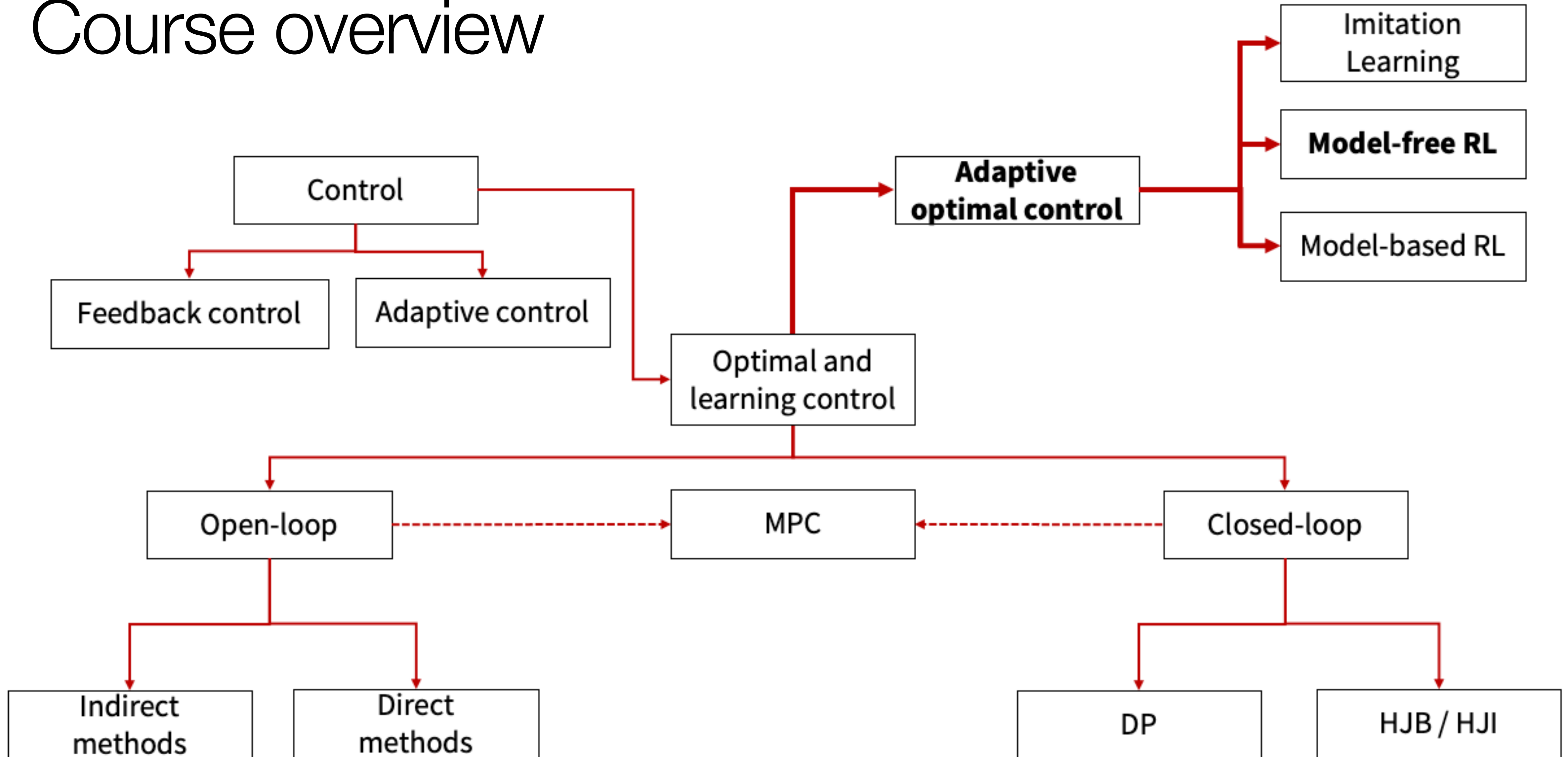


# AA203

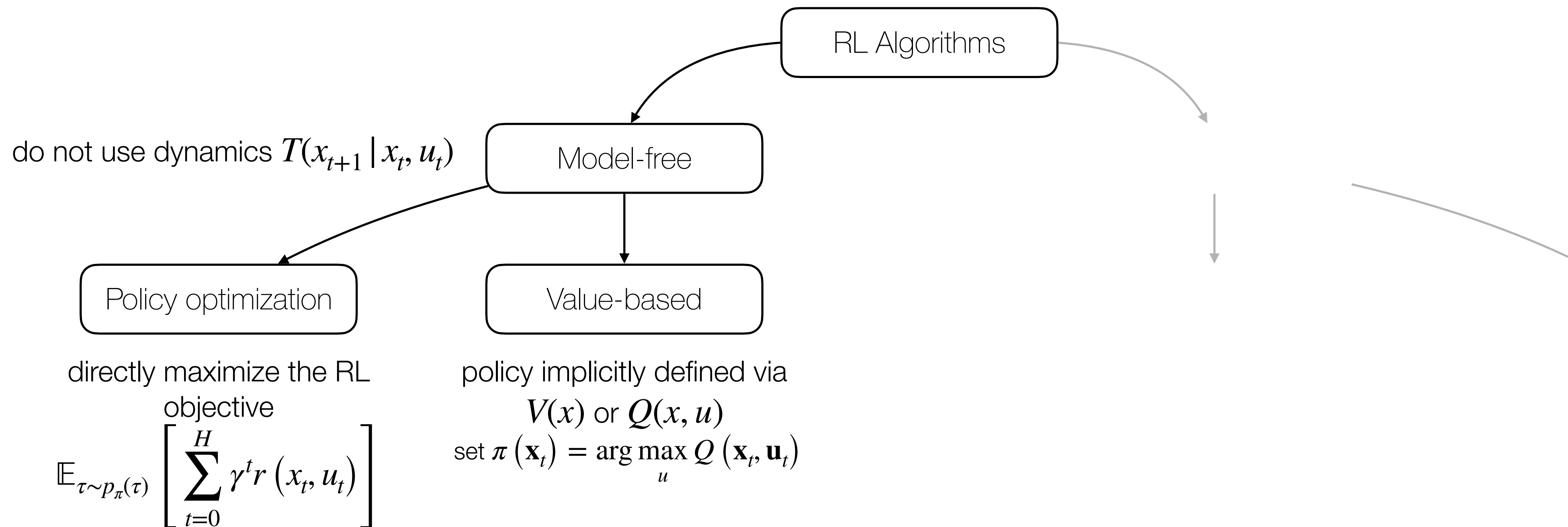
# Optimal and Learning-based Control

Model-free Reinforcement Learning: Policy Optimization

# Course overview



# A taxonomy of RL



# Outline

Intro to policy gradients

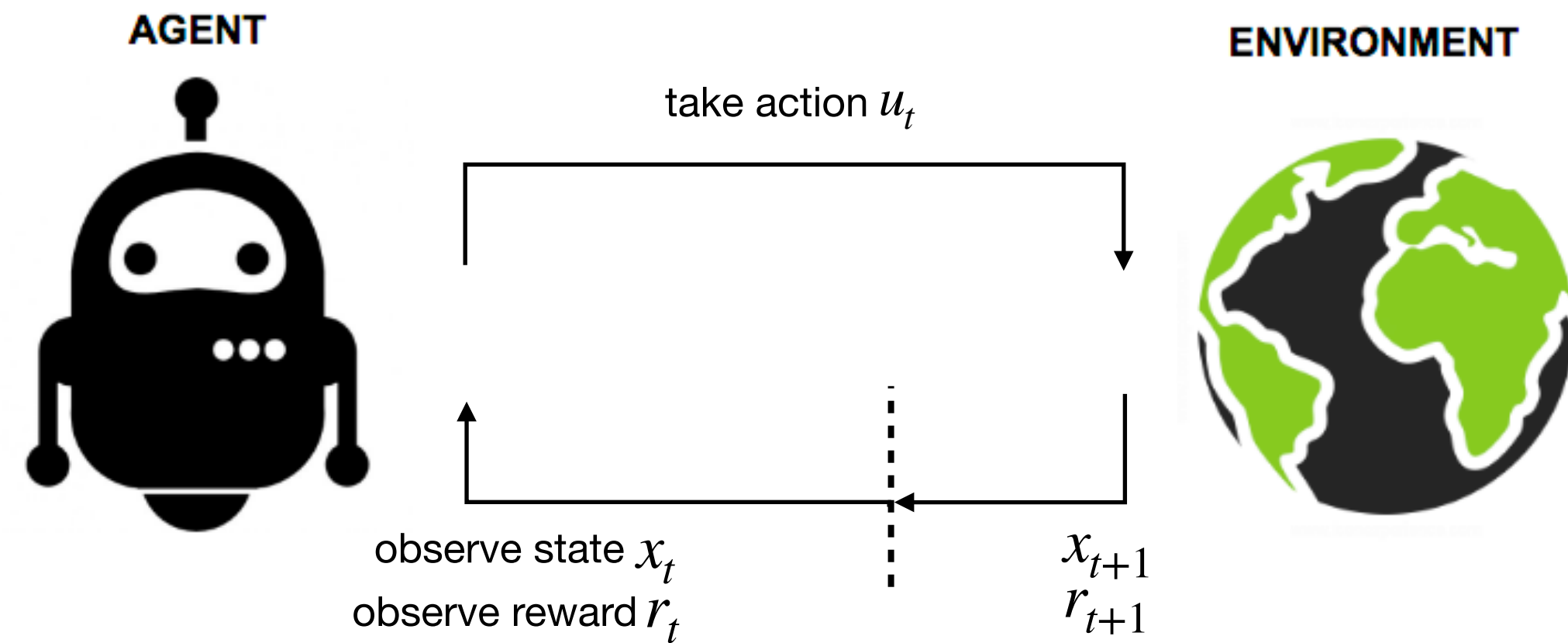
- REINFORCE algorithm
- Reducing variance of policy gradient

Actor-Critic methods

- Advantage
- Architecture design

Deep RL Algorithms & Applications

# The goal of reinforcement learning



- The agent interacts with the environment to generate trajectories  $\tau = (x_0, u_0, x_1, u_1, \dots, x_T)$
- We define the trajectory distribution

$$p(x_0, u_0, \dots, x_T) = p(\tau) = p(x_0) \prod_{t=1}^T \pi(u_t | x_t) p(x_{t+1} | x_t, u_t)$$

- We can express the RL objective as an expectation under the trajectory distribution

$$\pi^* = \arg \max_{\pi} \mathbb{E}_{\tau \sim p(\tau)} \left[ \sum_{t \geq 0} \gamma^t R(x_t, u_t) \right]$$

# Policy Optimization

- In policy optimization, we care about learning an (explicit) parametric policy  $\pi_\theta$ , with parameters  $\theta$
- In light of this, we can re-write the Eqs from the previous slide w.r.t.  $\theta$ :

$$p(x_0, u_0, \dots, x_T) = p(\tau) = p(x_0) \prod_{t=1}^T \pi_\theta(u_t | x_t) p(x_{t+1} | x_t, u_t)$$
$$\theta^* = \arg \max_{\theta} \underbrace{\mathbb{E}_{\tau \sim p(\tau)} \left[ \sum_{t \geq 0} \gamma^t R(x_t, u_t) \right]}_{J(\theta)}$$

To simplify the notation, we'll ignore discounting for now ( $\gamma = 1$ ) and consider

$$\theta^* = \arg \max_{\theta} \underbrace{\mathbb{E}_{\tau \sim p(\tau)} \left[ \sum_{t \geq 0} R(x_t, u_t) \right]}_{J(\theta)}$$

# Evaluating the objective

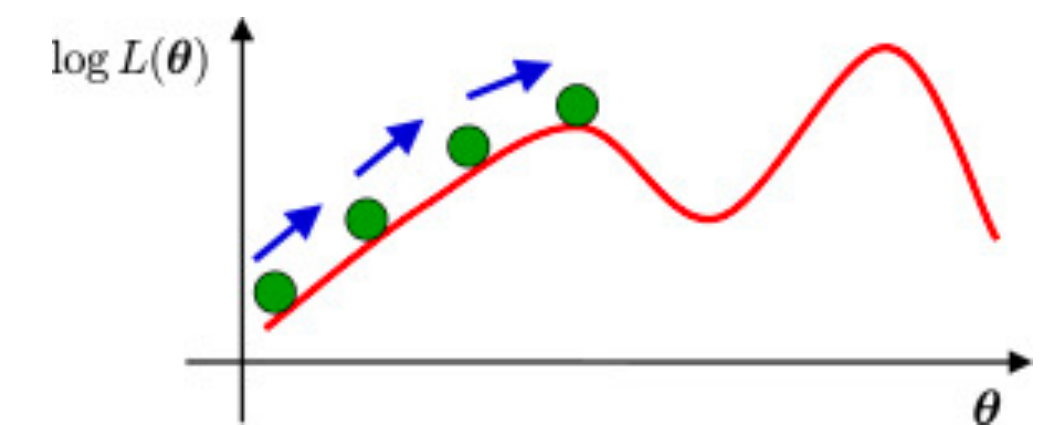
- Opposed to value-based methods, policy optimization attempts to learn the policy directly (i.e., optimize  $J(\theta)$  w.r.t.  $\theta$ )

$$J(\theta) = \mathbb{E}_{\tau \sim p(\tau)} \left[ \sum_{t \geq 0} R(x_t, u_t) \right]$$

One of the most direct ways to optimize this objective is to:

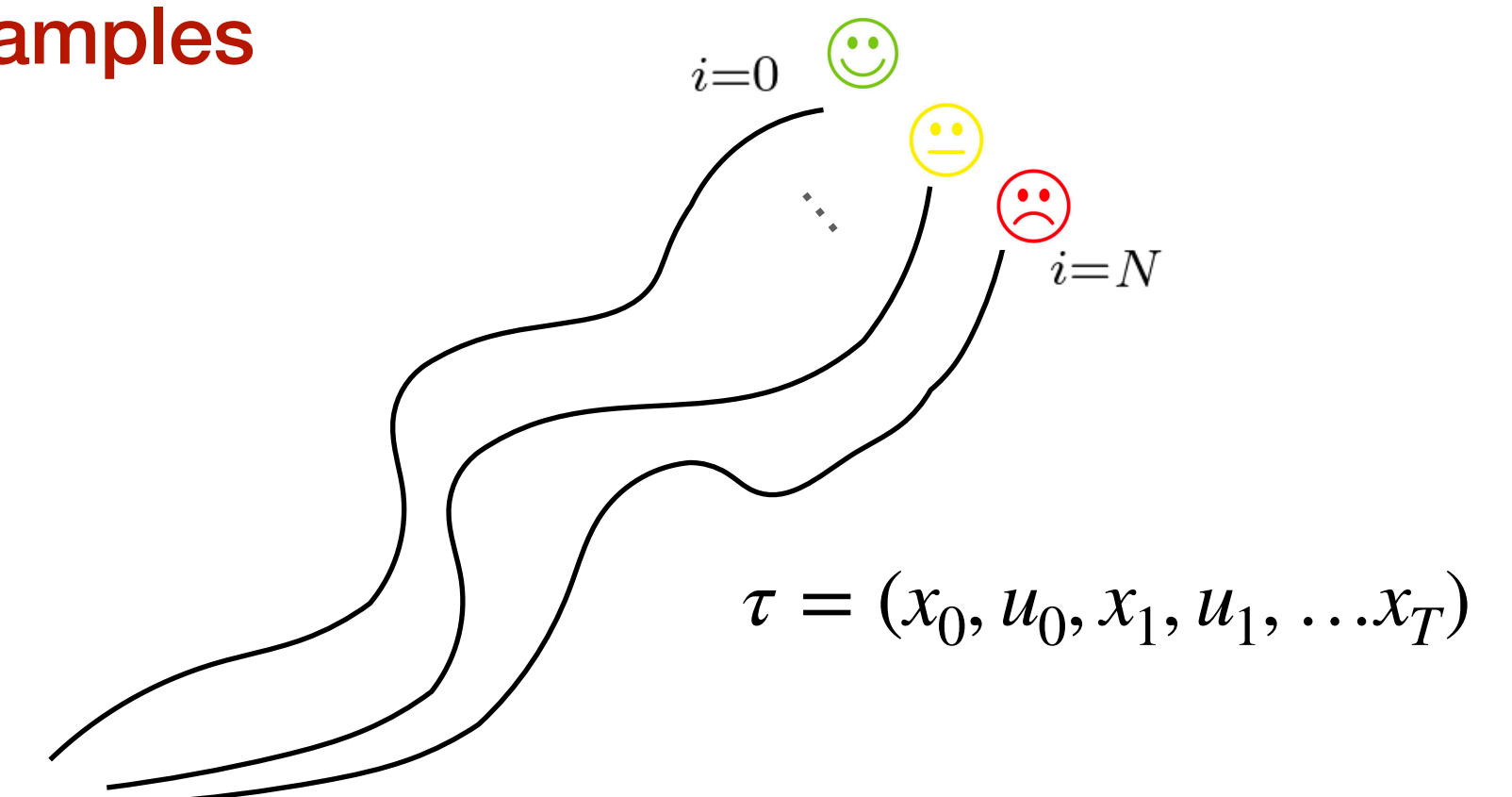
(1) estimate its gradient  $\nabla_{\theta} J(\theta)$

(2) cast the learning process as approximate gradient ascent on  $J(\theta)$        $\theta \leftarrow \theta + \alpha \nabla_{\theta} J(\theta)$



How can we evaluate the expectation in the objective? As usual in RL, **through samples**

$$J(\theta) = \mathbb{E}_{\tau \sim p(\tau)} \left[ \sum_{t \geq 0} R(x_t, u_t) \right] \approx \frac{1}{N} \sum_i \sum_t R(x_{i,t}, u_{i,t})$$



# Direct policy gradient

- In order to solve the problem through gradient-based optimization we need to compute  $\nabla_{\theta} J(\theta)$

- Let us define the compact notation  $r(\tau) = \sum_{t=1}^T R(x_t, u_t)$

- By definition of expectation  $J(\theta) = \mathbb{E}_{\tau \sim p(\tau)} [r(\tau)] = \int p_{\theta}(\tau) r(\tau) d\tau$

- We can then write the gradient  $\nabla_{\theta} J(\theta) = \int \nabla_{\theta} p_{\theta}(\tau) r(\tau) d\tau$

**Problem:** gradient depends on unknown dynamics and initial state distribution through  $p_{\theta}(\tau)$

**Useful identity:**

$$p_{\theta}(\tau) \nabla_{\theta} \log p_{\theta}(\tau) = p_{\theta}(\tau) \frac{\nabla_{\theta} p_{\theta}(\tau)}{p_{\theta}(\tau)} = \nabla_{\theta} p_{\theta}(\tau)$$

$$\nabla_{\theta} J(\theta) = \int \nabla_{\theta} p_{\theta}(\tau) r(\tau) d\tau = \int p_{\theta}(\tau) \nabla_{\theta} \log p_{\theta}(\tau) r(\tau) d\tau = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} [\nabla_{\theta} \log p_{\theta}(\tau) r(\tau)]$$

On the right track since we can evaluate expectations through samples... but we still have  $\nabla_{\theta} \log p_{\theta}(\tau)$  in the equation



# Direct policy gradient

Let us recall the trajectory distribution

$$p(x_0, u_0, \dots, x_T) = p(\tau) = p(x_0) \prod_{t=1}^T \pi(u_t | x_t) p(x_{t+1} | x_t, u_t)$$

$$\log p(\tau) = \log p(x_0) + \sum_{t=1}^T \log \pi_{\theta}(u_t | x_t) + \log p(x_{t+1} | x_t, u_t)$$

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} [\nabla_{\theta} \log p_{\theta}(\tau) r(\tau)] = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} \left[ \nabla_{\theta} \left[ \cancel{\log p(x_0)} + \sum_{t=1}^T \log \pi_{\theta}(u_t | x_t) + \cancel{\log p(x_{t+1} | x_t, u_t)} \right] r(\tau) \right]$$

- When taking the gradient w.r.t.  $\theta$ ,  $\log p(x_0)$ ,  $\log p(x_{t+1} | x_t, u_t)$  do not depend on  $\theta$
- While we can evaluate the log probability under our parametric policy  $\pi_{\theta}$
- This enable us to re-write the gradient  $\nabla_{\theta} J(\theta)$  as:

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_t | x_t) \right) \left( \sum_{t=1}^T R(x_t, u_t) \right) \right]$$

**Everything inside this expectation is known**

# Direct policy gradient

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_t | x_t) \right) \left( \sum_{t=1}^T R(x_t, u_t) \right) \right]$$

Everything inside this expectation is known

- Recall how we use samples to evaluate the objective:  $J(\theta) = \mathbb{E}_{\tau \sim p(\tau)} \left[ \sum_{t \geq 0} R(x_t, u_t) \right] \approx \frac{1}{N} \sum_i \sum_t R(x_{i,t}, u_{i,t})$
- We can use the same idea to evaluate the gradient:

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_t | x_t) \right) \left( \sum_{t=1}^T R(x_t, u_t) \right) \right] \approx \frac{1}{N} \sum_{i=1}^N \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \right) \left( \sum_{t=1}^T R(x_{i,t}, u_{i,t}) \right) \right]$$

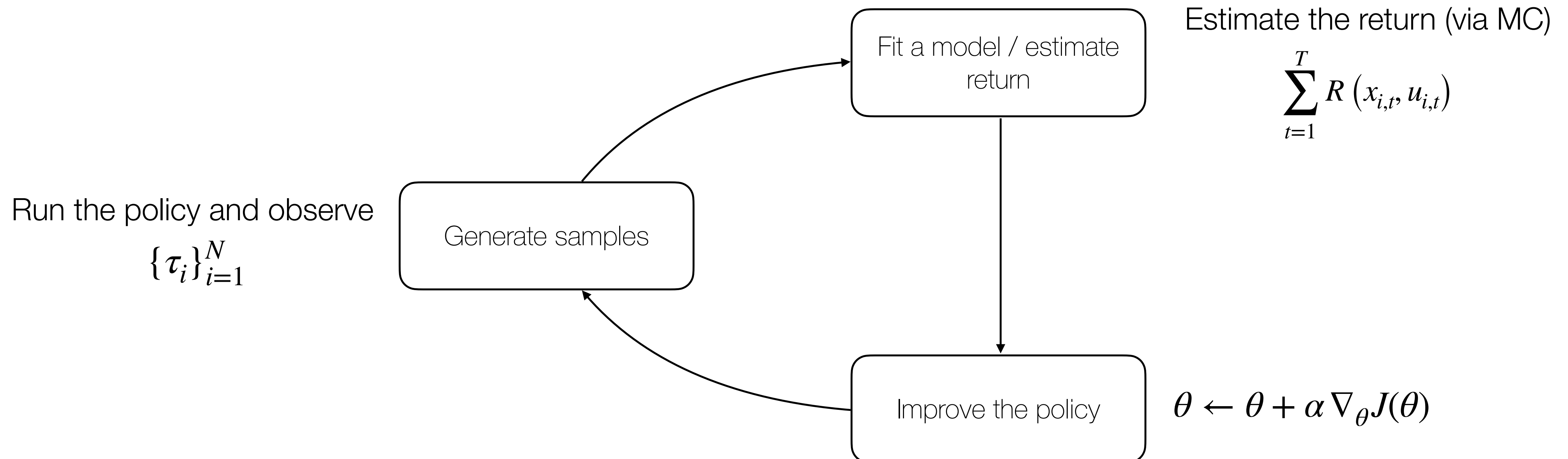
# REINFORCE algorithm

The procedure described so far gives us the basic policy gradient algorithms, a.k.a. REINFORCE:

1. Sample trajectories  $\{\tau_i\}_{i=1}^N$  from  $\pi_\theta(u_t | x_t)$ , i.e. run the policy in the environment

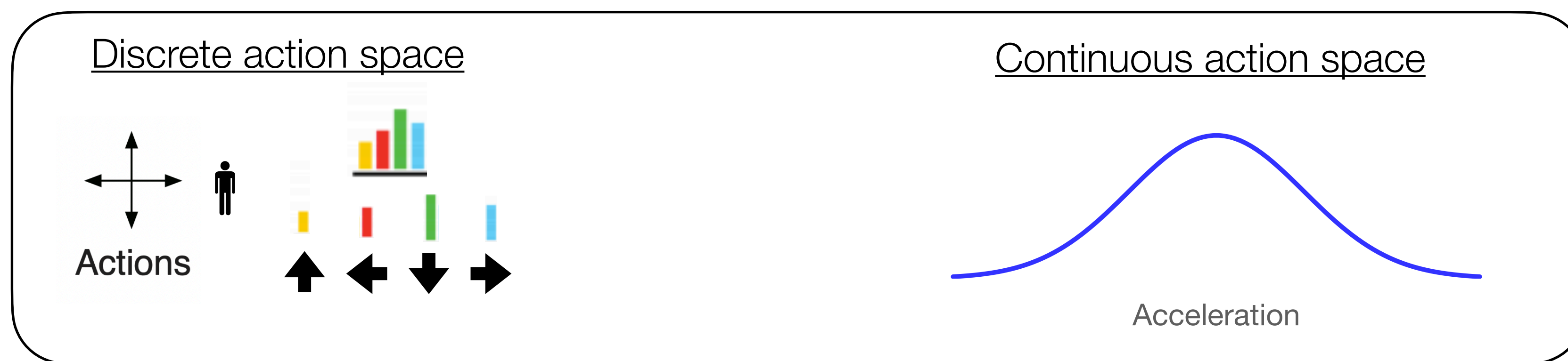
2. Evaluate the policy gradient  $\nabla_\theta J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left[ \left( \sum_{t=1}^T \nabla_\theta \log \pi_\theta(u_{i,t} | x_{i,t}) \right) \left( \sum_{t=1}^T R(x_{i,t}, u_{i,t}) \right) \right]$

3. Take a gradient step to update the policy  $\theta \leftarrow \theta + \alpha \nabla_\theta J(\theta)$

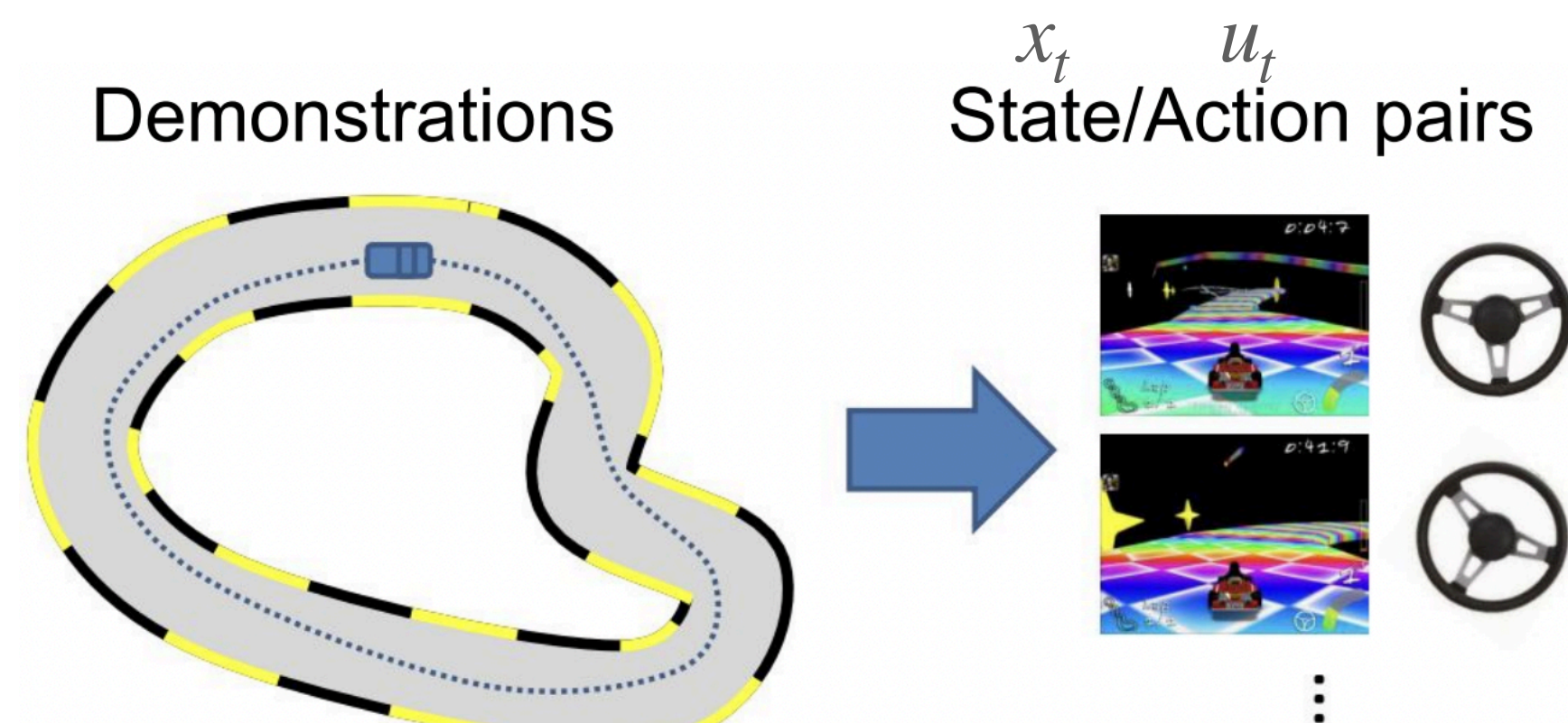


# Intuition: “what is PG doing?”

Consider the expression we derived for the policy gradient  $\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \right) \left( \sum_{t=1}^T R(x_{i,t}, u_{i,t}) \right) \right]$



Let's compare it with the expression of the gradient when performing maximum likelihood (e.g., supervised learning):



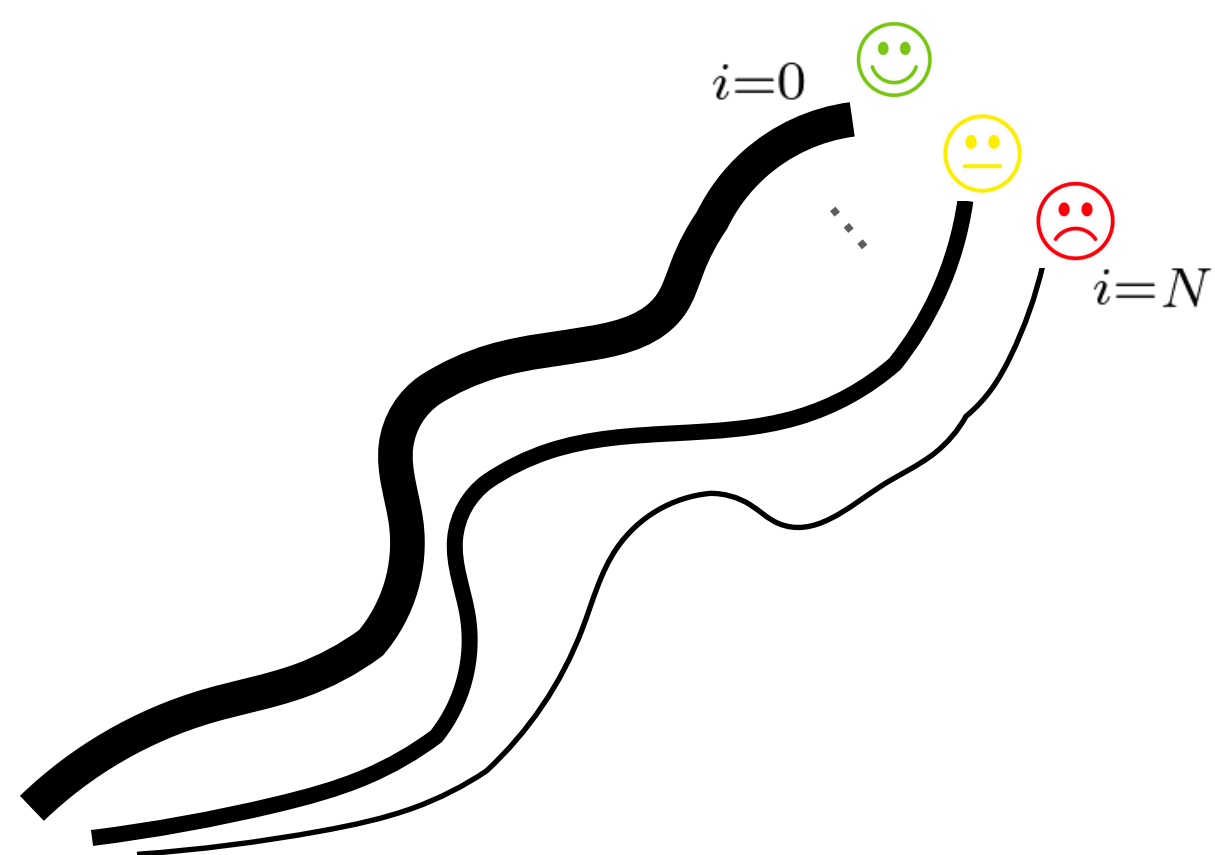
$$\nabla_{\theta} J_{MLE}(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \right) \right]$$

**The policy gradient is a weighted version of the MLE gradient**

# Intuition: “what is PG doing?”

Policy gradient:  $\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \right) \left( \sum_{t=1}^T R(x_{i,t}, u_{i,t}) \right) \right]$

Maximum Likelihood:  $\nabla_{\theta} J_{MLE}(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \right) \right]$



Taking a step in the direction the policy gradient essentially means:

$$\theta \leftarrow \theta + \alpha \nabla_{\theta} J(\pi_{\theta})$$

*“Change parameters  $\theta$  s.t. trajectories with higher reward have higher probability”*

**PG formalizes the idea of learning by “trial and error”**



# Outline

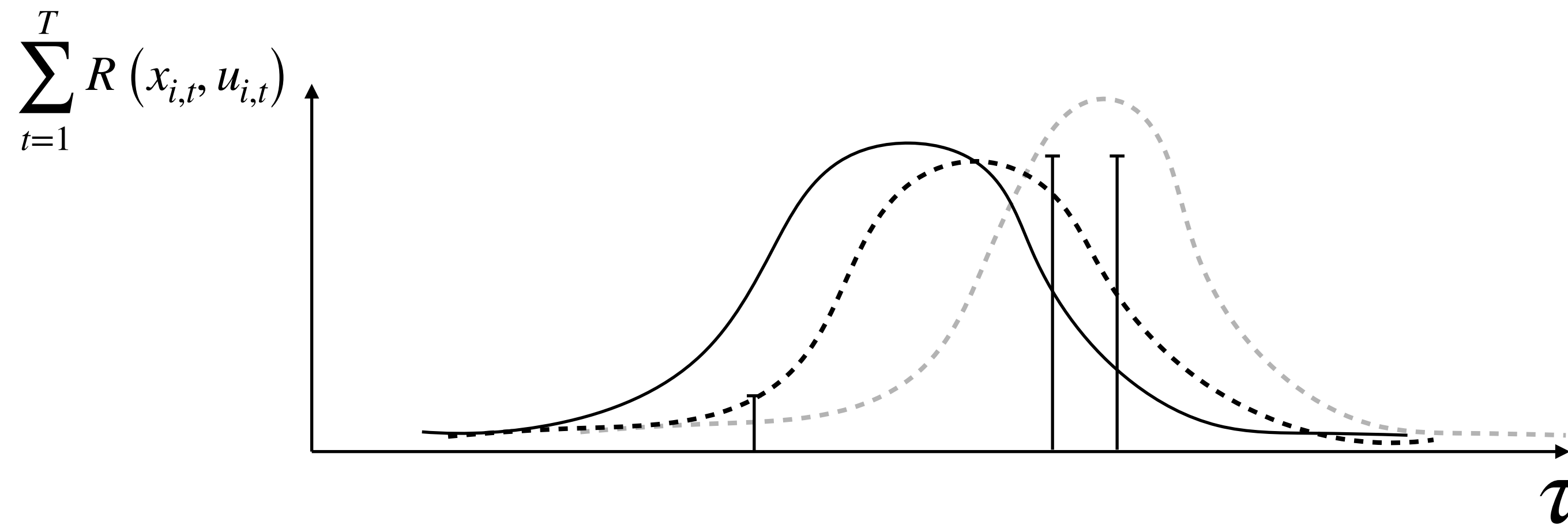
Intro to policy gradients

- REINFORCE algorithm
- Reducing variance of policy gradient

# Problem: high variance of the PG

Policy gradient:  $\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \right) \left( \sum_{t=1}^T R(x_{i,t}, u_{i,t}) \right) \right]$

Let's consider the following example:



- Depending on the sample, the policy gradient can vary wildly: PG estimator has **high variance**
- This negatively affects learning: worse performance, slower convergence

A lot of research in the domain of Policy Optimization revolves around finding ways to lower the variance of the policy gradient

# Reducing the variance

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left[ \left( \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \right) \left( \sum_{t=1}^T R(x_{i,t}, u_{i,t}) \right) \right]$$

A first simple approach to reduce the variance entails using *causality*: “policy at time  $t'$  cannot affect reward at time  $t < t'$ ”

Consider this equivalent expression:

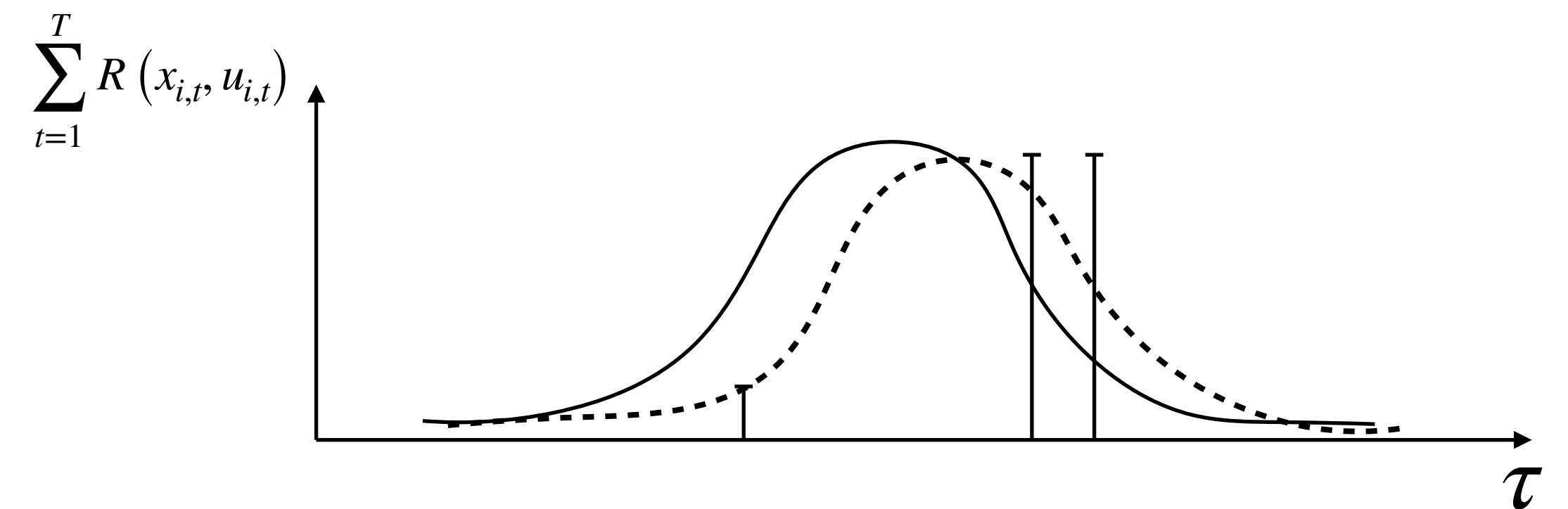
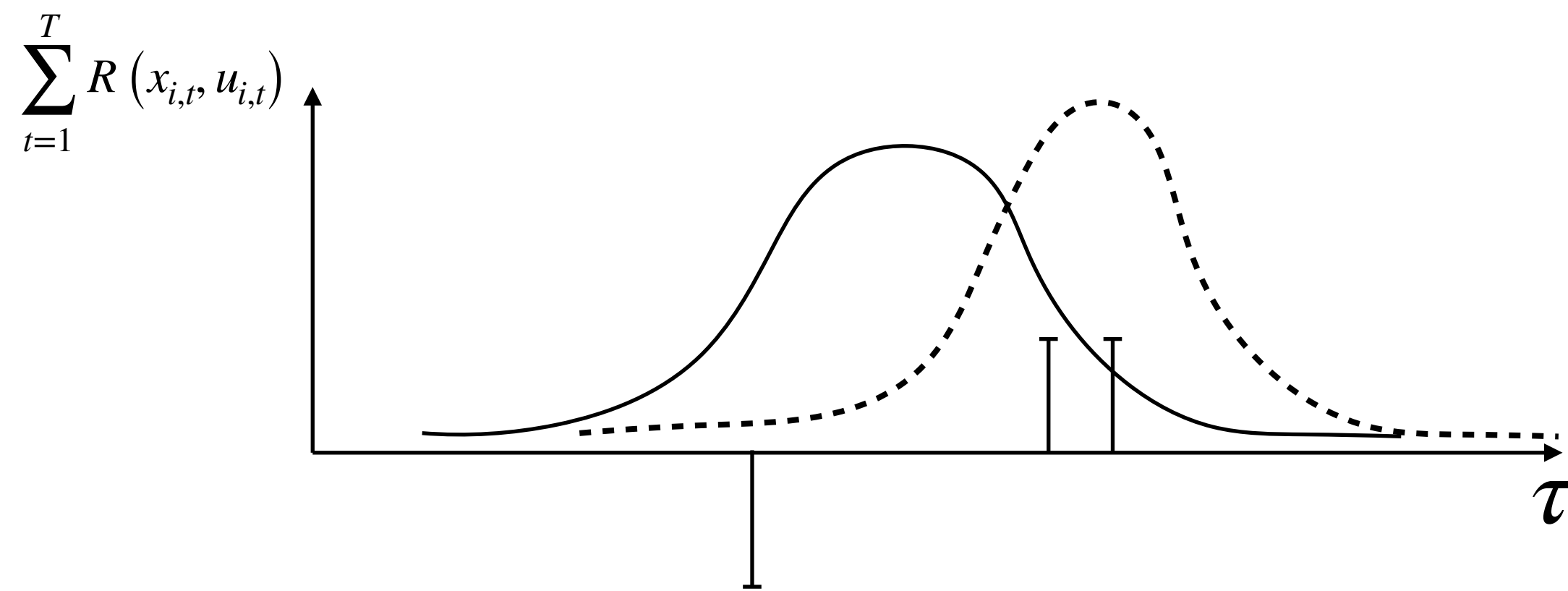
$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \left( \sum_{\substack{t'=t \\ \text{red}}}^T R(x_{i,t'}, u_{i,t'}) \right)$$



# Baseline

A second (and extremely important) approach to reduce variance of PG estimators relates with the concept of **baseline**

Let's reconsider our intuition on PG, i.e., “making good behavior more likely”



However, PG will only do this **if the returns are centered** (e.g., consider the counter-example on the right)

Intuitively, we want to “center” our returns, such that:

- The probability of behavior that is better than average gets increased
- The probability of behavior that is worse than average gets decreased

We are going to subtract a baseline  $b$  from the expression of the PG

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \nabla_{\theta} \log \pi_{\theta}(\tau) [r(\tau) - b]$$

# A closer look at the baseline

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} [\nabla_{\theta} \log p_{\theta}(\tau) r(\tau)] \rightarrow \nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} [\nabla_{\theta} \log p_{\theta}(\tau) (r(\tau) - b)]$$

**Claim:** adding the baseline does not change the value of the expected gradient

- To prove that, let's consider the following expectation:

**Useful identity:**

$$p_{\theta}(\tau) \nabla_{\theta} \log p_{\theta}(\tau) = p_{\theta}(\tau) \frac{\nabla_{\theta} p_{\theta}(\tau)}{p_{\theta}(\tau)} = \nabla_{\theta} p_{\theta}(\tau)$$

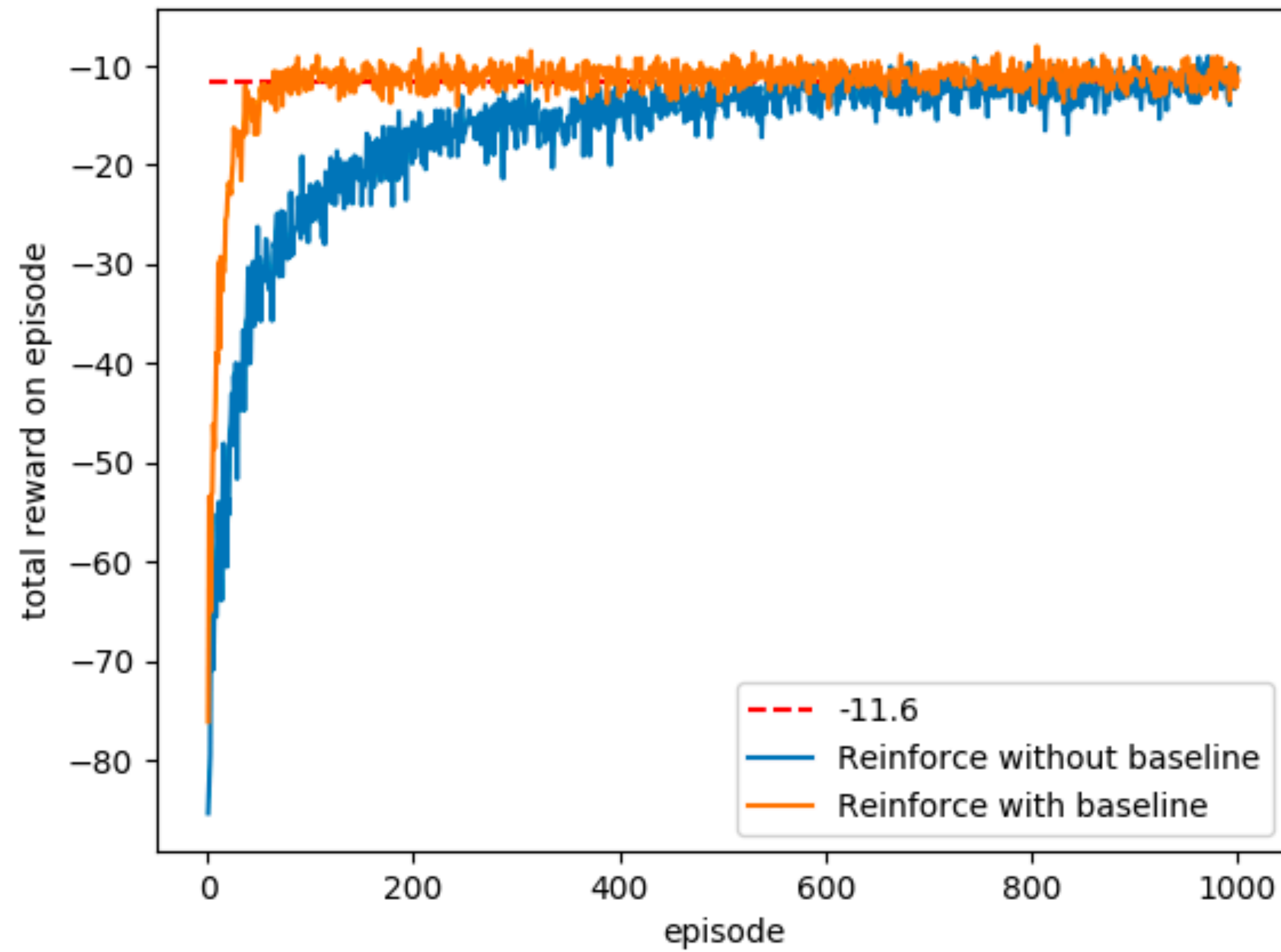
$$\mathbb{E} [\nabla_{\theta} \log p_{\theta}(\tau) b] = \int p_{\theta}(\tau) \nabla_{\theta} \log p_{\theta}(\tau) b d\tau = \int \nabla_{\theta} p_{\theta}(\tau) b d\tau = b \nabla_{\theta} \int p_{\theta}(\tau) d\tau = b \nabla_{\theta} 1 = 0$$

which makes our estimate of the gradient (with baseline) *unbiased* in expectation

- An extremely effective choice of the baseline is the **average return**,  $b = \frac{1}{N} \sum_{i=1}^N r(\tau_i)$

(We'll see how this motivates many popular RL algorithms...)

# Example



# Properties of policy gradient

At a high-level, we've been defining a scheme where:

- Given the RL objective  $J(\theta) = \mathbb{E}_{\tau \sim p(\tau)} [r(\tau)] = \int p_{\theta}(\tau) r(\tau) d\tau$
- We maximize the objective w.r.t.  $\theta$  by:
  - Computing the gradient  $\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} [\nabla_{\theta} \log p_{\theta}(\tau) r(\tau)]$
  - Taking a gradient step to update the policy  $\theta \leftarrow \theta + \alpha \nabla_{\theta} J(\theta)$

## Question:

Is this on- or off-policy? And why?

# Outline

## Actor-Critic methods

- Advantage
- Architecture design

# From PG to Actor-Critic methods

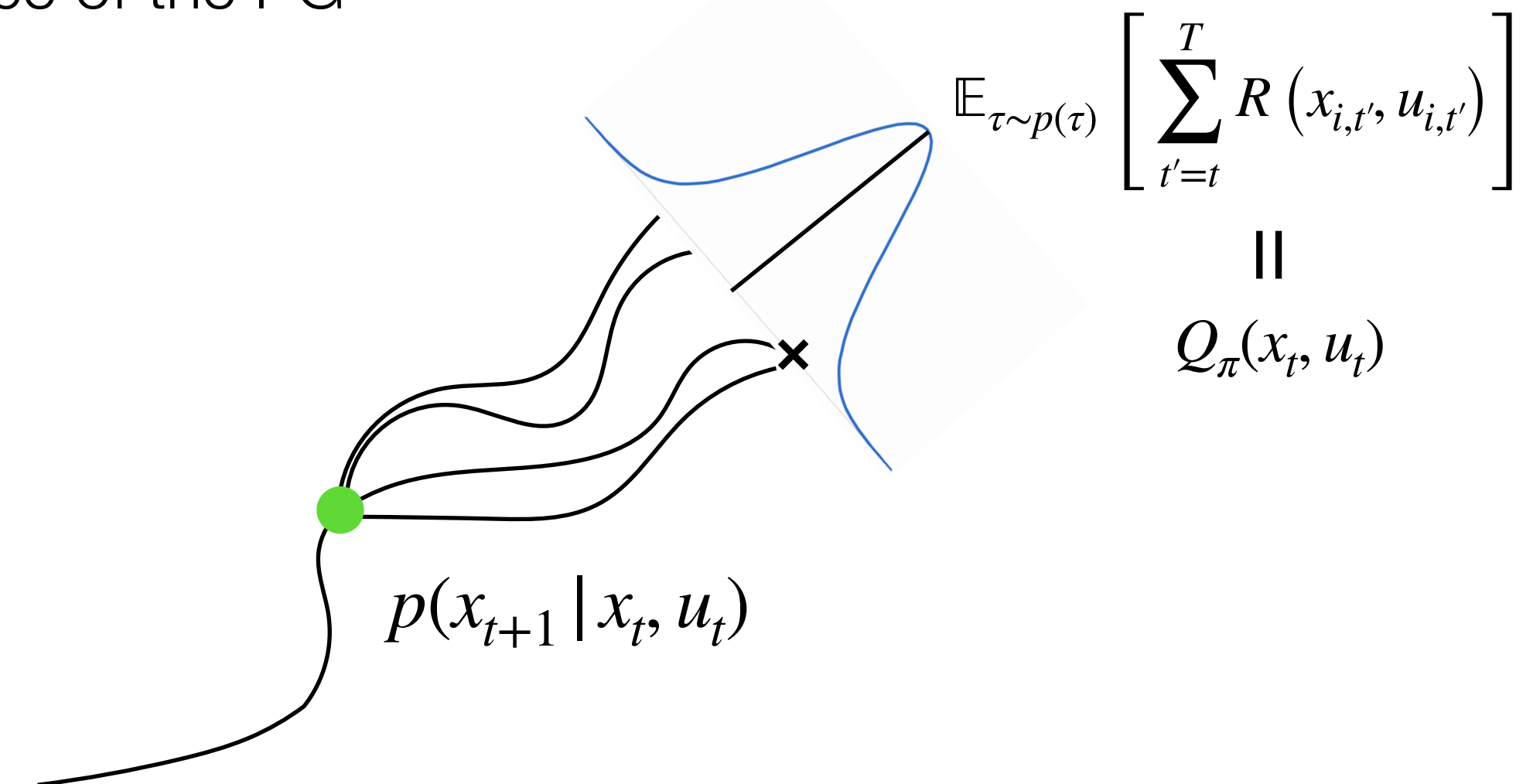
Once again, let's consider the policy gradient  $\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \left( \sum_{t'=t}^T R(x_{i,t'}, u_{i,t'}) \right)$

“reward-to-go”

This one-sample estimate of the reward-to-go contributes to the high variance of the PG

The idea of actor-critic methods is to define:

- An “actor”, i.e., a policy  $\pi_{\theta}(u_t | x_t)$
- A “critic” to better estimate the “reward-to-go”,  
e.g., estimate Q-values through function approximation  $Q_{\phi}(x_t, u_t)$



By using this better estimate of the reward-to-go we can get a lower variance policy gradient:

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) Q_{\phi}(x_t, u_t)$$

# What about the baseline?

Can we use a baseline when using the approximate reward-to-go and reduce the variance even further?

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \left( Q_{\phi}(x_{i,t}, u_{i,t}) - b \right)$$

- An effective choice for  $b$  is a state-dependent baseline  $b(x_t) = \mathbb{E}_{u_t \sim \pi(u_t | x_t)} [Q(x_t, u_t)] = V(x_t)$
- We can thus rewrite:

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) \left( Q_{\phi}(x_{i,t}, u_{i,t}) - V(x_{i,t}) \right)$$

“How much  $u_t$  is  
better than the  
average action  
in  $x_t$ ”



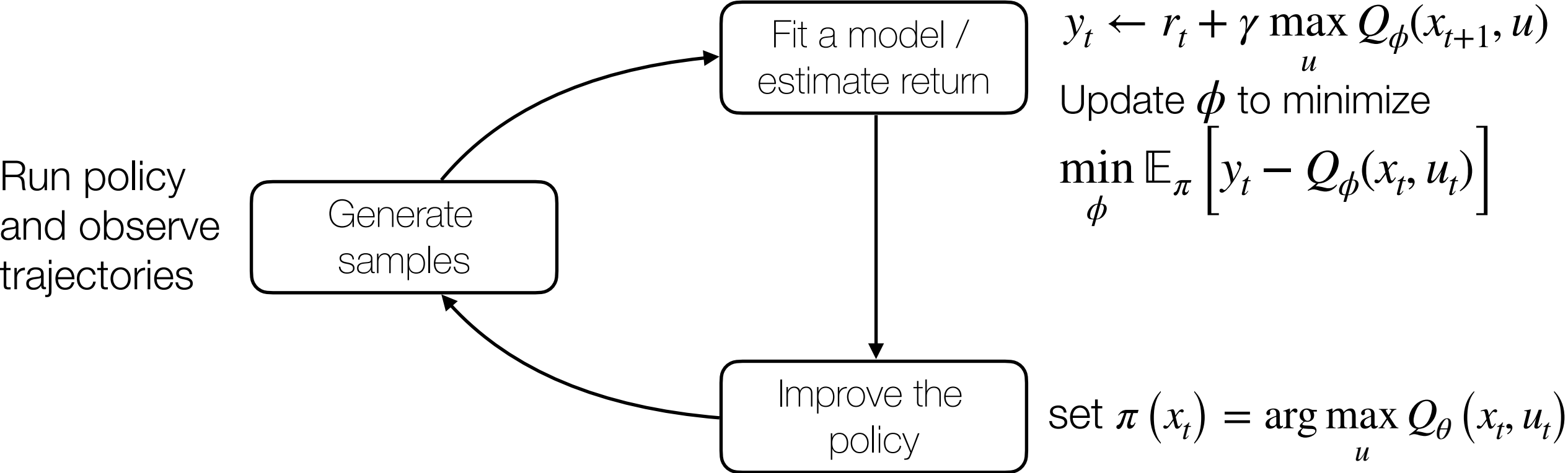
Following this gradient:

- increases the probability of actions that have returns better than average
- decreases the probability of actions that have returns worse than average

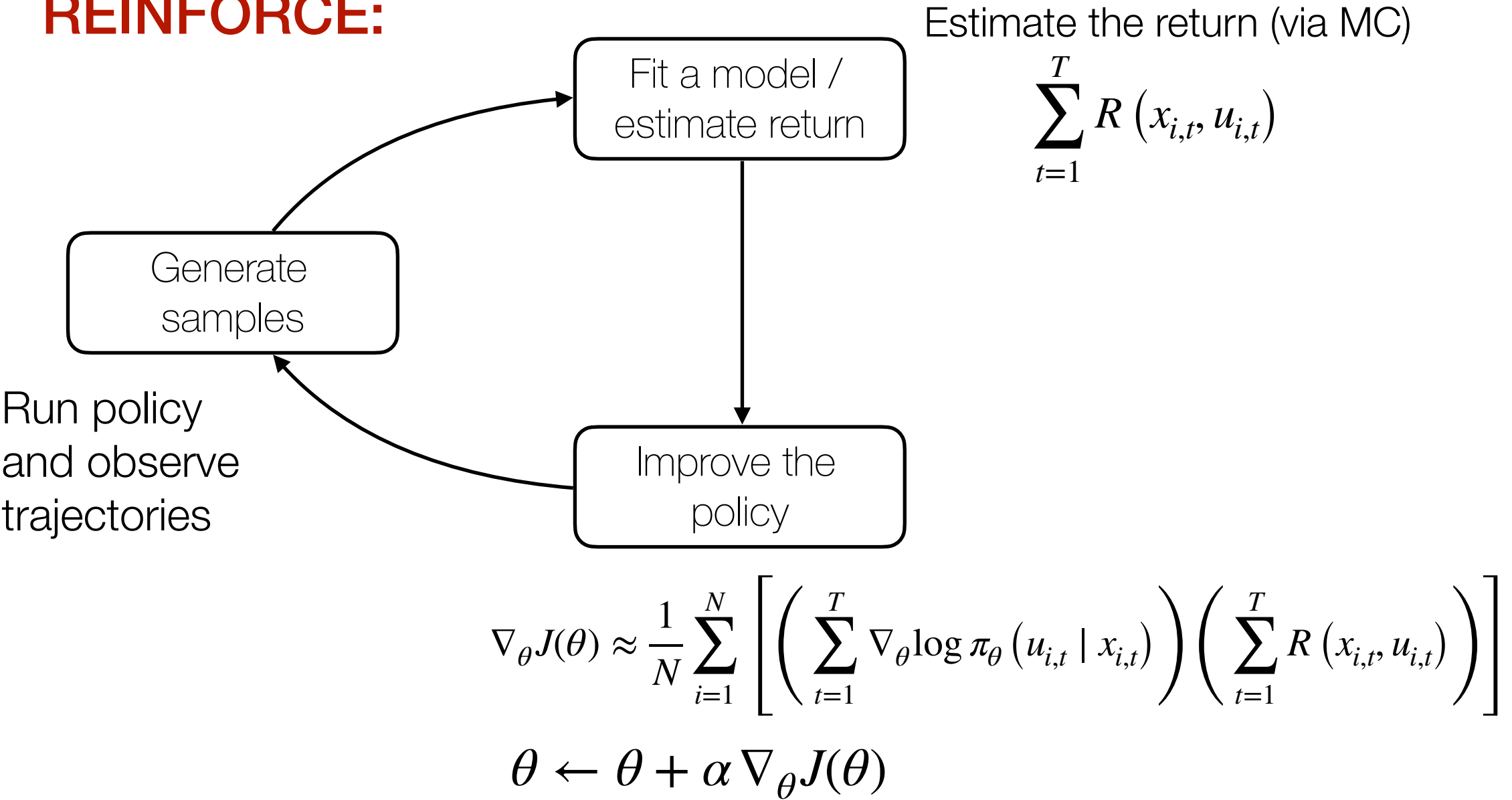
- The function  $A(x_t, u_t) = Q_{\phi}(x_t, u_t) - V(x_t)$  is usually referred to as **advantage function**



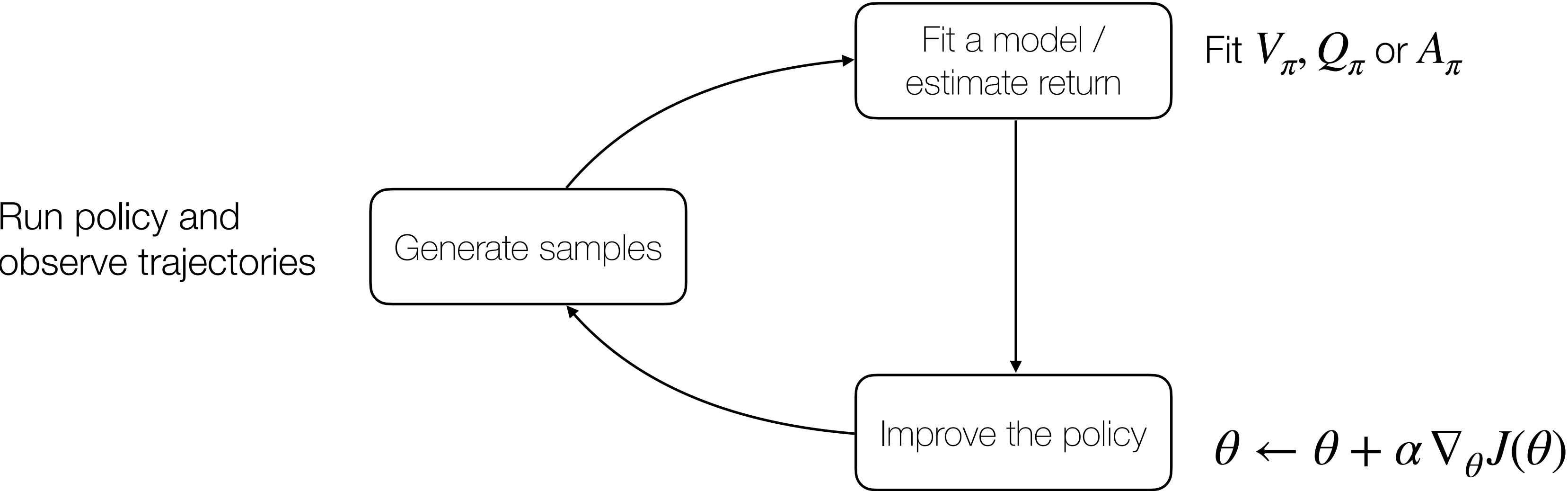
Fitted Q-learning:



REINFORCE:

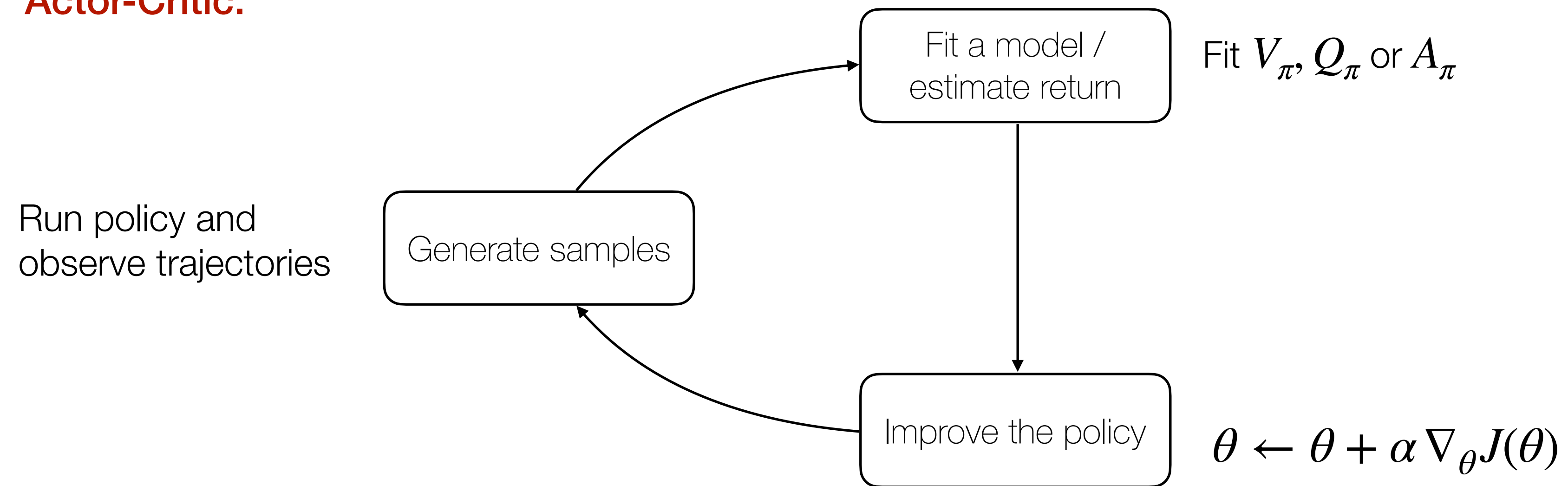


Actor-Critic:





## Actor-Critic:



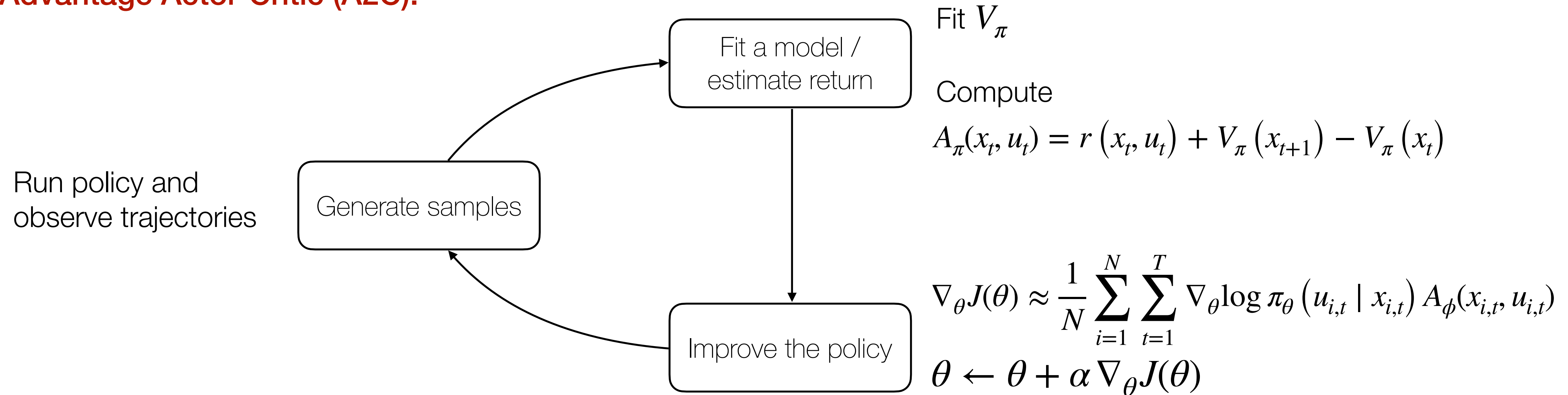
What quantity should we estimate? What are the trade-offs between estimating  $V_\pi(x_t)$ ,  $Q_\pi(x_t, u_t)$  or  $A_\pi$ ? No wrong/right, answer, it depends. For now, let's consider the complexity of the estimation problem (i.e., fitting  $V_\pi$  is easier: only  $x_t$  as input)

$$Q_\pi(x_t, u_t) = \mathbb{E}_{\tau \sim p(\tau)} \left[ \sum_{t'=t}^T R(x_{i,t'}, u_{i,t'}) \right] = r(x_t, u_t) + \mathbb{E}_{x_{t+1} \sim p(x_{t+1} | x_t, u_t)} \left[ V_\pi(x_{t+1}) \right] \approx r(x_t, u_t) + V_\pi(x_{t+1})$$

$$A_\pi(x_t, u_t) = Q_\pi(x_t, u_t) - V_\pi(x_t) \approx r(x_t, u_t) + V_\pi(x_{t+1}) - V_\pi(x_t)$$

This enables us to “only” fit  $V_\pi$

## Advantage Actor-Critic (A2C):



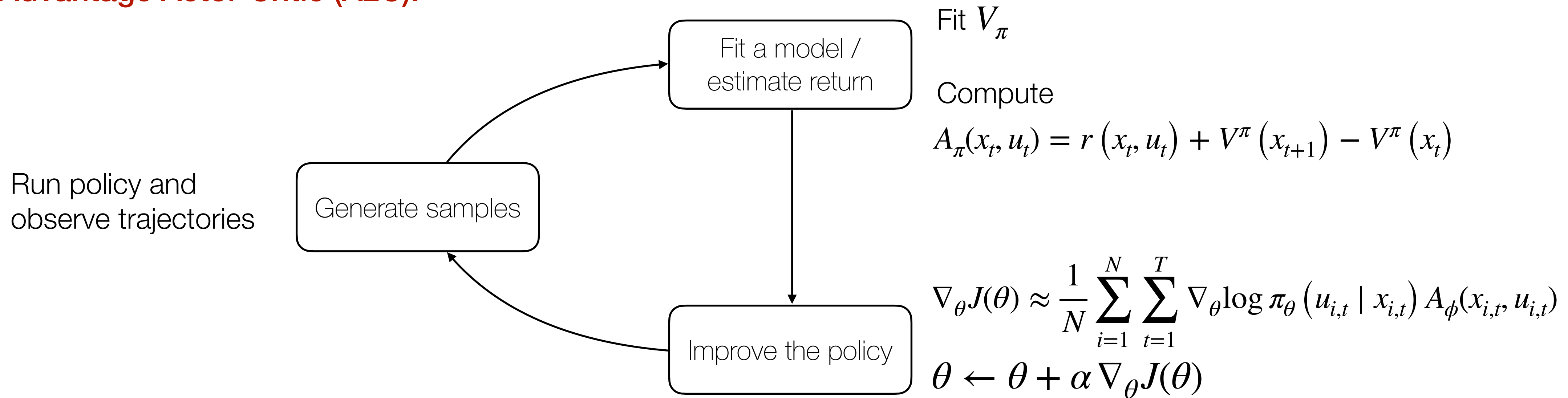
What quantity should we estimate? What are the trade-offs between estimating  $V_\pi(x_t)$ ,  $Q_\pi(x_t, u_t)$  or  $A_\pi$ ? No wrong/right, answer, it depends. For now, let's consider the complexity of the estimation problem (i.e., fitting  $V_\pi$  is easier: only  $x_t$  as input)

$$Q_\pi(x_t, u_t) = \mathbb{E}_{\tau \sim p(\tau)} \left[ \sum_{t'=t}^T R(x_{i,t'}, u_{i,t'}) \right] = r(x_t, u_t) + \mathbb{E}_{x_{t+1} \sim p(x_{t+1} | x_t, u_t)} \left[ V_\pi(x_{t+1}) \right] \approx r(x_t, u_t) + V_\pi(x_{t+1})$$

$$A_\pi(x_t, u_t) = Q_\pi(x_t, u_t) - V_\pi(x_t) \approx r(x_t, u_t) + V_\pi(x_{t+1}) - V_\pi(x_t)$$

This enables us to “only” fit  $V_\pi$

## Advantage Actor-Critic (A2C):



When fitting  $V_\pi$ , we can use different *targets* to define the supervised learning labels

### Question:

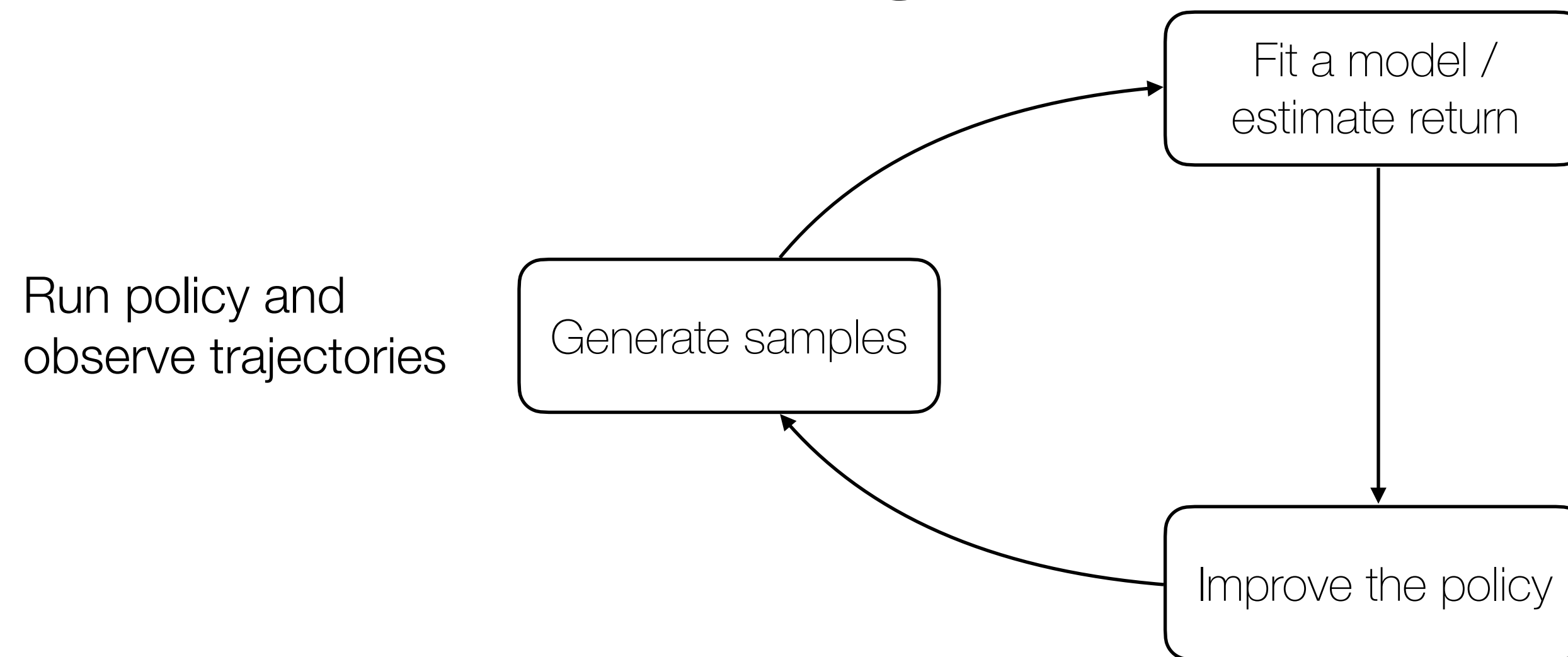
How to fit with MC target?

1. Collect dataset  $\mathcal{D} = \{(x_t, G_t)\}$ ,  $G_t = \sum_{t'=t}^T R(x_{t'}, u_{t'})$
2. Supervised regression on  $\mathcal{D}$

How to fit with TD target?

1. Collect dataset  $\mathcal{D} = \{(x_t, r_t + \gamma \hat{V}_\theta(x_t))\}$
2. Supervised regression on  $\mathcal{D}$

# Architecture design



Fit  $V_\pi$

Compute

$$A_\pi(x_t, u_t) = r(x_t, u_t) + V_\pi(x_{t+1}) - V_\pi(x_t)$$

$$\nabla_\theta J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \nabla_\theta \log \pi_\theta(u_{i,t} | x_{i,t}) A_\phi(x_{i,t}, u_{i,t})$$

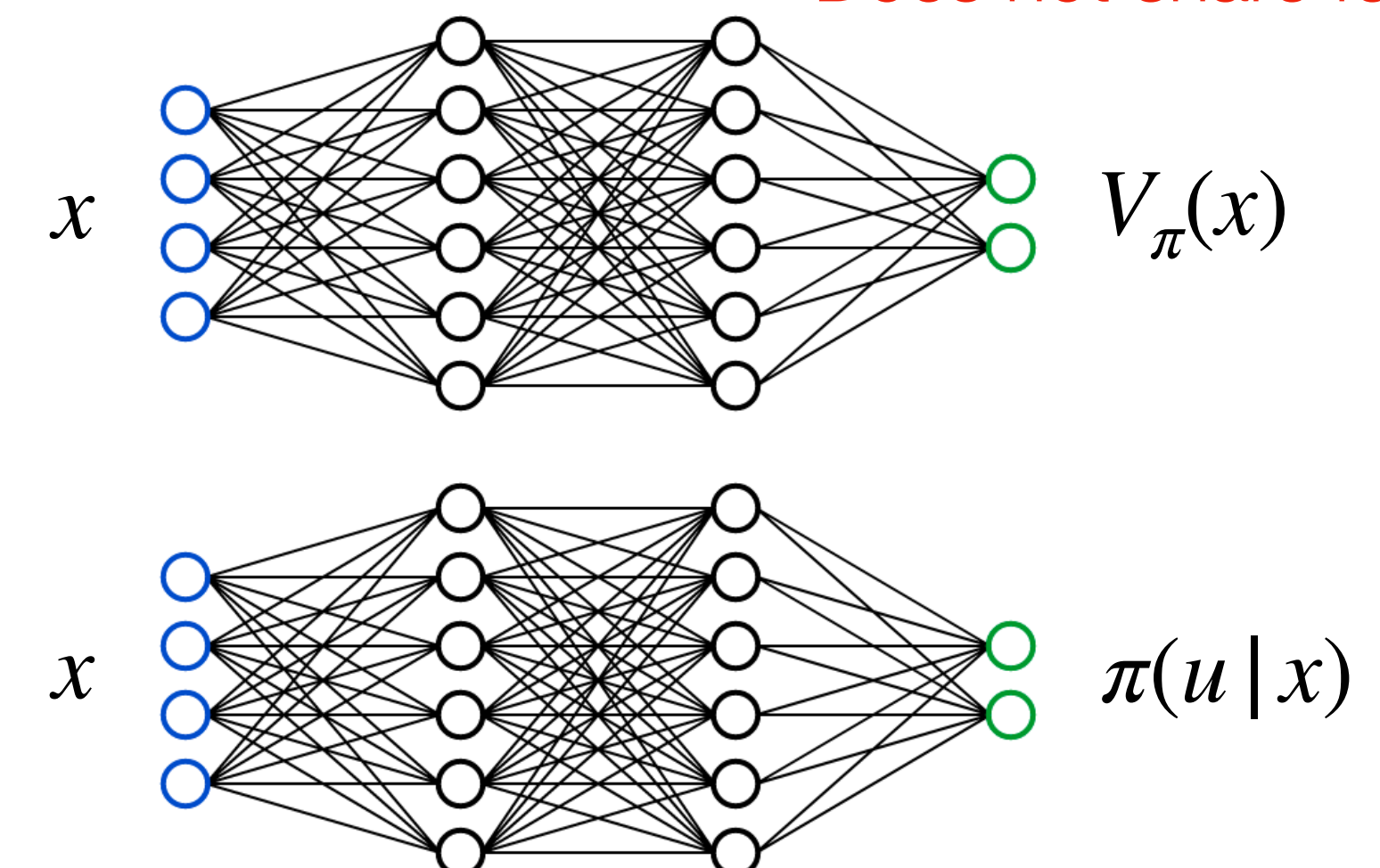
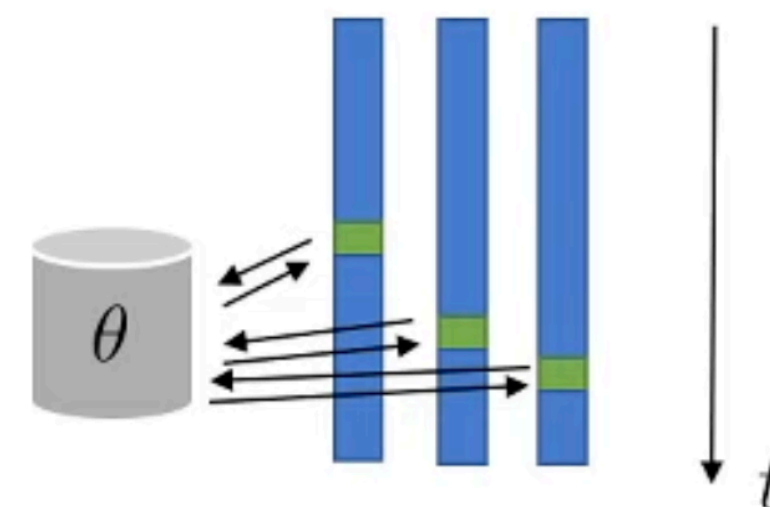
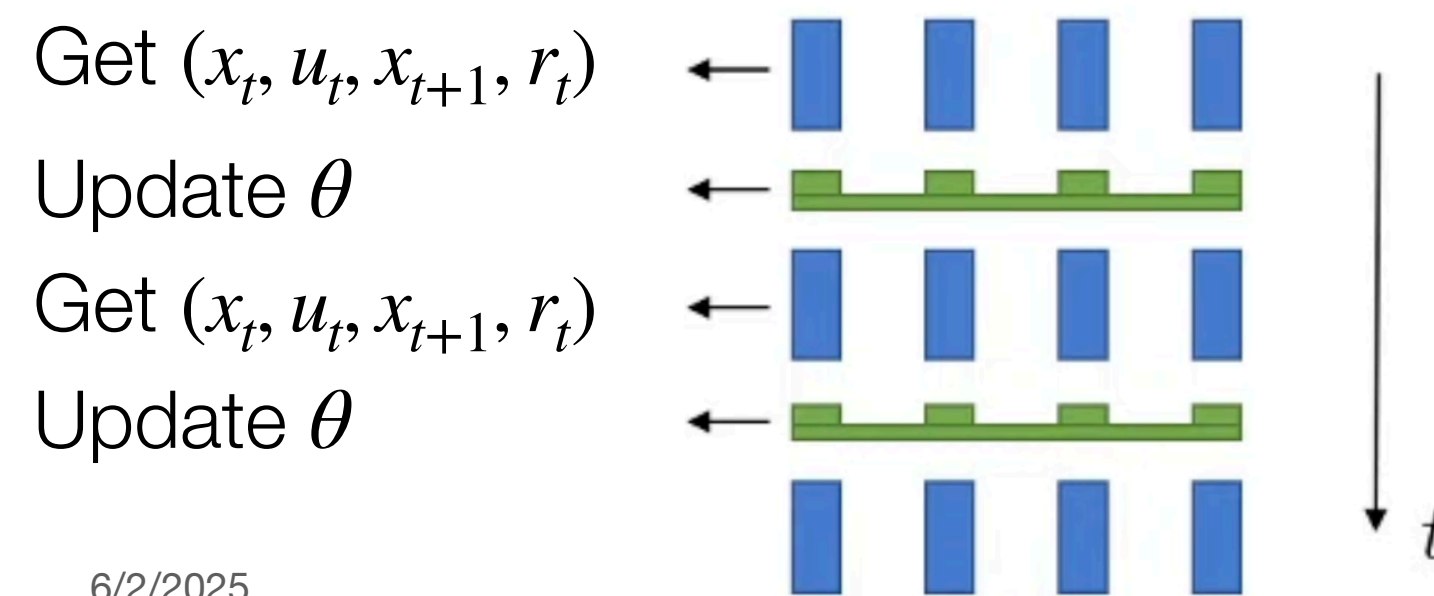
$$\theta \leftarrow \theta + \alpha \nabla_\theta J(\theta)$$

Simple, typically more stable

Does not share features

In practice, one could opt for different designs of this same algorithms, e.g.,:

- Two network vs shared network
- Parallel processing: synchronized vs asynchronous



# Outline

Deep RL Algorithms & Applications

# Reducing RL to optimization

- Much of modern ML entails reducing learning to a numerical optimization problem
  - Supervised learning as *training error minimization*
- This is different from what we have seen so far in RL:
  - Q-learning: fixed-point iteration  $\rightarrow$  can (in principle) include all transitions seen so far, however, it optimizes for the wrong objective
  - Policy gradient: yes, stochastic gradients of the RL objective, but no optimization problem
- We'll discuss approaches that define an optimization problem that allows us to do a small update to policy  $\pi$ , based on data sampled from  $\pi$



# Defining the objective

- We discussed how, in PO, we want to compute the following gradient  $\nabla_{\theta} J(\theta) = \mathbb{E} \left[ \nabla_{\theta} \log \pi_{\theta}(u_t | x_t) A(x_t, u_t) \right]$
- To implement this using modern auto-diff tools (e.g., Torch, Jax, Tensorflow), this means writing the following loss function:

$$L^{PG}(\theta) = \mathbb{E} \left[ \log \pi_{\theta}(u_t | x_t) A(x_t, u_t) \right]$$

- But we don't want to optimize it too far, since we are not working with the *true* advantage, rather with a noisy estimate
- Equivalently differentiate

$$L^{IS}(\theta) = \mathbb{E}_{\tau \sim p_{\theta}(\tau)} \left[ \frac{\pi_{\theta}(u_t | x_t)}{\pi_{\theta_{old}}(u_t | x_t)} A(\tau) \right]$$

- If we take the derivative of  $L^{IS}$  and evaluate at  $\theta = \theta_{old}$ , we get the same gradient

$$\nabla_{\theta} \log f(\theta) \Big|_{\theta_{old}} = \frac{\nabla_{\theta} f(\theta) \Big|_{\theta_{old}}}{f(\theta_{old})} = \nabla_{\theta} \left( \frac{f(\theta)}{f(\theta_{old})} \right) \Big|_{\theta_{old}}$$

# Trust Region Policy Optimization (TRPO)

$$\begin{aligned} & \underset{\theta}{\text{maximize}} \quad \hat{\mathbb{E}}_t \left[ \frac{\pi_{\theta}(u_t | x_t)}{\pi_{\theta_{old}}(u_t | x_t)} \hat{A}_t \right] \\ & \text{subject to} \quad \hat{\mathbb{E}}_t \left[ \text{KL}[\pi_{\theta_{old}}(\cdot | x_t), \pi_{\theta}(\cdot | x_t)] \right] \leq \delta \end{aligned}$$

- Main idea: use trust region to constrain change **in distribution space** (opposed to e.g., parameter space)
- Hard to use with architectures with multiple outputs, e.g., policy and value function
- Empirically performs poorly on tasks requiring deep nets, e.g., deep CNNs, RNNs
- Conjugate gradient makes implementation more complicated



# Proximal Policy Optimization (PPO)

- Can we solve the problem defined in TRPO without second-order optimization?

## PPO v1 - Surrogate loss with Lagrange multipliers

$$\underset{\theta}{\text{maximize}} \quad \hat{\mathbb{E}}_t \left[ \frac{\pi_{\theta}(u_t | x_t)}{\pi_{\theta_{old}}(u_t | x_t)} \hat{A}_t \right] + \beta \left( \hat{\mathbb{E}}_t \left[ \text{KL}[\pi_{\theta_{old}}(\cdot | x_t), \pi_{\theta}(\cdot | x_t)] \right] - \delta \right)$$

- Run SGD on the above objective
- Do dual descent update for  $\beta$

## PPO v2 - Clipped surrogate loss

$$r(\theta) = \frac{\pi_{\theta}(u_t | x_t)}{\pi_{\theta_{old}}(u_t | x_t)}, \quad r(\theta_{old}) = 1$$

$$\underset{\theta}{\text{maximize}} \quad \hat{\mathbb{E}}_t \left[ \min(r(\theta)A(\tau), \text{clip}(r(\theta), 1 - \epsilon, 1 + \epsilon)A(\tau)) \right]$$

- Heuristically replicates constraint in the objective
- One of the (if not the) most popular PO algorithm

# Examples: Maze Navigation

- Mnih et al. 2016 “Asynchronous Methods for Deep Reinforcement Learning”
- Advantage Actor-Critic
- Asynchronous parallel workers
- Policy and Value networks: CNNs & RNNs



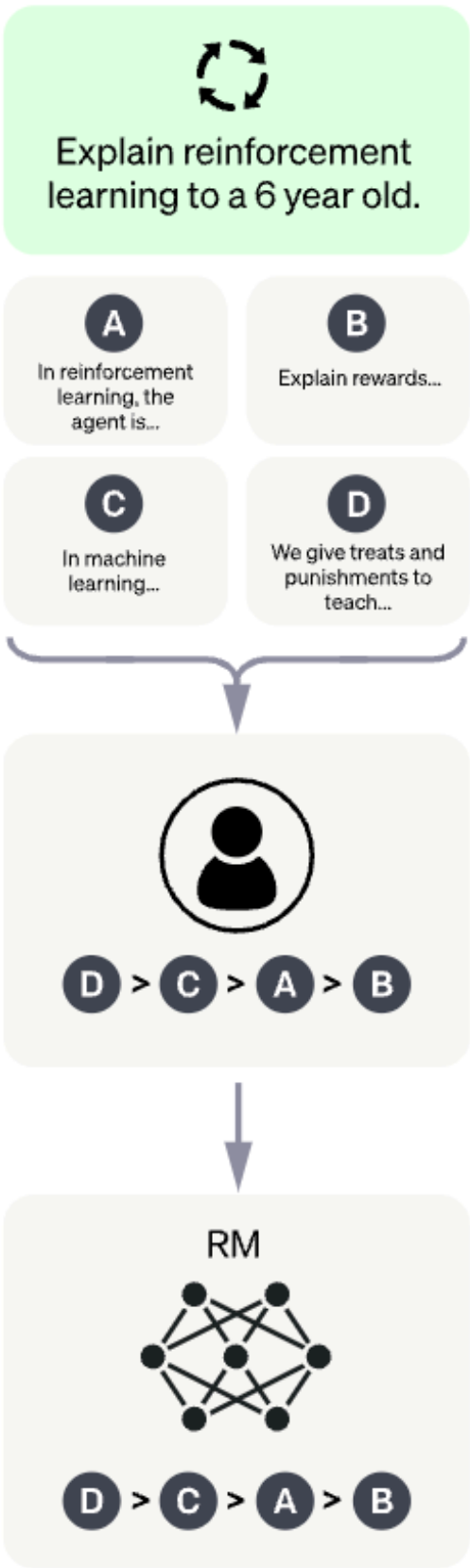


# Examples: Alignment of ChatGPT

Step 2

Collect comparison data and train a reward model.

A prompt and several model outputs are sampled.



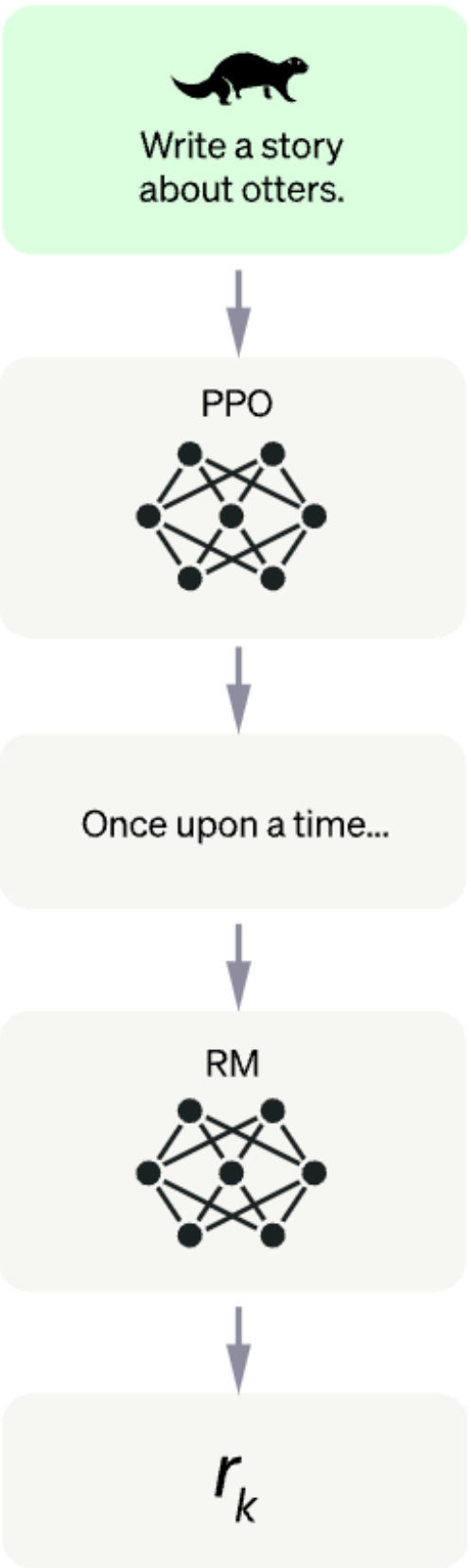
A labeler ranks the outputs from best to worst.

This data is used to train our reward model.

Step 3

Optimize a policy against the reward model using the PPO reinforcement learning algorithm.

A new prompt is sampled from the dataset.



The PPO model is initialized from the supervised policy.

The policy generates an output.

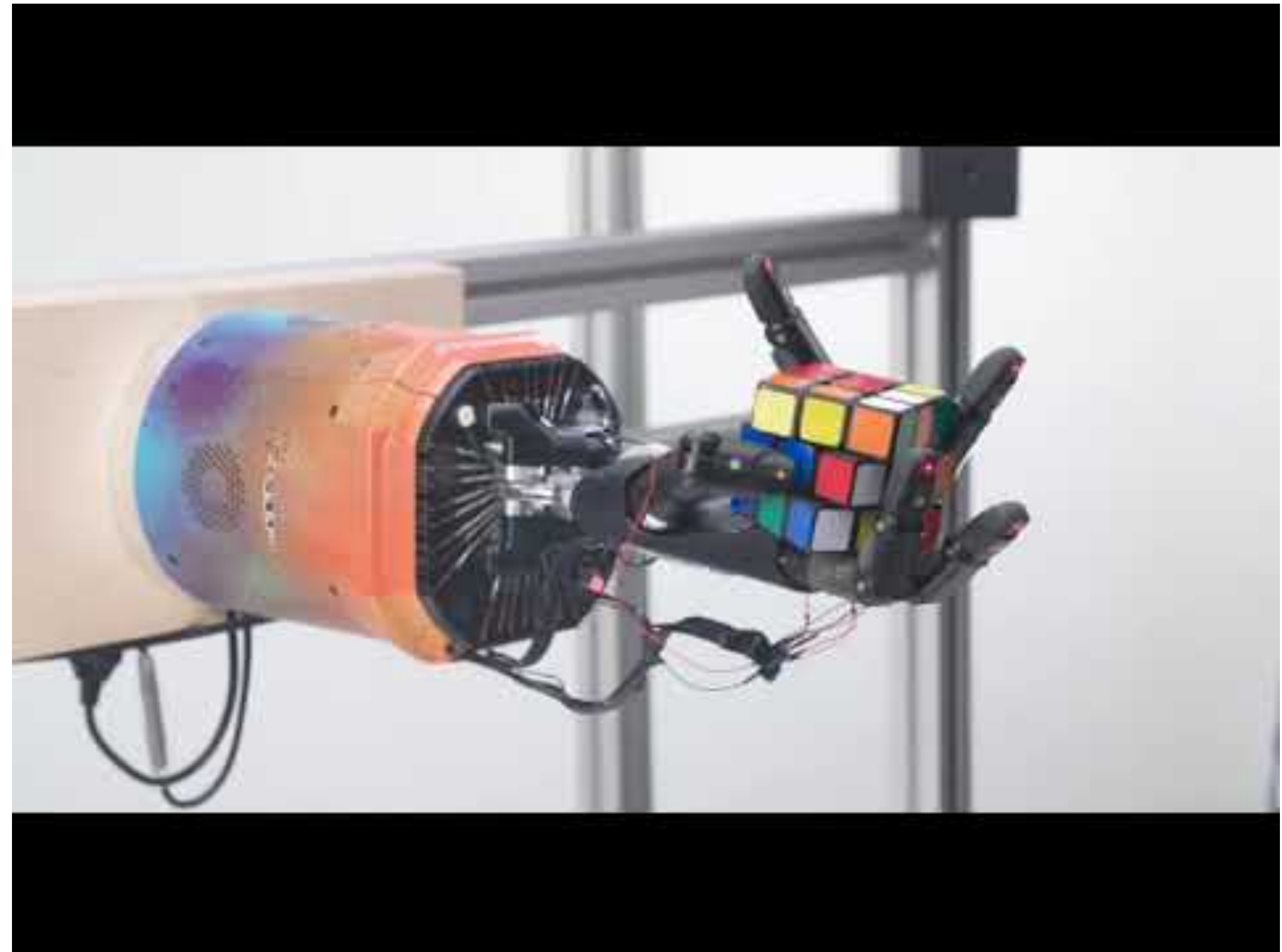
The reward model calculates a reward for the output.

The reward is used to update the policy using PPO.

ChatGPT

# Examples: Robot manipulation

- PPO
- Trained entirely in Sim



# Next time

- Model-based RL