

Introduction to Reinforcement Learning and Policy-Gradients with Tensor-Flow

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Slides adapted from (Berkeley [CS 294](#): Deep Reinforcement Learning by Sergey Levine)

Why Reinforcement Learning?

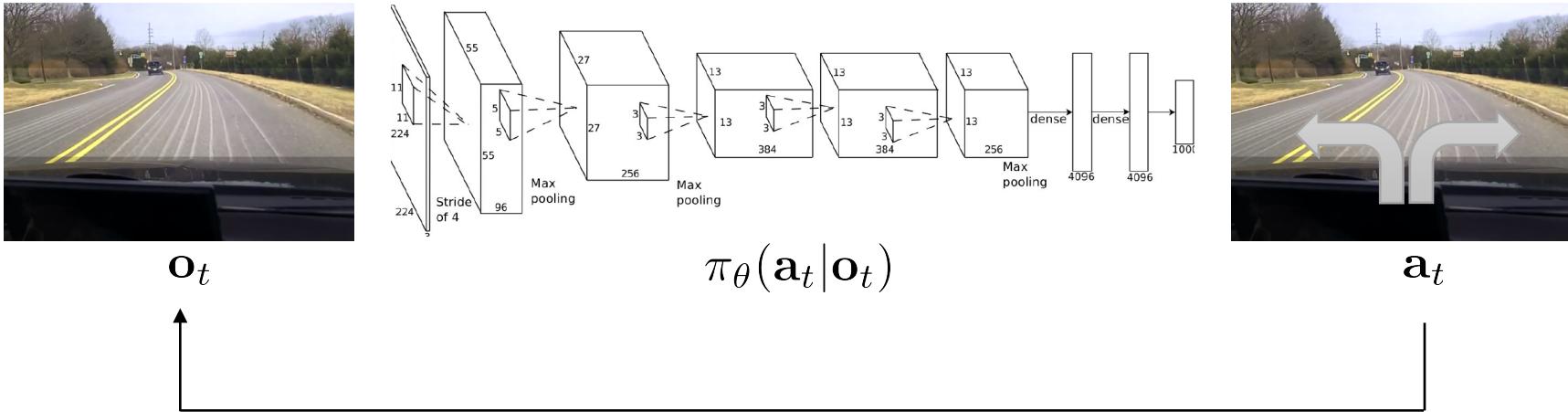


Today's Lecture

1. Definition of reinforcement learning problem
 2. Brief overview of RL algorithm types
 3. Introduction to policy gradient algorithms
 4. Implementation of policy gradient algorithms in TF
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- Goals:
 - Understand definitions & notation
 - Get an overview of different reinforcement learning algorithms
 - Understand how the policy gradient RL-algorithm can be implemented in TF

Definitions

Terminology & notation



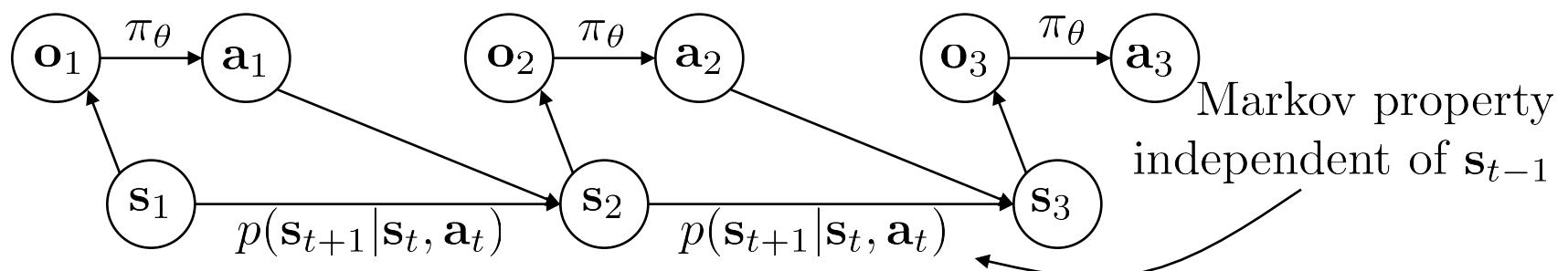
\mathbf{s}_t – state

\mathbf{o}_t – observation

\mathbf{a}_t – action

$\pi_\theta(\mathbf{a}_t | \mathbf{o}_t)$ – policy

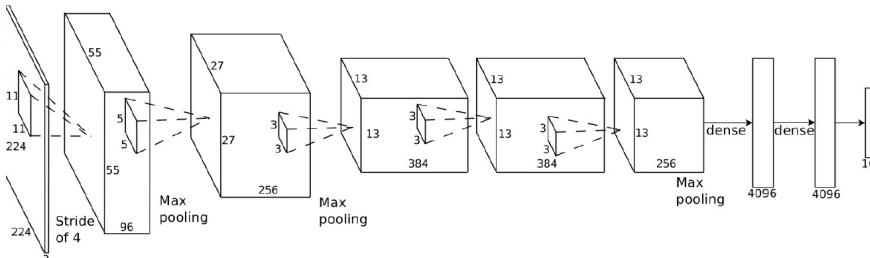
$\pi_\theta(\mathbf{a}_t | \mathbf{s}_t)$ – policy (fully observed)



Reward functions



\mathbf{o}_t



$$\pi_\theta(\mathbf{a}_t | \mathbf{o}_t)$$



\mathbf{a}_t

which action is better or worse?

$r(\mathbf{s}, \mathbf{a})$: reward function

tells us which states and actions are better



high reward

\mathbf{s} , \mathbf{a} , $r(\mathbf{s}, \mathbf{a})$, and $p(\mathbf{s}'|\mathbf{s}, \mathbf{a})$ define
Markov decision process



low reward

Definitions

partially observed Markov decision process $\mathcal{M} = \{\mathcal{S}, \mathcal{A}, \mathcal{O}, \mathcal{T}, \mathcal{E}, r\}$

\mathcal{S} – state space states $s \in \mathcal{S}$ (discrete or continuous)

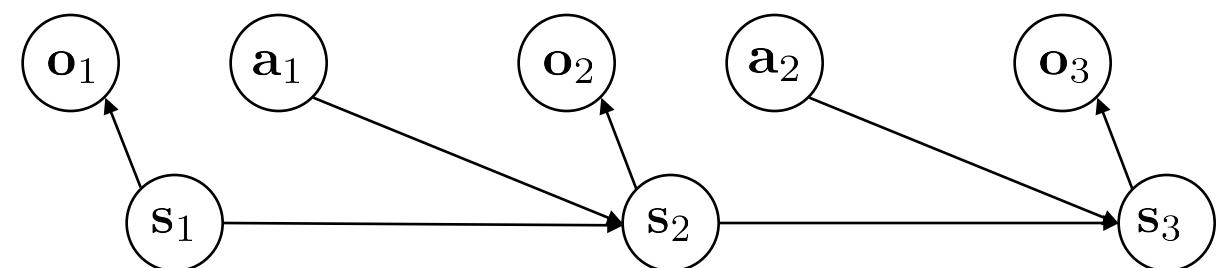
\mathcal{A} – action space actions $a \in \mathcal{A}$ (discrete or continuous)

\mathcal{O} – observation space observations $o \in \mathcal{O}$ (discrete or continuous)

\mathcal{T} – transition operator (like before)

\mathcal{E} – emission probability $p(o_t | s_t)$

r – reward function $r : \mathcal{S} \times \mathcal{A} \rightarrow \mathbb{R}$



Expectations and stochastic systems

$$\theta^* = \arg \max_{\theta} E_{(\mathbf{s}, \mathbf{a}) \sim p_{\theta}(\mathbf{s}, \mathbf{a})}[r(\mathbf{s}, \mathbf{a})]$$

infinite horizon case

$$\theta^* = \arg \max_{\theta} \sum_{t=1}^T E_{(\mathbf{s}_t, \mathbf{a}_t) \sim p_{\theta}(\mathbf{s}_t, \mathbf{a}_t)}[r(\mathbf{s}_t, \mathbf{a}_t)]$$

finite horizon case

In RL, we almost always care about *expectations*



$r(\mathbf{s}, \mathbf{a})$ – *not smooth*

ψ – probability of falling

$E_{(\mathbf{s}, \mathbf{a}) \sim p_{\psi}(\mathbf{s}, \mathbf{a})}[r(\mathbf{s}, \mathbf{a})]$ – *smooth* in ψ !

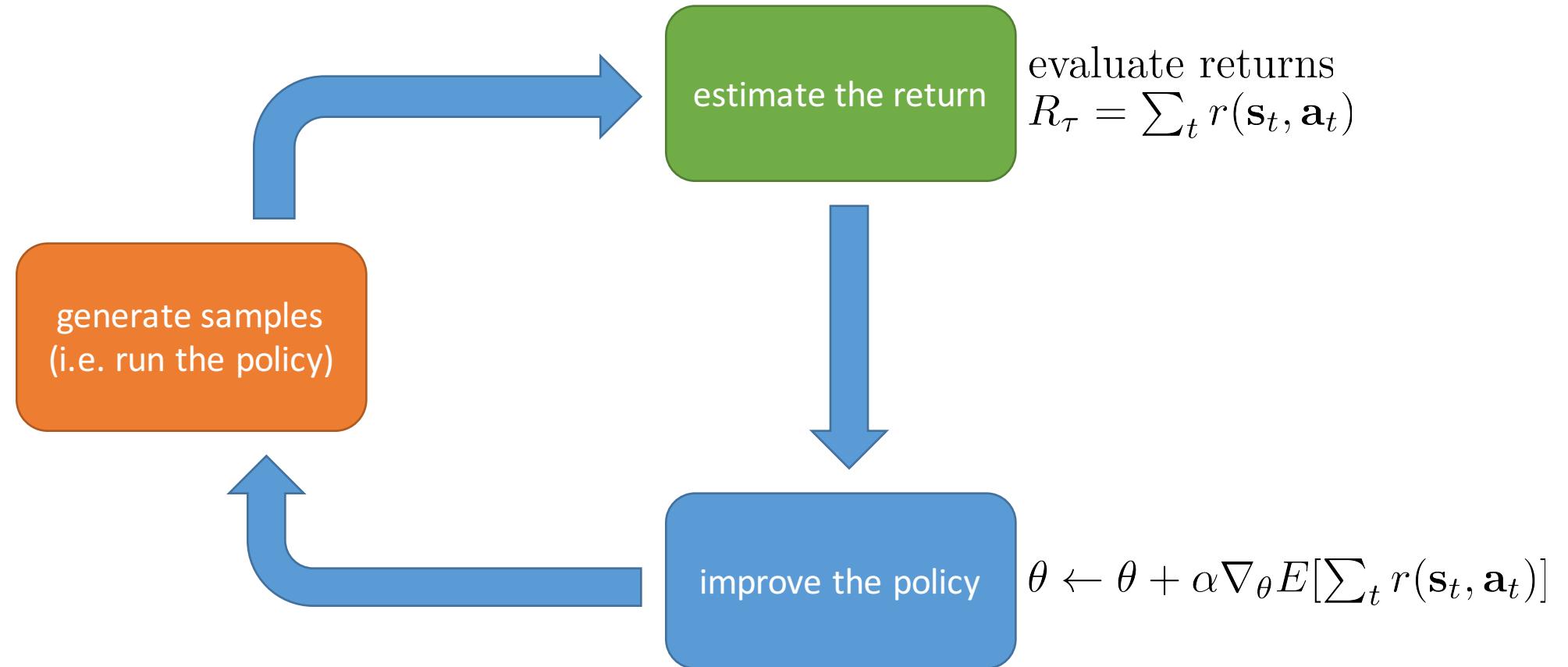
Algorithms

Types of RL algorithms

$$\theta^* = \arg \max_{\theta} E_{\tau \sim p_{\theta}(\tau)} \left[\sum_t r(s_t, a_t) \right]$$

- Policy gradients: directly differentiate the above objective
- Value-based: estimate value function or Q-function of the optimal policy (no explicit policy)
- Actor-critic: estimate value function or Q-function of the current policy, use it to improve policy
- Model-based RL: estimate the transition model, and then...
 - Use it for planning (no explicit policy)
 - Use it to improve a policy
 - Something else

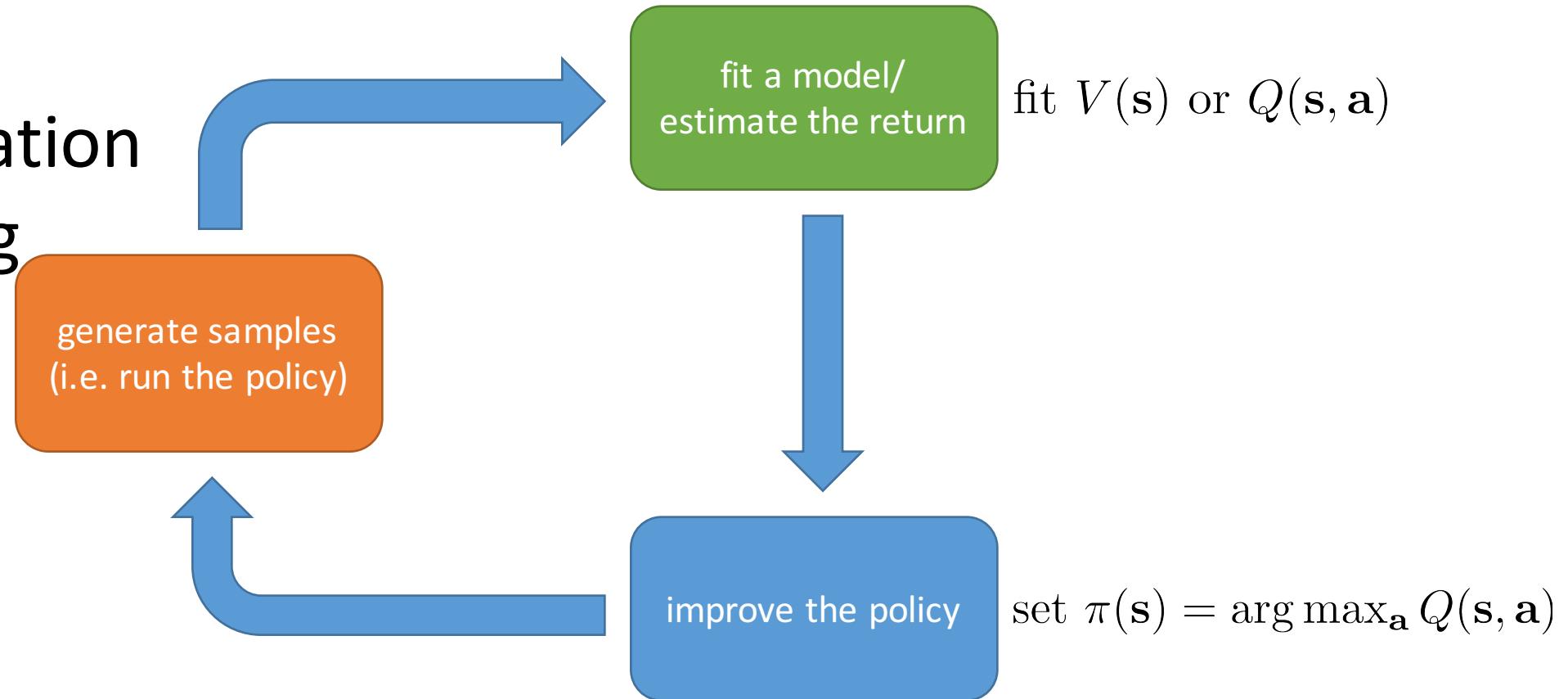
Direct policy gradients



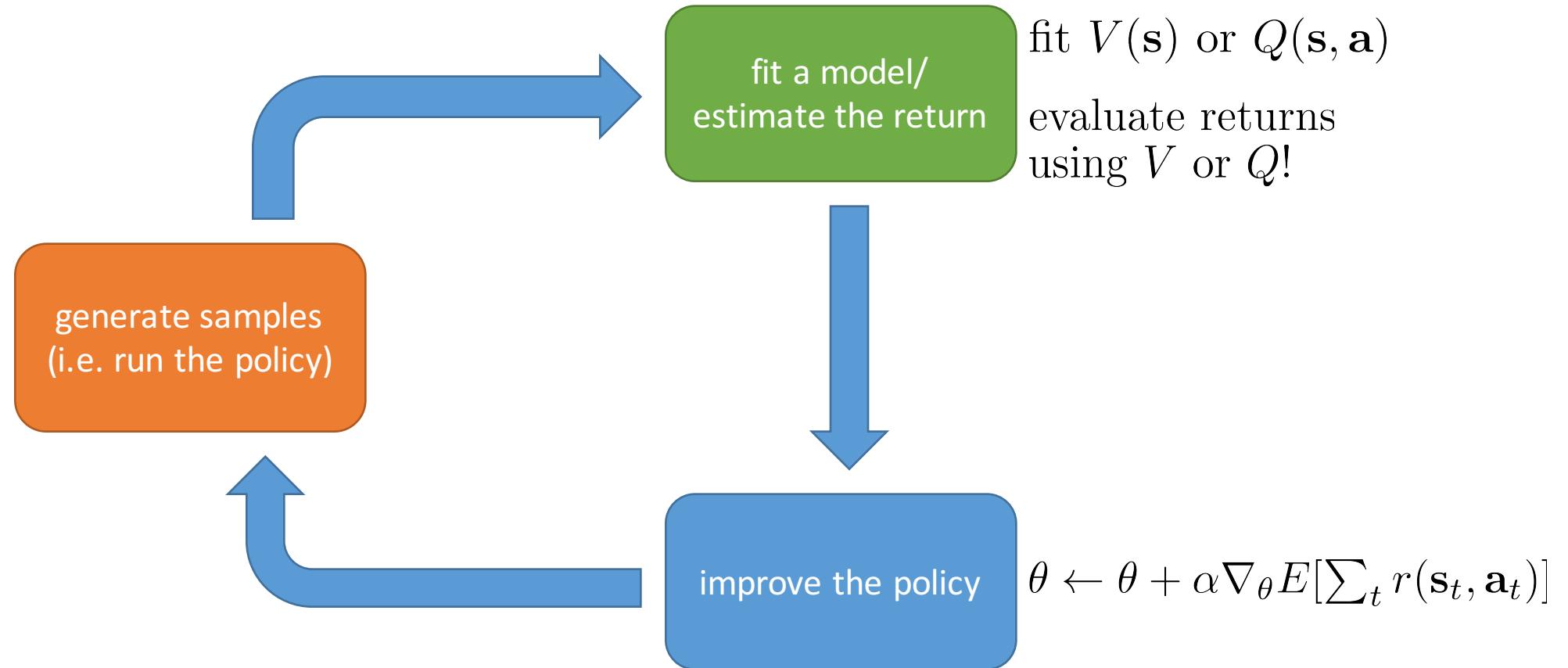
Value function based algorithms

Examples:

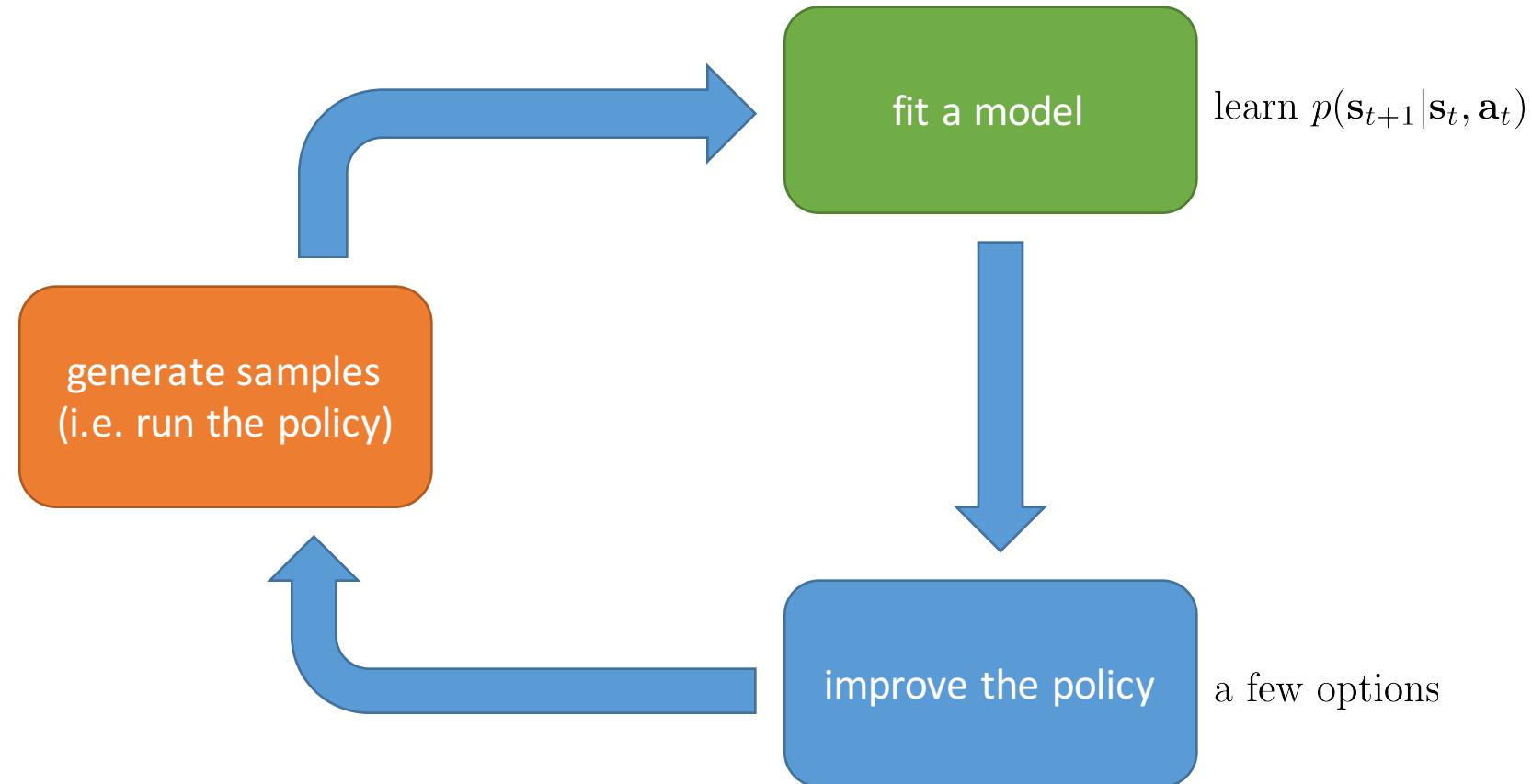
- Value-Iteration
- Q-Learning
- DQN



Actor-critic: value functions + policy gradients

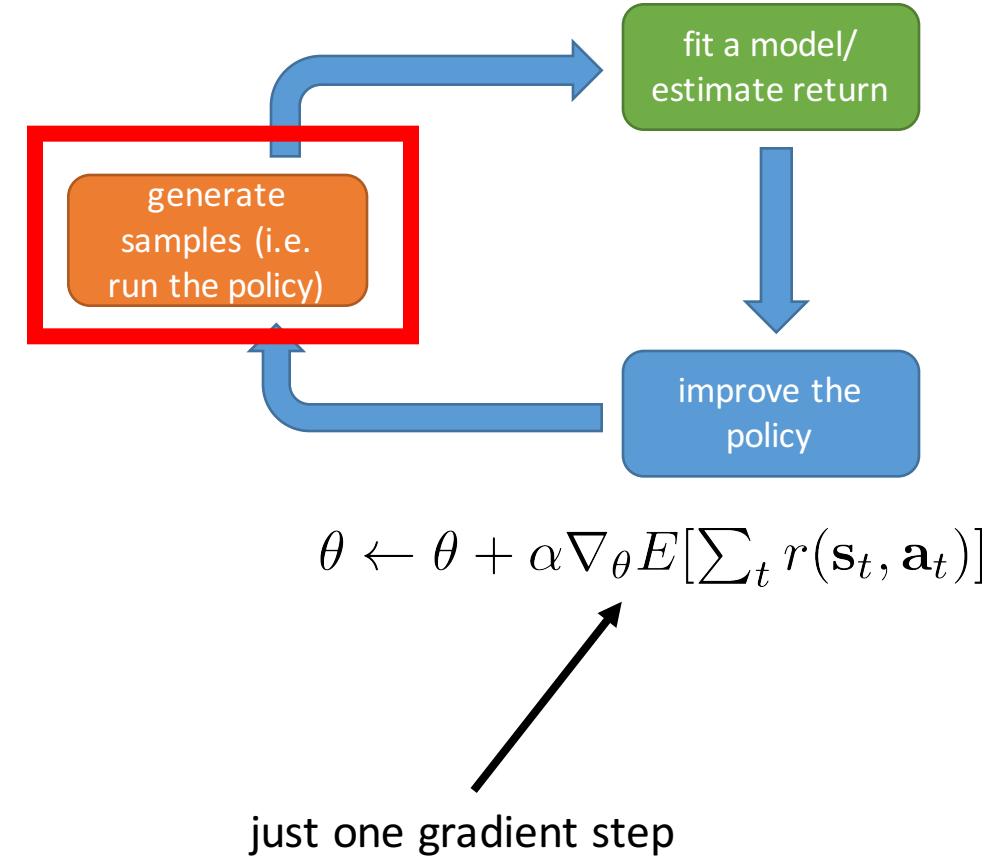


Model-based RL algorithms

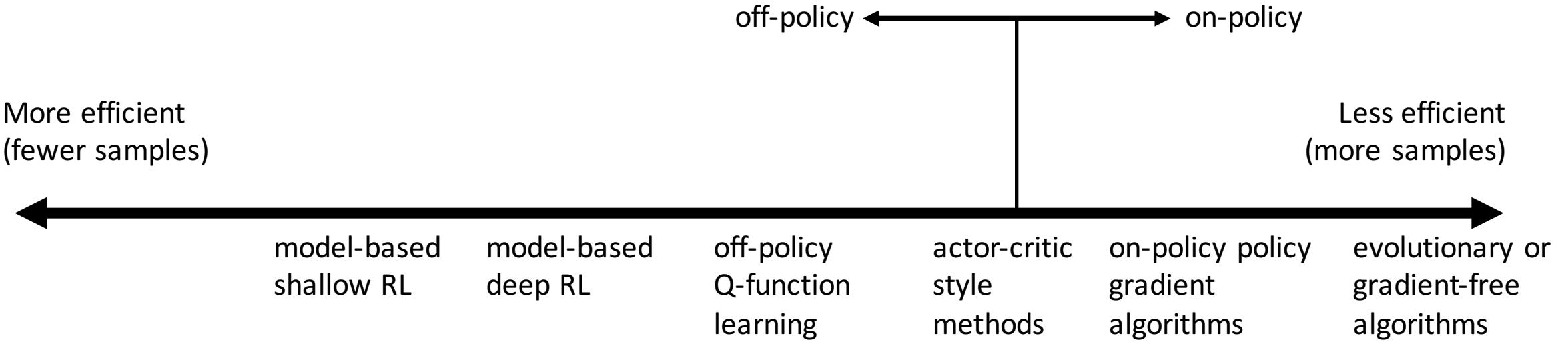


Comparison: sample efficiency

- Sample efficiency = how many samples do we need to get a good policy?
- Most important question: is the algorithm *off policy*?
 - Off policy: able to improve the policy without generating new samples from that policy
 - On policy: each time the policy is changed, even a little bit, we need to generate new samples



Comparison: sample efficiency



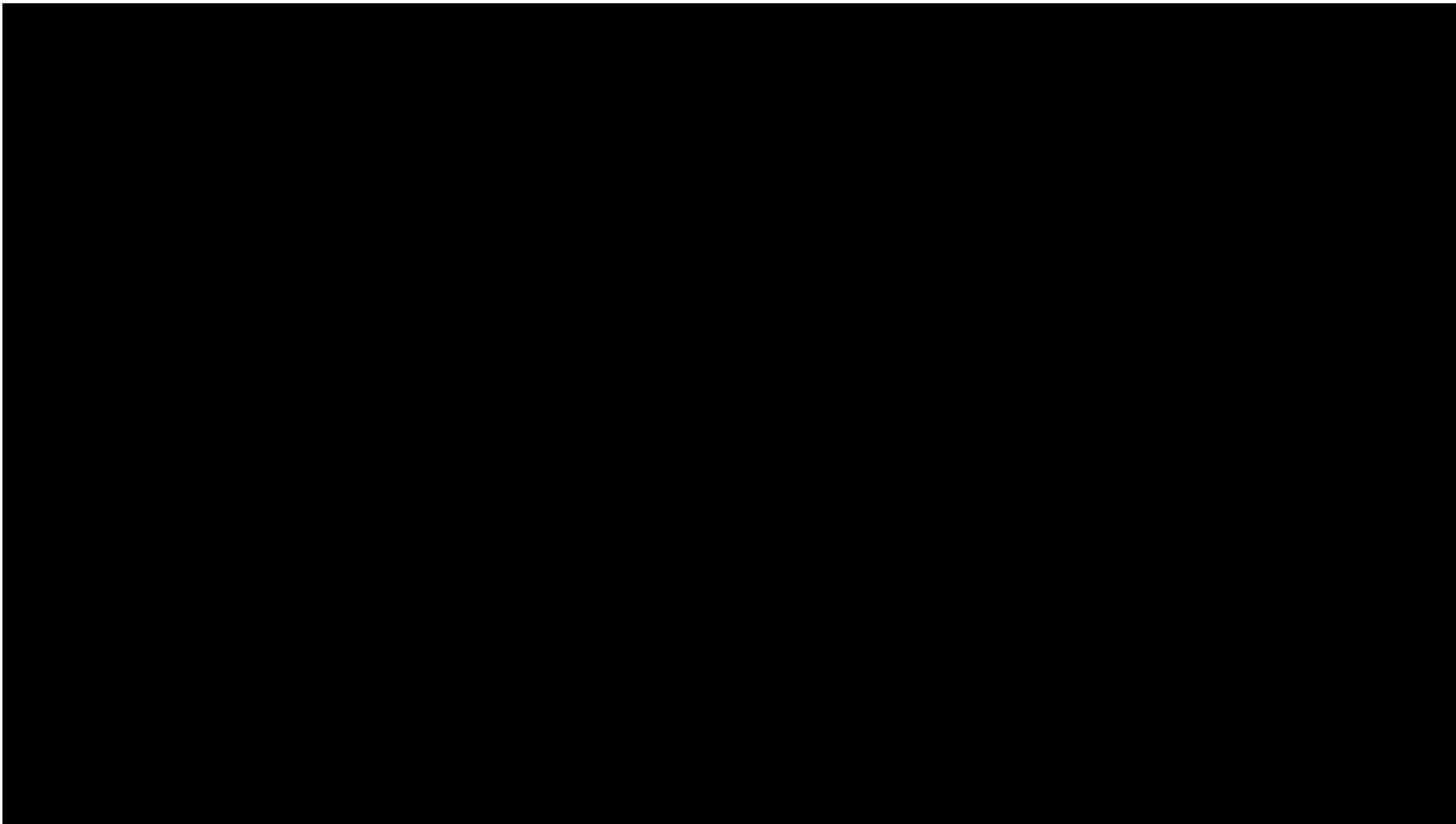
Why would we use a *less* efficient algorithm?

Wall clock time is not the same as efficiency!

Comparison: stability and ease of use

- Value function fitting
 - At best, minimizes error of fit (“Bellman error”)
 - Not the same as expected reward
 - At worst, doesn’t optimize anything
 - Many popular deep RL value fitting algorithms are not guaranteed to converge to *anything* in the nonlinear case
- Model-based RL
 - Model minimizes error of fit
 - This will converge
 - No guarantee that better model = better policy
- Policy gradient
 - The only one that actually performs gradient descent (ascent) on the true objective

Example: Robotic Manipulation with value function based algorithm



For detail see the Normalized Advantage Function ([NAF](#)) algorithm

Introduction to Policy Gradients

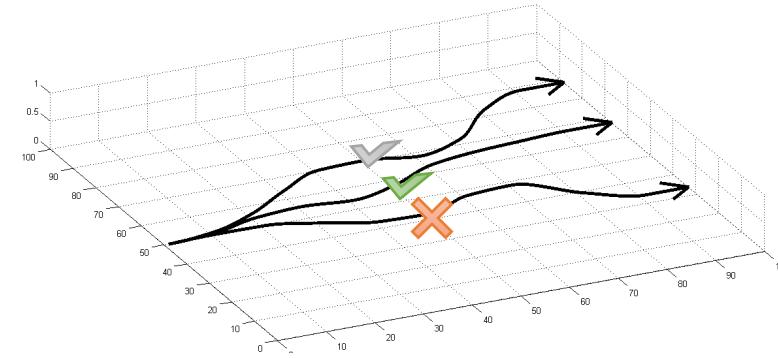
Evaluating the objective

$$\theta^* = \arg \max_{\theta} E_{\tau \sim p_{\theta}(\tau)} \left[\sum_t r(\mathbf{s}_t, \mathbf{a}_t) \right]$$

$J(\theta)$

$$J(\theta) = E_{\tau \sim p_{\theta}(\tau)} \left[\sum_t r(\mathbf{s}_t, \mathbf{a}_t) \right] \approx \frac{1}{N} \sum_i \sum_t r(\mathbf{s}_{i,t}, \mathbf{a}_{i,t})$$

sum over samples from π_{θ}



Direct policy differentiation

$$\theta^* = \arg \max_{\theta} E_{\tau \sim p_{\theta}(\tau)} \left[\underbrace{\sum_t r(\mathbf{s}_t, \mathbf{a}_t)}_{J(\theta)} \right]$$

a convenient identity

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

$$J(\theta) = E_{\tau \sim \pi_{\theta}(\tau)} [r(\tau)] = \int \pi_{\theta}(\tau) r(\tau) d\tau$$
$$\sum_{t=1}^T r(\mathbf{s}_t, \mathbf{a}_t)$$

$$\nabla_{\theta} J(\theta) = \int \underline{\nabla_{\theta} \pi_{\theta}(\tau)} r(\tau) d\tau = \int \underline{\pi_{\theta}(\tau) \nabla_{\theta} \log \pi_{\theta}(\tau)} r(\tau) d\tau = E_{\tau \sim \pi_{\theta}(\tau)} [\nabla_{\theta} \log \pi_{\theta}(\tau) r(\tau)]$$

Evaluating the policy gradient

recall: $J(\theta) = E_{\tau \sim p_\theta(\tau)} \left[\sum_t r(\mathbf{s}_t, \mathbf{a}_t) \right] \approx \frac{1}{N} \sum_i \sum_t r(\mathbf{s}_{i,t}, \mathbf{a}_{i,t})$

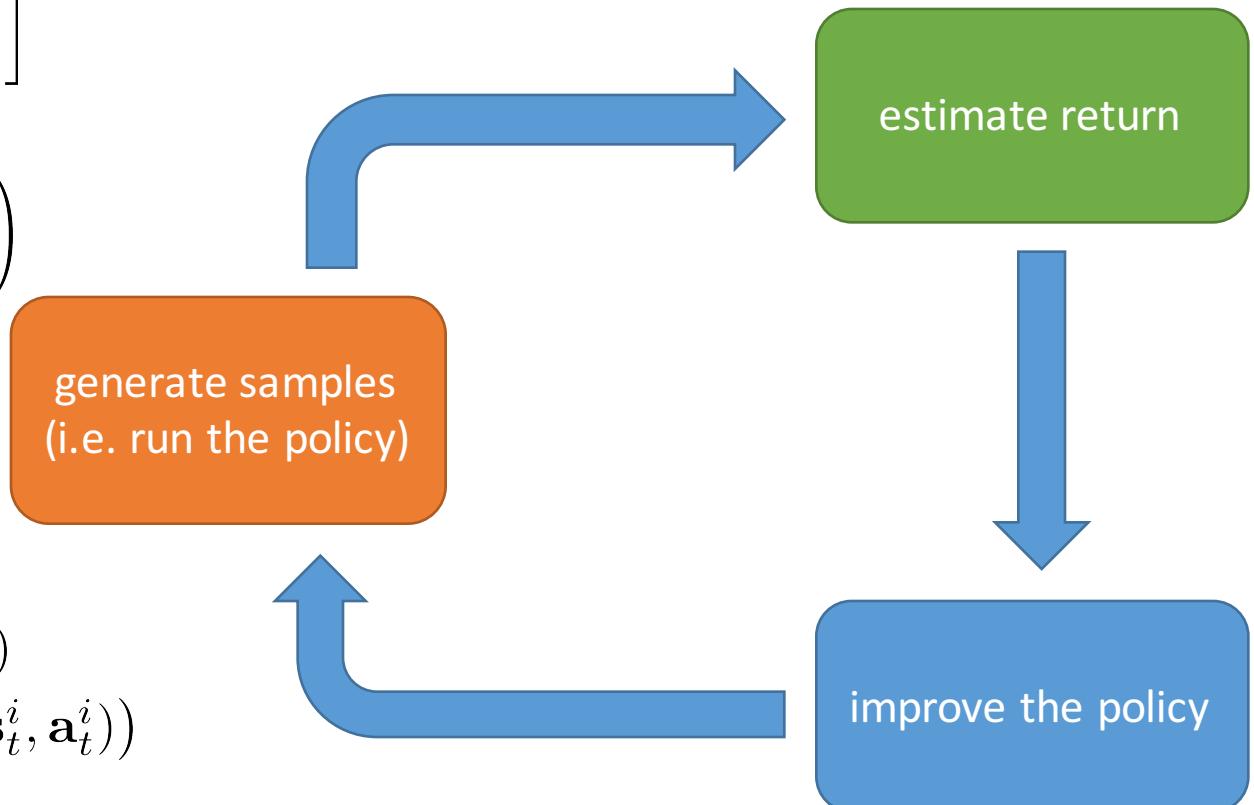
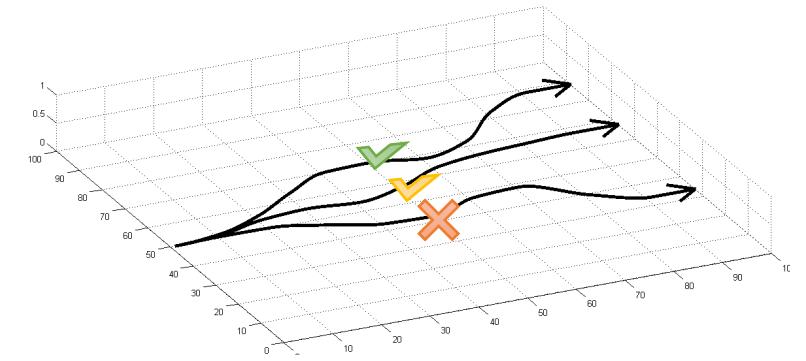
$$\nabla_\theta J(\theta) = E_{\tau \sim \pi_\theta(\tau)} \left[\left(\sum_{t=1}^T \nabla_\theta \log \pi_\theta(\mathbf{a}_t | \mathbf{s}_t) \right) \left(\sum_{t=1}^T r(\mathbf{s}_t, \mathbf{a}_t) \right) \right]$$

$$\nabla_\theta J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left(\sum_{t=1}^T \nabla_\theta \log \pi_\theta(\mathbf{a}_{i,t} | \mathbf{s}_{i,t}) \right) \left(\sum_{t=1}^T r(\mathbf{s}_{i,t}, \mathbf{a}_{i,t}) \right)$$

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

REINFORCE algorithm:

- 1. sample $\{\tau^i\}$ from $\pi_\theta(\mathbf{a}_t | \mathbf{s}_t)$ (run the policy)
- 2. $\nabla_\theta J(\theta) \approx \sum_i \left(\sum_t \nabla_\theta \log \pi_\theta(\mathbf{a}_t^i | \mathbf{s}_t^i) \right) \left(\sum_t r(\mathbf{s}_t^i, \mathbf{a}_t^i) \right)$
- 3. $\theta \leftarrow \theta + \alpha \nabla_\theta J(\theta)$



Example: Gaussian policies

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left(\sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t}) \right) \left(\sum_{t=1}^T r(\mathbf{s}_{i,t}, \mathbf{a}_{i,t}) \right)$$

example: $\pi_{\theta}(\mathbf{a}_t | \mathbf{s}_t) = \mathcal{N}(f_{\text{neural network}}(\mathbf{s}_t); \Sigma)$

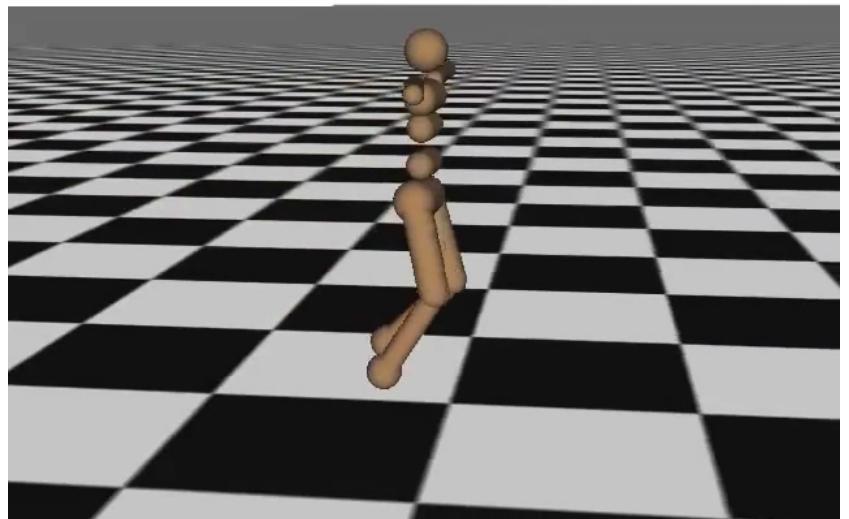
$$\log \pi_{\theta}(\mathbf{a}_t | \mathbf{s}_t) = -\frac{1}{2} \|f(\mathbf{s}_t) - \mathbf{a}_t\|_{\Sigma}^2 + \text{const}$$

$$\nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_t | \mathbf{s}_t) = -\frac{1}{2} \Sigma^{-1} (f(\mathbf{s}_t) - \mathbf{a}_t) \frac{df}{d\theta}$$

REINFORCE algorithm:

- 1. sample $\{\tau^i\}$ from $\pi_{\theta}(\mathbf{a}_t | \mathbf{s}_t)$ (run it on the robot)
- 2. $\nabla_{\theta} J(\theta) \approx \sum_i (\sum_t \nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_t^i | \mathbf{s}_t^i)) (\sum_t r(\mathbf{s}_t^i, \mathbf{a}_t^i))$
- 3. $\theta \leftarrow \theta + \alpha \nabla_{\theta} J(\theta)$

Iteration 2000



What did we just do?

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left(\sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t}) \right) \left(\sum_{t=1}^T r(\mathbf{s}_{i,t}, \mathbf{a}_{i,t}) \right)$$

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \underbrace{\sum_{t=1}^T \nabla_{\theta} \log_{\theta} \pi_{\theta}(\tau_i)}_{\sum_{t=1}^T \nabla_{\theta} \log_{\theta} \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t})} r(\tau_i)$$

maximum likelihood: $\nabla_{\theta} J_{\text{ML}}(\theta) \approx \frac{1}{N} \sum_{i=1}^N \nabla_{\theta} \log \pi_{\theta}(\tau_i)$

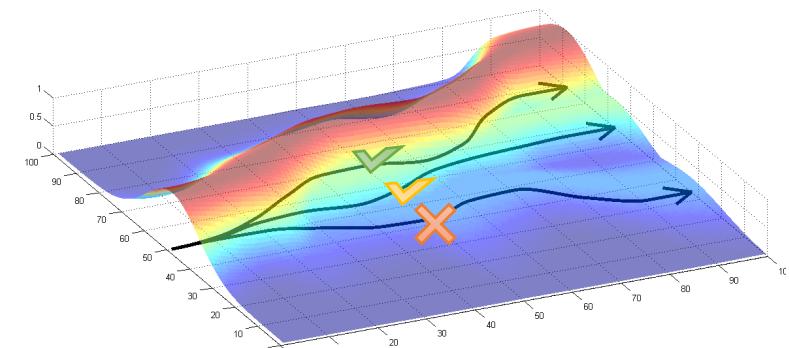
good stuff is made more likely

bad stuff is made less likely

simply formalizes the notion of “trial and error”!

REINFORCE algorithm:

- 1. sample $\{\tau^i\}$ from $\pi_{\theta}(\mathbf{a}_t | \mathbf{s}_t)$ (run it on the robot)
- 2. $\nabla_{\theta} J(\theta) \approx \sum_i (\sum_t \nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_t^i | \mathbf{s}_t^i)) (\sum_t r(\mathbf{s}_t^i, \mathbf{a}_t^i))$
- 3. $\theta \leftarrow \theta + \alpha \nabla_{\theta} J(\theta)$



Reducing variance

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \left(\sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t}) \right) \left(\sum_{t=1}^T r(\mathbf{s}_{i,t}, \mathbf{a}_{i,t}) \right)$$

What you do now does **not** affect the reward of the past!

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t}) \underbrace{\left(\sum_{t'=1}^T r(\mathbf{s}_{i,t''}, \mathbf{a}_{i,t''}) \right)}_{\text{reward to go}}$$

“reward to go”

$$\hat{Q}_{i,t}$$

Baselines

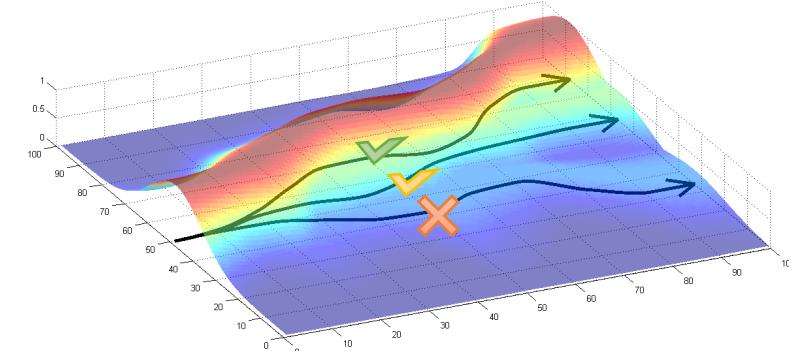
a convenient identity

$$\pi_\theta(\tau) \nabla_\theta \log \pi_\theta(\tau) = \nabla_\theta \pi_\theta(\tau)$$

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

$$b = \frac{1}{N} \sum_{i=1}^N r(\tau)$$

but... are we *allowed* to do that??



$$E[\nabla_\theta \log \pi_\theta(\tau) b] = \int \pi_\theta(\tau) \nabla_\theta \log \pi_\theta(\tau) b d\tau = \int \nabla_\theta \pi_\theta(\tau) b d\tau = b \nabla_\theta \int \pi_\theta(\tau) d\tau = b \nabla_\theta 1 = 0$$

subtracting a baseline is *unbiased* in expectation!

average reward is *not* the best baseline, but it's pretty good!

Implementation of Policy Gradients

Policy gradient with automatic differentiation

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \underline{\nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t}) \hat{Q}_{i,t}}$$



pretty inefficient to compute these explicitly!

How can we compute policy gradients with automatic differentiation?

We need a graph such that its gradient is the policy gradient!

maximum likelihood: $\nabla_{\theta} J_{\text{ML}}(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t})$ $J_{\text{ML}}(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \log \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t})$

Just implement “pseudo-loss” as a weighted maximum likelihood:

$$\tilde{J}(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \log \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t}) \hat{Q}_{i,t}$$



cross entropy (discrete) or squared error (Gaussian)

Policy gradient with automatic differentiation

Pseudocode example (with discrete actions):

Policy gradient:

```
# Given:  
# actions - (N*T) x Da tensor of actions  
# states - (N*T) x Ds tensor of states  
# rew_to_go - (N*T) x 1 tensor of estimated reward to go  
# Build the graph:  
logits = policy.predictions(states) # This should return (N*T) x Da tensor of action logits  
negative_likelihoods = tf.nn.softmax_cross_entropy_with_logits(labels=actions, logits=logits)  
weighted_negative_likelihoods = tf.multiply(negative_likelihoods, rew_to_go)  
loss = tf.reduce_mean(weighted_negative_likelihoods)  
gradients = loss.gradients(loss, variables)
```

$$\tilde{J}(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \log \pi_\theta(\mathbf{a}_{i,t} | \mathbf{s}_{i,t}) \hat{Q}_{i,t}$$

Reward to go

Policy gradient with automatic differentiation

Pseudocode example (with discrete actions):

Policy gradient:

```
# Given:  
# actions - (N*T) x Da tensor of actions  
# states - (N*T) x Ds tensor of states  
# rew_to_go - (N*T) x 1 tensor of estimated reward to go  
# Build the graph:  
mean = policy.predictions(states) # This should return (N*T) x Da tensor of action logits  
negative_likelihoods = gaussian_log_prob(sy_ac_na, mean, sy_logstd)  
weighted_negative_likelihoods = tf.multiply(negative_likelihoods, rew_to_go)  
loss = tf.reduce_mean(weighted_negative_likelihoods)  
gradients = loss.gradients(loss, variables)
```

$$\log \pi_{\theta}(\mathbf{a}_t | \mathbf{s}_t) = -\frac{1}{2} \|f(\mathbf{s}_t) - \mathbf{a}_t\|_{\Sigma}^2 + \text{const}$$

$$\tilde{J}(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \log \pi_{\theta}(\mathbf{a}_{i,t} | \mathbf{s}_{i,t}) \hat{Q}_{i,t}$$

Reward to go

Policy gradient in practice

- Remember that the gradient has high variance
 - This isn't the same as supervised learning!
 - Gradients will be really noisy!
- Consider using much larger batches
- Tweaking learning rates is very hard
 - Adaptive step size rules like ADAM can be OK-ish
 - There exist algorithms that adjust the gradient stepsize to obtain more stability, such as Trust-Region Policy Optimization (TRPO) and Proximal Policy Optimization (PPO)

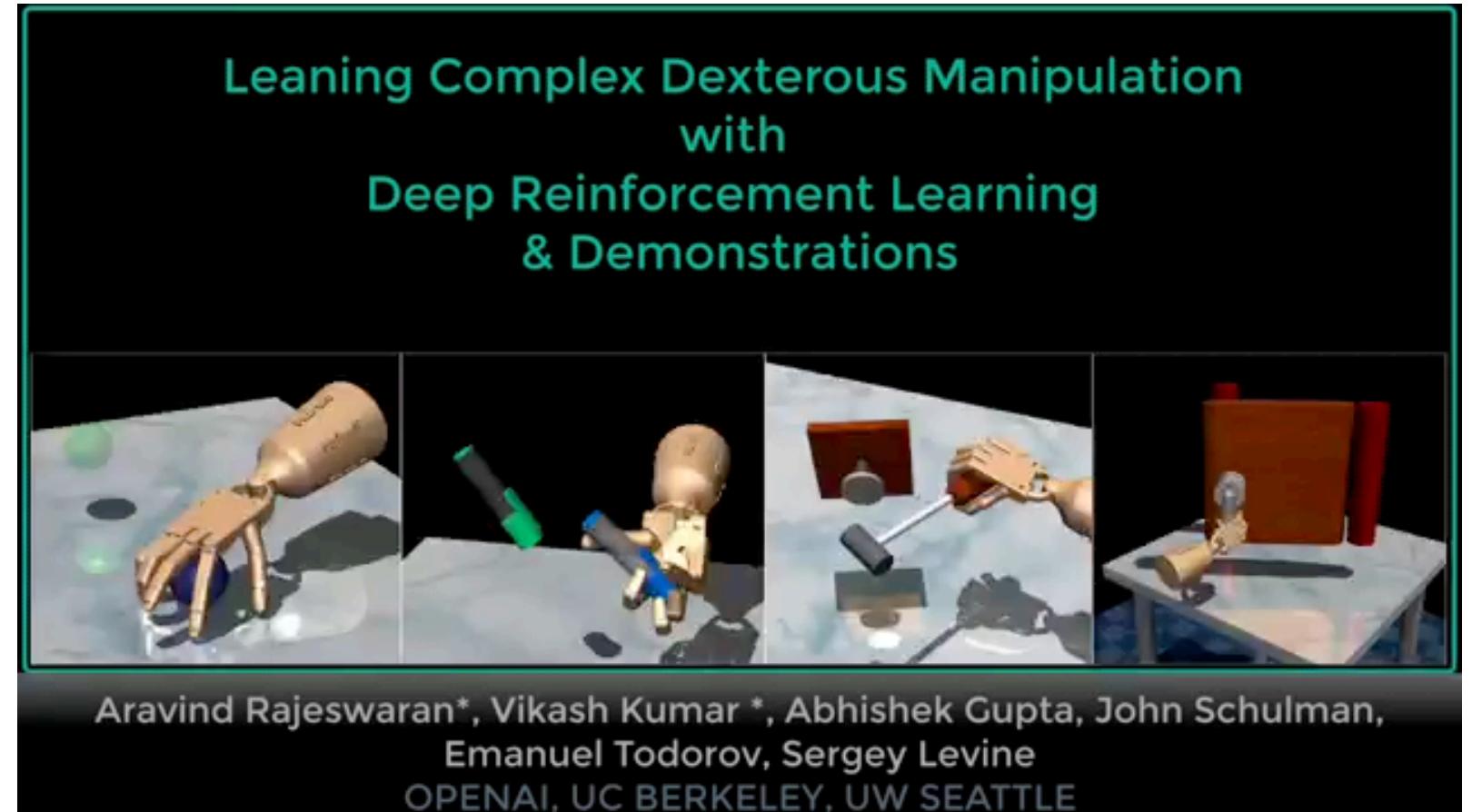
Suggested Project

- Implement policy gradient as in [homework 2 of CS 294: DeepRL, Fall 2017](#)
 - Vanilla policy gradient algorithm in Tensorflow
 - Add baseline for variance reduction
 - Agents trained for Inverted Pendulum and Cheetah environments
(for Cheetah Mujoco physics engine necessary, 30 day trial license available)
 - Most of the code is prepared, you only need to fill in a couple of blanks

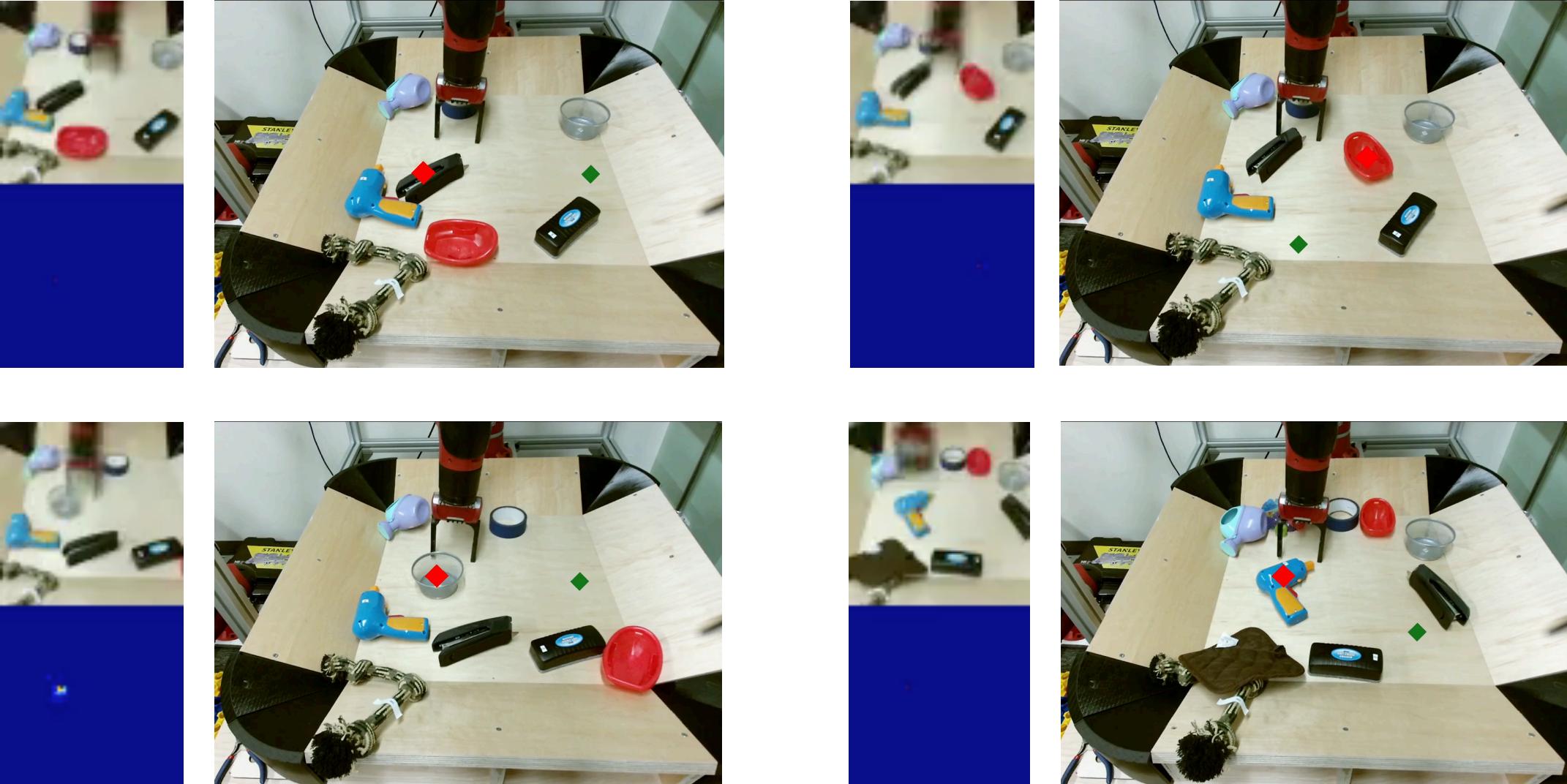
The material was prepared by [Abhishek Gupta](#) and Josh Aicham.

Example: trust region policy optimization, policies initialized from demonstration

- Natural gradient with automatic step adjustment
- Discrete and continuous actions
- Using a small number of demonstrations to overcome exploration problem.



Beyond RL: Self-supervised Learning with Video-Prediction and Sampling Based Planning



Policy gradients suggested lectures and readings

- Lectures online: Berkeley [CS 294](#), Course at UCL by David Silver
- Classic papers
 - Williams (1992). Simple statistical gradient-following algorithms for connectionist reinforcement learning: introduces REINFORCE algorithm
 - Baxter & Bartlett (2001). Infinite-horizon policy-gradient estimation: temporally decomposed policy gradient (not the first paper on this! see actor-critic section later)
 - Peters & Schaal (2008). Reinforcement learning of motor skills with policy gradients: very accessible overview of optimal baselines and natural gradient
- Deep reinforcement learning policy gradient papers
 - Levine & Koltun (2013). Guided policy search: deep RL with importance sampled policy gradient (unrelated to later discussion of guided policy search)
 - Schulman, L., Moritz, Jordan, Abbeel (2015). Trust region policy optimization: deep RL with natural policy gradient and adaptive step size
 - Schulman, Wolski, Dhariwal, Radford, Klimov (2017). Proximal policy optimization algorithms: deep RL with importance sampled policy gradient