(X)] = \frac{1}{1 - \gamma} \mathbb{E}_{X \sim \rho_\gamma^\pi} [r(X)].$$

Policy Gradient Theorem

Theorem (Policy Gradient Theorem – Sutton et al. 2000)

Assume that π_θ is differentiable w.r.t. $\theta \in \Theta$. We have

$$\begin{aligned} \nabla_\theta J_\rho(\pi_\theta) &= \\ & \sum_{k \geq 0} \gamma^k \int d\rho(x) \mathcal{P}^{\pi_\theta}(dx'|x; k) \int \nabla_\theta \pi_\theta(a'|x') Q^{\pi_\theta}(x', a') da' = \\ & \frac{1}{1 - \gamma} \int \rho_\gamma^{\pi_\theta}(dx) \int \pi_\theta(a|x) \nabla_\theta \pi_\theta(a|x) Q^{\pi_\theta}(x, a) da = \\ & \frac{1}{1 - \gamma} \mathbb{E} [\nabla_\theta \log \pi_\theta(A|X) Q^{\pi_\theta}(X, A)], \end{aligned}$$

with $X \sim \rho_\gamma^{\pi_\theta}$, $A \sim \pi_\theta(\cdot|X)$.

Policy Gradient Theorem – Proof

We write the value function at state $x \in \mathcal{X}$ as the expected value of the action-value function, i.e.,

$$V^{\pi_\theta}(x) = \int \pi_\theta(a|x) Q^{\pi_\theta}(x, a) da.$$

We take its derivative w.r.t. θ and use the product rule to get

$$\nabla_\theta V^{\pi_\theta}(x) = \int [\nabla_\theta \pi_\theta(a|x) Q^{\pi_\theta}(x, a) + \pi_\theta(a|x) \nabla_\theta Q^{\pi_\theta}(x, a)] da. \quad (20)$$

Policy Gradient Theorem – Proof

$$\text{As } Q^{\pi_\theta}(x, a) = r(x, a) + \gamma \int \mathcal{P}(dx'|x, a)V^{\pi_\theta}(x'),$$

$$\nabla_\theta Q^{\pi_\theta}(x, a) = \gamma \int \mathcal{P}(dx'|x, a)\nabla_\theta V^{\pi_\theta}(x').$$

This alongside with (20) gives us the recursive Bellman-like equation for the gradient of $V^{\pi_\theta}(x)$:

$$\nabla_\theta V^{\pi_\theta}(x) = \int \nabla_\theta \pi_\theta(a|x)Q^{\pi_\theta}(x, a)da + \gamma \int \mathcal{P}^{\pi_\theta}(dx'|x)\nabla_\theta V^{\pi_\theta}(x'). \quad (21)$$

Policy Gradient Theorem – Proof

Expanding $\nabla_{\theta} V^{\pi_{\theta}}(x')$ likewise, we get that

$$\begin{aligned} \nabla_{\theta} V^{\pi_{\theta}}(x) &= \int \nabla_{\theta} \pi_{\theta}(a|x) Q^{\pi_{\theta}}(x, a) da + \\ &\quad \gamma \int \mathcal{P}^{\pi_{\theta}}(dx'|x) \left[\nabla_{\theta} \pi_{\theta}(a'|x') Q^{\pi_{\theta}}(x', a') da' + \right. \\ &\quad \left. \gamma \int \mathcal{P}^{\pi_{\theta}}(dx''|x') \nabla_{\theta} V^{\pi_{\theta}}(x'') \right]. \end{aligned}$$

Following this pattern recursively, we get that

$$\begin{aligned} \nabla_{\theta} V^{\pi_{\theta}}(x) &= \sum_{k \geq 0} \gamma^k \int \mathcal{P}^{\pi_{\theta}}(dx'|x; k) \int \nabla_{\theta} \pi_{\theta}(a'|x') Q^{\pi_{\theta}}(x', a') da' \\ &= \frac{1}{1 - \gamma} \int \rho_{\gamma}^{\pi_{\theta}}(dx'|x) \int \nabla_{\theta} \pi_{\theta}(a'|x') Q^{\pi_{\theta}}(x', a') da'. \end{aligned}$$

Policy Gradient Theorem – Proof

Also since $\nabla_{\theta}\pi_{\theta}(a'|x') = \pi_{\theta}(a'|x')\nabla_{\theta}\log\pi_{\theta}(a'|x')$, we can write the gradient as

$$\begin{aligned}\nabla_{\theta}V^{\pi_{\theta}}(x) &= \\ \frac{1}{1-\gamma} \int \rho_{\gamma}^{\pi_{\theta}}(dx'|x) \int \pi_{\theta}(a'|x') \nabla_{\theta}\log\pi_{\theta}(a'|x') Q^{\pi_{\theta}}(x', a') da' &= \\ \frac{1}{1-\gamma} \int \rho_{\gamma}^{\pi_{\theta}}(dx'|x) \mathbb{E}_{A' \sim \pi_{\theta}(\cdot|X')} [\nabla_{\theta}\log\pi_{\theta}(A'|X') Q^{\pi_{\theta}}(X', A')] &.\end{aligned}$$

Policy Gradient Theorem – Proof

As $J_\rho(\pi_\theta) = \int V^{\pi_\theta}(x) d\rho(x)$, taking the average of x w.r.t. ρ , we get that

$$\begin{aligned}\nabla_\theta J_\rho(\pi_\theta) &= \frac{1}{1-\gamma} \int \rho_\gamma^{\pi_\theta}(dx) \int \pi_\theta(a|x) \nabla_\theta \pi_\theta(a|x) Q^{\pi_\theta}(x, a) da \\ &= \frac{1}{1-\gamma} \mathbb{E}_{\substack{X \sim \rho_\gamma^{\pi_\theta} \\ A \sim \pi_\theta(\cdot|X)}} [\nabla_\theta \log \pi_\theta(A|X) Q^{\pi_\theta}(X, A)],\end{aligned}$$

which is the desired result.

Policy Gradient Theorem

- This theorem provides an elegant formula for the PG.
- It relates the PG to the discounted future-state distribution $\rho_\gamma^{\pi_\theta}$, the action-value function $Q^{\pi_\theta}(x, a)$, and the gradient of π_θ .
- To compute the PG in the RL setting, we have to estimate it using samples. If we get
 - a state X sampled from $\rho_\gamma^{\pi_\theta}$,
 - an action A sampled from $\pi_\theta(\cdot|X)$, and
 - know action-value $Q^{\pi_\theta}(X, A)$,

the random variables

$$\nabla_\theta \log \pi_\theta(A|X) Q^{\pi_\theta}(X, A)$$

is an unbiased estimate of the PG (cf. (13)).

- We can then use it in an SGD scheme to improve the policy.

Sampling X from $\rho_\gamma^{\pi_\theta}$

Sampling from $\rho_\gamma^{\pi_\theta}$ is relatively straightforward in the on-policy sampling scenario when the agent follows π_θ .

- The agent starts an episode from $X_0 \sim \rho$ and follows π_θ .
- We get a sequence of states X_0, X_1, \dots .
- These would be samples from $\int d\rho(x) \mathcal{P}^{\pi_\theta}(\cdot|x; k)$ for $k = 0, 1, \dots$.
- The distribution $\rho_\gamma^{\pi_\theta}$, however, has a γ^k factor for the k -th time step, see (19).
- Its effect is that the contribution to the gradient from X_k , which is

$$\mathbb{E} [\nabla_\theta \log \pi_\theta(A|X) Q^{\pi_\theta}(X, A)] = \int \pi_\theta(a|x) \nabla_\theta \pi_\theta(a|x) Q^{\pi_\theta}(x, a) da,$$

should be weighted by γ^k .

Sampling X from $\rho_{\gamma}^{\pi_{\theta}}$

- Another way to directly sample from $\rho_{\gamma}^{\pi_{\theta}}$ is to follow π , but at each step terminate the episode with probability $1 - \gamma$.

Sampling A from $\pi_{\theta}(\cdot|X)$

An action A sampled from $\pi_{\theta}(\cdot|X)$ is automatically generated when the agent follows policy π_{θ} (on-policy).

Computation of $Q^{\pi_\theta}(X, A)$

- The remaining issue is the computation of $Q^{\pi_\theta}(X, A)$ for $X \sim \rho_\gamma^{\pi_\theta}$ and $A \sim \pi_\theta(\cdot|X)$ using data.
- This is essentially a PE problem, and we may use various action-value function estimators that we have developed so far.

Computation of $Q^{\pi_\theta}(X, A)$

- A simple approach: Use the MC estimate $Q^{\pi_\theta}(X, A)$. This would lead to what is known as the REINFORCE algorithm by Williams [1992].¹
- In the on-policy setting when the agent follows π_θ , it generates the sequence $X_0, A_0, R_0, X_1, A_1, R_1, \dots$ with $A_t \sim \pi_\theta(\cdot | X_t)$.
- The return $G_t^\pi = \sum_{k \geq t} \gamma^{k-t} R_k$ is an unbiased estimate of $Q^{\pi_\theta}(X_t, A_t)$.
- We replace the action-value function at that state-action with this return from time t onward.
- The return, however, is a high variance estimate of the action-value function.
- One approach is to use a baseline in order to reduce the variance of this MC estimate.

¹REINFORCE stands for REward Increment \times Nonnegative Factor \times Offset Reinforcement \times Characteristic Eligibility.

Computation of $Q^{\pi_{\theta}}(X, A)$ – Actor-Critic Methods

- Another approach is to use an action-value function estimator instead.
 - TD methods
 - LSTD
 - Fitted Value Iteration (for PE, and not for Control)
- Such a method is called **actor-critic** method
 - The actor refers to the policy (and often PG method to improve it)
 - The critic refers to the value function estimator used to criticize the policy (actor).

Computation of $Q^{\pi_\theta}(X, A)$ – Actor-Critic Methods

- The use of a critic, however, may induce a bias as $\mathbb{E} \left[\hat{Q}^{\pi_\theta}(X, A) | X, A \right]$ may be different from $Q^{\pi_\theta}(X, A)$, especially if we use a TD method (which introduces bias because of bootstrapping) or a function approximator (for large state-action spaces).
- Such a method would explicitly represent both policy and value function.
- Actor-critic methods bring together some of the benefits of both value-based and policy search methods.

Summary

- Policy Search Methods: Explicit representation of the policy and searching within the policy space
- The search might be guided by the zero-order or first-order methods
- We may sometimes constraint the change of the policy update
 - Policy gradient is only a local information. We should not make a large move.

References

- Richard S. Sutton, David McAllester, Satinder Singh, and Yishay Mansour. Policy gradient methods for reinforcement learning with function approximation. In *Advances in Neural Information Processing Systems (NIPS - 12)*, 2000.
- Ronald J. Williams. Simple statistical gradient-following algorithms for connectionist reinforcement learning. *Machine Learning*, 8 (3-4):229–256, 1992.