

# Deep Q Networks

# Deep Reinforcement Learning

University of Cambridge

• Review

- Review
- State of the field

- Review
- State of the field
- Implement Deep Q Networks (DQN) (Mnih et al.)

- Review
- State of the field
- Implement Deep Q Networks (DQN) (Mnih et al.)

#### **Review**:

• Finished up MDPs

- Finished up MDPs
- The return and optimal policies

- Finished up MDPs
- The return and optimal policies
- Deep Q learning

- Finished up MDPs
- The discounted return and optimal policies
- Deep Q learning

- Finished up MDPs
- The return and optimal policies
- Deep Q learning

#### The **discounted return** (G)

$$G = \sum_{t=0}^{\infty} \gamma^t R(s_{t+1}) = \sum_{t=0}^{\infty} \gamma^t r_t$$

#### The discounted return (G)

$$G = \sum_{t=0}^{\infty} \gamma^t R(s_{t+1}) = \sum_{t=0}^{\infty} \gamma^t r_t$$

With a reward of 1 at each timestep and  $\gamma = 0.9$ 

$$G = \sum_{t=0}^{\infty} \gamma^t r_t = 1 + 0.9 + 0.81 + \ldots = \frac{1}{1 - \gamma} = 10$$

#### The expected discounted return ( $G_{\pi}$ )

$$G_{\pi} = \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t ~|~ a_t \sim \pi(s_t)\right]$$

$$G_{\pi} = \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

Transitions and policy are stochastic. Consider uncertainty in the reward.

$$G_{\pi} = \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

Transitions and policy are stochastic. Consider uncertainty in the reward.

$$r_t \sim \left[\underbrace{\frac{R(s_{t+1})}{T(s_{t+1} \mid s_t, a_t)}}_{\text{reward fn}} \overline{T(s_{t+1} \mid s_t, a_t)} \underbrace{\pi(a_t \mid s_t)}_{\text{action probs}}\right]$$



The **expectation** turns that distribution into a single number. This tells us what reward to expect "on average"

$$\mathbb{E}[r_t] = \int_{s_{t+1}} \int_A \underbrace{\frac{R(s_{t+1})}_{\text{reward fn}} \overline{T(s_{t+1} \mid s_t, a_t)}}_{\text{reward fn}} \underbrace{\frac{\pi(a_t \mid s_t)}{\pi(a_t \mid s_t)}}_{\text{action probs}}$$

$$G_{\pi} = \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t ~|~ a_t \sim \pi(s_t)\right]$$

With the expected discounted return, we can define the optimal policy

$$\pi_* = \max_{\pi} \ \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

- Finished up MDPs
- The return and optimal policies
- Deep Q learning

# The Plan:

- 1. Derive the value function  ${\cal V}$
- 2. Derive Q function from V
- 3. Figure out a behavior policy using  ${\cal Q}$
- 4. Learn to train Q

# The Plan:

- 1. Derive the value function V
- 2. Derive Q function from V
- 3. Figure out a behavior policy using  ${\cal Q}$
- 4. Learn to train Q

With the expected return

$$G_{\pi} = \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

We derived the Value Function  $(V_{\pi})$   $V_{\pi}: S \to \mathbb{R}$ 

With the expected return

$$G_{\pi} = \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

We derived the Value Function  $(V_{\pi})$   $V_{\pi}: S \to \mathbb{R}$ 

$$V_{\pi}(s_0) = \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

With the expected return

$$G_{\pi} = \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

We derived the Value Function  $(V_{\pi})$   $V_{\pi}: S \to \mathbb{R}$ 

$$V_{\pi}(s_0) = \mathbb{E}\left[\sum_{t=0}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

Difference between  $G_{\pi}$  and  $V_{\pi}$  is dependence on  $s_0$ 

# The Plan:

- 1. Derive the value function  ${\cal V}$
- 2. Derive Q function from V
- 3. Figure out a behavior policy using  ${\cal Q}$
- 4. Learn to train Q

Factor out first term from the return to introduce a dependence on  $a_0$ 

$$V_{\pi}(s_0, a_0) = \mathbb{E}[r_0 \mid a_0] + \mathbb{E}\left[\sum_{t=1}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

Factor out first term from the return to introduce a dependence on  $a_0$ 

$$V_{\pi}(s_0, a_0) = \mathbb{E}[r_0 \mid a_0] + \mathbb{E}\left[\sum_{t=1}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

When *V* depends on a specific action, we call it the **Q** function:

$$S \times A \to \mathbb{R}$$

$$Q_{\pi}(s_0, a_0) = \mathbb{E}[r_0 \mid a_0] + \mathbb{E}\left[\sum_{t=1}^{\infty} \gamma^t r_t \mid a_t \sim \pi(s_t)\right]$$

# The Plan:

- 1. Derive the value function  ${\cal V}$
- 2. Derive Q function from V
- 3. Figure out a behavior policy using Q
- 4. Learn to train Q

$$\pi_*(s) = \operatorname*{argmax}_{a \in A} Q_*(s,a)$$

**In English:** Compute Q value for all possible actions and pick the action with the biggest Q value. Repeat at each timestep.

# The Plan:

- 1. Derive the value function  ${\cal V}$
- 2. Derive Q function from V
- 3. Figure out a behavior policy using  ${\cal Q}$
- 4. Learn to train Q

$$Q(s,a) = r + \gamma \cdot \max_{\{a' \in A\}} Q(s',a')$$

$$Q(s,a) = r + \gamma \cdot \max_{\{a' \in A\}} Q(s',a')$$

This course is **Deep** RL, so we need to use a neural network, parameterized by  $\theta$ 

$$Q(s,a) = r + \gamma \cdot \max_{\{a' \in A\}} Q(s',a')$$

This course is **Deep** RL, so we need to use a neural network, parameterized by  $\theta$ 

$$Q(s,a,\theta) = r + \gamma \cdot \max_{\{a' \in A\}} Q(s',a',\theta)$$

$$Q(s,a) = r + \gamma \cdot \max_{\{a' \in A\}} Q(s',a')$$

This course is **Deep** RL, so we need to use a neural network, parameterized by  $\theta$ 

$$Q(s, a, \theta) = r + \gamma \cdot \max_{\{a' \in A\}} Q(s', a', \theta)$$

$$Q(s,a,\theta) - \left(r + \gamma \cdot \max_{\{a' \in A\}} Q(s',a',\theta)\right) = 0$$

$$Q(s,a) = r + \gamma \cdot \max_{\{a' \in A\}} Q(s',a')$$

This course is **Deep** RL, so we need to use a neural network, parameterized by  $\theta$ 

$$Q(s,a,\theta) - \left(r + \gamma \cdot \max_{\{a' \in A\}} Q(s',a',\theta)\right) = 0$$

Training objective

$$\min_{\theta} \left( Q(s, a, \theta) - \left( r + \gamma \cdot \max_{\{a' \in A\}} Q(s', a', \theta) \right) \right)^2$$

- Review
- State of the field
- Implement Deep Q Networks (DQN) (Mnih et al.)

# **Deep RL**

Like much of deep learning, Deep RL has a gap between theory and practice
## **Deep RL**

Like much of deep learning, Deep RL has a gap between theory and practice

Just like neural networks (1943) and backpropagation (1970), the theory of value functions (1957) and Q learning (1989) has been around for a long time

## **Deep RL**

Like much of deep learning, Deep RL has a gap between theory and practice

Just like neural networks (1943) and backpropagation (1970), the theory of value functions (1957) and Q learning (1989) has been around for a long time

Hardware advances and good ML software enabled us to take advantage of decades of theory. Mnih et al (2015) took Q learning theory and made it work well with neural networks

## **Deep RL**

Like much of deep learning, Deep RL has a gap between theory and practice

Just like neural networks (1943) and backpropagation (1970), the theory of value functions (1957) and Q learning (1989) has been around for a long time

Hardware advances and good ML software enabled us to take advantage of decades of theory. Mnih et al (2015) took Q learning theory and made it work well with neural networks

Since its inception, Deep RL has added "patches" to combine theory with deep networks to obtain better and better results

## Agenda

- Review
- State of the field
- Implement Deep Q Networks (DQN) (Mnih et al.)

```
dataset = load_dataset()
model = nn.Module(dataset.x.size, dataset.y.size)
theta = model.init(seed=0) # Functional
```

```
for update in range(num_updates):
    train_data = dataset.sample()
    theta = train(model, theta, train_data) # Functional
    metrics = evaluate(model, theta, dataset.val set)
```

```
==> dataset = load_dataset()
model = nn.Module(dataset.x.size, dataset.y.size)
theta = model.init(seed=0) # Functional
```

```
for update in range(num_updates):
    train_data = dataset.sample()
    theta = train(model, theta, train_data) # Functional
    metrics = evaluate(model, theta, dataset.val set)
```

Those familiar with deep learning know of the "training loop"

dataset = load\_dataset()

==> model = nn.Module(dataset.x.size, dataset.y.size)
 theta = model.init(seed=0) # Functional

for update in range(num\_updates):
 train\_data = dataset.sample()
 theta = train(model, theta, train\_data) # Functional
 metrics = evaluate(model, theta, dataset.val set)

```
dataset = load_dataset()
model = nn.Module(dataset.x.size, dataset.y.size)
==> theta = model.init(seed=0) # Functional
```

```
for update in range(num_updates):
    train_data = dataset.sample()
    theta = train(model, theta, train_data) # Functional
    metrics = evaluate(model, theta, dataset.val_set)
```

```
dataset = load_dataset()
model = nn.Module(dataset.x.size, dataset.y.size)
theta = model.init(seed=0) # Functional
```

```
==> for update in range(num_updates):
    train_data = dataset.sample()
    theta = train(model, theta, train_data) # Functional
    metrics = evaluate(model, theta, dataset.val_set)
```

```
dataset = load_dataset()
model = nn.Module(dataset.x.size, dataset.y.size)
theta = model.init(seed=0) # Functional
```

```
for update in range(num_updates):
    train_data = dataset.sample()
    theta = train(model, theta, train_data) # Functional
    metrics = evaluate(model, theta, dataset.val_set)
```

```
dataset = load_dataset()
model = nn.Module(dataset.x.size, dataset.y.size)
theta = model.init(seed=0) # Functional
```

```
for update in range(num_updates):
    train_data = dataset.sample()
==> theta = train(model, theta, train_data) # Functional
    metrics = evaluate(model, theta, dataset.val_set)
```

Those familiar with deep learning know of the "training loop"

```
dataset = load_dataset()
model = nn.Module(dataset.x.size, dataset.y.size)
theta = model.init(seed=0) # Functional
```

```
for update in range(num_updates):
    train_data = dataset.sample()
    theta = train(model, theta, train_data) # Functional
    metrics = evaluate(model, theta, dataset.val set)
```

==>

Deep RL has a training loop similar to the deep learning loop

Deep RL has a training loop similar to the deep learning loop

```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
pi = policy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

Main differences with the DL train loop

```
==> env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
==> pi = policy(Q, theta)
```

```
for update in range(num_updates):
```

```
==> collected_data = collect_training_data(env, pi)
==> dataset += collected_data
train_data = dataset.sample()
theta = train(Q, theta, train_data)
metrics = evaluate(env, pi)
```

**Today:** Go through the loop line by line to implement DQN

```
==> env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
==> pi = policy(Q, theta)
```

for update in range(num\_updates):

=> collected\_data = collect\_training\_data(env, pi)
=>> dataset += collected\_data
train\_data = dataset.sample()
theta = train(Q, theta, train\_data)
metrics = evaluate(env, pi)

Instead of loading a static dataset, we collect data from an environment

```
==> env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
pi = policy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

The Plan:

- 1. Terminal states
- 2. The Gymnasium interface

The Plan:

- 1. Terminal states
- 2. The Gymnasium interface

**Question:** How can we represent a Mario Bros Game Over screen in an MDP?



**Question:** How can we represent a Mario Bros Game Over screen in an MDP?

# Answer: We enter a **terminal state** that we cannot leave





How do we model the return in these terminal states?

How do we model the return in these terminal states?

Recall the discounted return

$$G = \sum_{t=0}^{\infty} \gamma^t r_t$$

How do we model the return in these terminal states?

Recall the discounted return

$$G = \sum_{t=0}^{\infty} \gamma^t r_t$$

After entering the terminal state at t = n all future rewards are zero. We can write the discounted return as

$$G = \sum_{t=0}^\infty \gamma^t r_t \cdot (t \le n) = \sum_{t=0}^n \gamma^t r_t$$

$$G = \sum_{t=0}^\infty \gamma^t r_t \cdot (t \le n) = \sum_{t=0}^n \gamma^t r_t$$

$$G = \sum_{t=0}^\infty \gamma^t r_t \cdot (t \le n) = \sum_{t=0}^n \gamma^t r_t$$

Many environments introduce the **done flag (d)** to simplify data collection and training

$$G = \sum_{t=0}^\infty \gamma^t r_t \cdot (t \le n) = \sum_{t=0}^n \gamma^t r_t$$

Many environments introduce the **done flag (d)** to simplify data collection and training

$$\begin{pmatrix} s_0 & s_1 & s_2 & \dots & s_n \\ d_0 = 0 & d_1 = 0 & d_2 = 0 & \dots & d_n = 1 \end{pmatrix}$$

$$G = \sum_{t=0}^\infty \gamma^t r_t \cdot (t \le n) = \sum_{t=0}^n \gamma^t r_t$$

Many environments introduce the **done flag (d)** to simplify data collection and training

$$\begin{pmatrix} s_0 & s_1 & s_2 & \dots & s_n \\ d_0 = 0 & d_1 = 0 & d_2 = 0 & \dots & d_n = 1 \end{pmatrix}$$

We call the states from the initial to terminal state an **episode** 

The Plan:

- 1. Terminal states
- 2. The Gymnasium interface

#### **Environment Libraries**

The gymnasium library contains many popular test environments





#### **Environment Libraries**

The gymnasium library contains many popular test environments





#### gymnasium also defines the standard environment interface

Launching environments is very easy

env = gymnasium.make("LunarLander-v2")

Launching environments is very easy

- env = gymnasium.make("LunarLander-v2")
- What is S, A for LunarLander?
- S, A = env.observation\_space, env.action\_space

Launching environments is very easy

env = gymnasium.make("LunarLander-v2")

What is S, A for LunarLander?

S, A = env.observation\_space, env.action\_space

**Observations** are states that are not guaranteed to be Markov. For LunarLander, they are Markov.

env = gymnasium.make("LunarLander-v2")
```
env = gymnasium.make("LunarLander-v2")
```

The environments start "off". We must reset the environment, which returns an initial state.

state, \_ = env.reset(seed=0)

```
env = gymnasium.make("LunarLander-v2")
```

The environments start "off". We must reset the environment, which returns an initial state.

state, \_ = env.reset(seed=0)

Step the environment in time by feeding an action (transition function T)

next\_state, reward, terminated, truncated, \_ = env.step(action)

```
env = gymnasium.make("LunarLander-v2")
```

The environments start "off". We must reset the environment, which returns an initial state.

state, \_ = env.reset(seed=0)

Step the environment in time by feeding an action (transition function T)

next\_state, reward, terminated, truncated, \_ = env.step(action)  $d = \text{terminated} \lor \text{truncated}$ 

You can write your own environments using gymnasium

```
class RoboParking(gymnasium.Env):
  observation_space = spaces.Box(
    # x, y, xdot, ydot
    low = (0, 0, -1, -1),
    high=(4, 4, 1, 1),
    dtype=np.float32
  # left, right, forward, backward
  action space = spaces.Discrete(4)
  def R(self, pos): # R(s')
```

```
# Our goal is 0,0,0,0
return norm(sensor.state())
```

def T(self, action):
 # We don't know the true T
 wheels.apply\_torque(action)
 next\_state = sensor.state()
 return next\_state

```
def step(self, action):
    next_state = T(action) # s'
    return (
        next_state,
        R(next_state), # reward R(s')
        norm(next_state) < 0.01 # d
        False, {} # trunc, extra info</pre>
```

We discussed the environment

```
==> env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
pi = policy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

Next, let us define the deep Q function

```
env = LunarLander()
```

=> Q = nn.Module(env.state\_space, env.action\_space)
 theta = Q.init(seed=0)
 pi = policy(Q, theta)

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

#### Recall the type signature of the Q function $Q: S \times A \rightarrow \mathbb{R}$

Recall the type signature of the Q function  $Q: S \times A \to \mathbb{R}$ And recall the optimal policy  $\pi_*(s) = \operatorname{argmax}_{a \in A} Q_*(s, a, \theta)$ 

Recall the type signature of the Q function  $Q: S \times A \to \mathbb{R}$ And recall the optimal policy  $\pi_*(s) = \operatorname{argmax}_{a \in A} Q_*(s, a, \theta)$ We would need to evaluate Q function |A| times for each state

Recall the type signature of the Q function  $Q: S \times A \to \mathbb{R}$ And recall the optimal policy  $\pi_*(s) = \operatorname{argmax}_{a \in A} Q_*(s, a, \theta)$ We would need to evaluate Q function |A| times for each state

 $egin{array}{ll} Q(s,a=1) \ Q(s,a=2) \ & \vdots \end{array}$ 

Recall the type signature of the Q function  $Q: S \times A \to \mathbb{R}$ And recall the optimal policy  $\pi_*(s) = \operatorname{argmax}_{a \in A} Q_*(s, a, \theta)$ We would need to evaluate Q function |A| times for each state

$$egin{array}{ll} Q(s,a=1) \ Q(s,a=2) \ dots \end{array}$$

Inefficient: |A| = 100 means 100 forward passes for each timestep

Recall the type signature of the Q function  $Q: S \times A \to \mathbb{R}$ And recall the optimal policy  $\pi_*(s) = \operatorname{argmax}_{a \in A} Q_*(s, a, \theta)$ 

We instead represent the Q network as

 $Q:S\to \mathbb{R}^{|A|}$ 

Recall the type signature of the Q function  $Q: S \times A \to \mathbb{R}$ And recall the optimal policy  $\pi_*(s) = \operatorname{argmax}_{a \in A} Q_*(s, a, \theta)$ We instead represent the Q network as

 $Q:S\to \mathbb{R}^{|A|}$ 

Compute the Q value for all actions in a single forward pass

Recall the type signature of the Q function  $Q: S \times A \to \mathbb{R}$ And recall the optimal policy  $\pi_*(s) = \operatorname{argmax}_{a \in A} Q_*(s, a, \theta)$ We instead represent the Q network as

 $Q:S\to \mathbb{R}^{|A|}$ 

Compute the Q value for all actions in a single forward pass

$$egin{aligned} Q(s, heta) &= Q(s, a = 1, heta), \ Q(s, a = 2, heta), \ &\vdots \end{aligned}$$

We instead represent the Q network as

 $Q:S\to \mathbb{R}^{|A|}$ 

#### We instead represent the Q network as

$$Q:S\to \mathbb{R}^{|A|}$$

**Architecture:** 2 layer MLP with hidden size of 256 is sufficient for standard benchmarks

#### We instead represent the Q network as

 $Q:S\to \mathbb{R}^{|A|}$ 

**Architecture:** 2 layer MLP with hidden size of 256 is sufficient for standard benchmarks

```
Q = Sequential(
  Linear(state_size, 256), LeakyReLU(),
  Linear(256, 256), LeakyReLU(),
  Linear(256, action_size)
```

#### We instead represent the Q network as

 $Q:S\to \mathbb{R}^{|A|}$ 

**Architecture:** 2 layer MLP with hidden size of 256 is sufficient for standard benchmarks

```
Q = Sequential(
```

Linear(state\_size, 256), LayerNorm(), LeakyReLU(), Linear(256, 256), LayerNorm(), LeakyReLU(), Linear(256, action\_size)

For states with structure (e.g., pixels), prepend encoders to Q

#### For states with structure (e.g., pixels), prepend encoders to Q



#### Init

Done with the Q function architecture, let's discuss init

```
env = LunarLander()
```

=> Q = nn.Module(env.state\_space, env.action\_space)
 theta = Q.init(seed=0)
 pi = policy(Q, theta)

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

#### Init

#### Done with the Q function architecture, let's discuss init

```
env = LunarLander()
      Q = nn.Module(env.state space, env.action space)
=> theta = Q.init(seed=0)
      pi = policy(Q, theta)
      for update in range(num updates):
        collected data = collect training data(env, pi)
        dataset += collected data
        train data = dataset.sample()
        theta = train(Q, theta, train data)
        metrics = evaluate(env, pi)
```

Not much to say for parameter initialization

Not much to say for parameter initialization

**Tip:** Initialize the final layer of your Q function to output values near 0

Not much to say for parameter initialization

**Tip:** Initialize the final layer of your Q function to output values near 0

nn.init.normal(std=1e-3, bias=0)

Not much to say for parameter initialization

**Tip:** Initialize the final layer of your Q function to output values near 0

nn.init.normal(std=1e-3, bias=0)

Prevents Q value overestimation (to be discussed in depth later)

Done with init, let's discuss policy  $\pi$ 

```
env = LunarLander()
     Q = nn.Module(env.state space, env.action space)
    theta = Q.init(seed=0)
==>
      pi = policy(Q, theta)
      for update in range(num updates):
        collected data = collect training data(env, pi)
        dataset += collected data
        train data = dataset.sample()
        theta = train(Q, theta, train data)
        metrics = evaluate(env, pi)
```

Done with init, let's discuss policy  $\pi$ 

```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
==> pi = policy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

Recall the policy  $\pi: S \to \Delta A$ 

Recall the policy  $\pi: S \to \Delta A$ 

In practice, we usually have two policies

Recall the policy  $\pi: S \to \Delta A$ 

In practice, we usually have two policies

1. The optimal policy we are trying to learn  $(\pi)$ 

Recall the policy  $\pi: S \to \Delta A$ 

In practice, we usually have two policies

- 1. The optimal policy we are trying to learn  $(\pi)$
- 2. An **exploration policy** ( $\pi_E$ ) for collecting training data

Recall the policy  $\pi: S \to \Delta A$ 

In practice, we usually have two policies

- 1. The optimal policy we are trying to learn  $(\pi)$
- 2. An **exploration policy** ( $\pi_E$ ) for collecting training data

Why do we need  $\pi_E$ ?

Recall the policy  $\pi: S \to \Delta A$ 

In practice, we usually have two policies

- 1. The optimal policy we are trying to learn  $(\pi)$
- 2. An **exploration policy** ( $\pi_E$ ) for collecting training data

Why do we need  $\pi_E$ ?



#### **Exploration Policy**

We use the **exploration policy** ( $\pi_E$ ) to explore and collect data

## **Exploration Policy**

We use the **exploration policy** ( $\pi_E$ ) to explore and collect data

We will use our collected data to train the Q function. What properties should our collected data have?
We use the **exploration policy** ( $\pi_E$ ) to explore and collect data

We will use our collected data to train the Q function. What properties should our collected data have?

1. To ensure Q is accurate everywhere, take every possible action in every possible state

We use the **exploration policy** ( $\pi_E$ ) to explore and collect data

We will use our collected data to train the Q function. What properties should our collected data have?

1. To ensure Q is accurate everywhere, take every possible action in every possible state

Given these requirements, we cannot do better than random exploration

 $\pi_E(s) = \mathcal{U}(A)$ 

$$\pi_E(s) = \mathcal{U}(A)$$

**Question:** Are there any downsides to this exploration policy?

$$\pi_E(s) = \mathcal{U}(A)$$

**Question:** Are there any downsides to this exploration policy?

Answer: It could take a really, really long time

$$\pi_E(s) = \mathcal{U}(A)$$

**Question:** Are there any downsides to this exploration policy?

Answer: It could take a really, really long time



$$\pi_E(s) = \mathcal{U}(A)$$

Alternative: Bias the policy towards states with known large Q values

$$\pi_E(s) = \mathcal{U}(A)$$

Alternative: Bias the policy towards states with known large Q values



$$\pi_E(s) = \mathcal{U}(A)$$

Alternative: Bias the policy towards states with known large Q values

One approach is the  $\epsilon$ -greedy policy

$$\pi_E(s) = \mathcal{U}(A)$$

Alternative: Bias the policy towards states with known large Q values

One approach is the  $\epsilon$ -greedy policy

$$\begin{split} \epsilon \in [0,1] \\ u \sim \mathcal{U}[0,1] \\ \pi_E(s) = \begin{cases} \mathcal{U}(A) \text{ if } u \leq \epsilon \\ \operatorname{argmax}_{a \in A} Q(s,a) \text{ if } u > \epsilon \end{cases} \end{split}$$

$$\pi_E(s) = \mathcal{U}(A)$$

Alternative: Bias the policy towards states with known large Q values

One approach is the  $\epsilon$ -greedy policy

$$\begin{split} \epsilon \in [0,1] \\ u \sim \mathcal{U}[0,1] \\ \pi_E(s) = \begin{cases} \mathcal{U}(A) \text{ if } u \leq \epsilon \\ \arg\max_{a \in A} Q(s,a) \text{ if } u > \epsilon \end{cases} \end{split}$$

This approach is simple and works surprisingly well in practice





Explore promising areas with large Q values more often and as  $t \to \infty$ , explore all state/action tuples



Explore promising areas with large Q values more often and as  $t \to \infty$ , explore all state/action tuples

**Question:** Any downsides to the  $\varepsilon$ -greedy approach?



**Question:** Any downsides to the  $\varepsilon$ -greedy approach?

**Answer:** Not independently distributed. The state/action distribution will be biased by the Q function. Seems to be ignored in practice?

**Summary:** Maintain two policies

Summary: Maintain two policies

 $\pi$ : The policy that approximates  $\pi_*$ 

Summary: Maintain two policies

 $\pi$ : The policy that approximates  $\pi_*$ 

 $\pi_E: \mathbf{A}$  stochastic policy used for exploring the environment and collecting data

Let us make a small change to the pseudocode

```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
==> pi = policy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

Now we have two policies, one for collection and one for evaluation

```
env = LunarLander()
```

Q = nn.Module(env.state\_space, env.action\_space)
theta = Q.init(seed=0)

=> pi, pi\_e = max\_q(Q, theta), e\_greedy(Q, theta)

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi_e)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

==>

Let's move onto data collection

env = LunarLander()
Q = nn.Module(env.state\_space, env.action\_space)
theta = Q.init(seed=0)
pi, pi\_e = max\_q(Q, theta), e\_greedy(Q, theta)

for update in range(num\_updates):
 collected\_data = collect\_training\_data(env, pi\_e)
 dataset += collected\_data
 train\_data = dataset.sample()
 theta = train(Q, theta, train\_data)
 metrics = evaluate(env, pi)

#### We must interact with the MDP to collect training data

We must interact with the MDP to collect training data

Recall the Q learning objective

$$\min_{\theta} \left( Q(\underline{s},\underline{a},\theta) - \left( \underline{r} + \gamma \cdot \operatorname*{argmax}_{\{a' \in A\}} Q(\underline{s'},a',\theta) \right) \right)^2$$

We must interact with the MDP to collect training data

Recall the Q learning objective

$$\min_{\theta} \left( Q(\underline{s}, \underline{a}, \theta) - \left( \underline{r} + \gamma \cdot \operatorname*{argmax}_{\{a' \in A\}} Q(\underline{s'}, a', \theta) \right) \right)^2$$

Many algorithms train using a transition tuple (s, a, r, s', d)

Collecting transitions correctly is deceptively tricky (off by one errors)

Collecting transitions correctly is deceptively tricky (off by one errors)

```
states, next states, rewards, actions, dones = [], [], ...
s, = env.reset(seed=0)
d = False
while not d:
 a = pi e(s, theta)
  next_s, r, trunc, term, = env.step(action) # r = R(s')
  d = trunc or term
  states.append(s), next states.append(next s), rewards...
  s = next s
# n+1 states total, but each list should be len n
episode = (states, next states, rewards, actions, dones)
return episode
```

We call the dataset a **replay buffer** ( $\mathcal{D}$ )

#### We call the dataset a **replay buffer** ( $\mathcal{D}$ )

$$\mathcal{D} = \begin{bmatrix} (s_0, a_0, r_0, s_0', d_0) \\ (s_1, a_1, r_1, s_1', d_1) \\ \vdots \end{bmatrix}$$

#### We call the dataset a **replay buffer** ( $\mathcal{D}$ )

$$\mathcal{D} = \begin{bmatrix} (s_0, a_0, r_0, s_0', d_0) \\ (s_1, a_1, r_1, s_1', d_1) \\ \vdots \end{bmatrix}$$

#### We call the dataset a **replay buffer** ( $\mathcal{D}$ )

$$\mathcal{D} = \begin{bmatrix} (s_0, a_0, r_0, s_0', d_0) \\ (s_1, a_1, r_1, s_1', d_1) \\ \vdots \end{bmatrix}$$

**Note:** We often enforce a max size of  $\mathcal D$  using a ring buffer

We populated the dataset, now let's sample from it

```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
pi, pi_e = max_q(Q, theta), e_greedy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi_e)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

==>

We populated the dataset, now let's sample from it

```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
pi, pi_e = max_q(Q, theta), e_greedy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi_e)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

==>

We sample training data from  ${\mathcal D}$ 

We sample training data from  ${\mathcal D}$ 

We will call the train data our training batch  ${\mathcal B}$ 

 $\mathcal{B} \sim [\mathcal{U}(\mathcal{D}), ..., \mathcal{U}(\mathcal{D})]$ 

We sample training data from  ${\mathcal D}$ 

We will call the train data our training batch  ${\mathcal B}$ 

 $\mathcal{B} \sim [\mathcal{U}(\mathcal{D}), ..., \mathcal{U}(\mathcal{D})]$ 

$$\mathcal{B} = \left[ \left(s_j, a_j, r_j, s_j', d_j\right), ..., \left(s_k, a_k, r_k, s_k', d_k\right) \right]$$

We sample training data from  ${\mathcal D}$ 

We will call the train data our training batch  ${\mathcal B}$ 

 $\mathcal{B} \sim [\mathcal{U}(\mathcal{D}), ..., \mathcal{U}(\mathcal{D})]$ 

$$\mathcal{B} = \left[ \left(s_j, a_j, r_j, s_j', d_j\right), ..., \left(s_k, a_k, r_k, s_k', d_k\right) \right]$$

Randomly sampling old data helps mitigate correlations between data, improving training stability

Randomly sampling old data helps mitigate correlations between data, improving training stability
# **Replay Buffers**

Randomly sampling old data helps mitigate correlations between data, improving training stability

Biased towards many prior policies instead of one

# **Replay Buffers**

Randomly sampling old data helps mitigate correlations between data, improving training stability

Biased towards many prior policies instead of one



```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
pi, pi_e = max_q(Q, theta), e_greedy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi_e)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
pi, pi_e = max_q(Q, theta), e_greedy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi_e)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, train_data)
    metrics = evaluate(env, pi)
```

We make a few modifications to the Q learning objective to improve performance

We make a few modifications to the Q learning objective to improve performance

1. Early termination using d

We make a few modifications to the Q learning objective to improve performance

- 1. Early termination using d
- 2. Target networks

We make a few modifications to the Q learning objective to improve performance

- 1. Early termination using d
- 2. Target networks

Recall the standard Q learning objective

$$\min_{\theta} \left( Q(s, a, \theta) - \left( r + \gamma \cdot \max_{\{a' \in A\}} Q(s', a', \theta) \right) \right)^2$$

Recall the standard Q learning objective

$$\min_{\boldsymbol{\theta}} \left( Q(s, a, \boldsymbol{\theta}) - \left( r + \gamma \cdot \max_{\{a' \in A\}} Q(s', a', \boldsymbol{\theta}) \right) \right)^2$$

Rather than learn to output 0 at terminal states, we modify the objective

$$\min_{\boldsymbol{\theta}} \left( Q(s, a, \boldsymbol{\theta}) - \left( r + \neg \boldsymbol{d} \cdot \boldsymbol{\gamma} \cdot \max_{\{a' \in A\}} Q(s', a', \boldsymbol{\theta}) \right) \right)^2$$

Recall the standard Q learning objective

$$\min_{\boldsymbol{\theta}} \left( Q(s, a, \boldsymbol{\theta}) - \left( r + \gamma \cdot \max_{\{a' \in A\}} Q(s', a', \boldsymbol{\theta}) \right) \right)^2$$

Rather than learn to output 0 at terminal states, we modify the objective

$$\min_{\boldsymbol{\theta}} \left( Q(s, a, \boldsymbol{\theta}) - \left( r + \neg \boldsymbol{d} \cdot \boldsymbol{\gamma} \cdot \max_{\{a' \in A\}} Q(s', a', \boldsymbol{\theta}) \right) \right)^2$$

For terminal transitions, this reduces to

$$\min_{\theta} \left( Q(s, a, \theta) - r \right)^2$$

We make a few modifications to the Q learning objective to improve performance

- 1. Early termination using d
- 2. Target networks

Traditional Q learning used a table

|       | s = 0 | s = 1 | s = 2 |
|-------|-------|-------|-------|
| a = 0 | 1     | 3     | 2     |
| a = 1 | 7     | 4     | 5     |
| a=2   | 1     | 0     | 0     |

Traditional Q learning used a table

|       | s = 0 | s = 1 | s = 2 |
|-------|-------|-------|-------|
| a = 0 | 1     | 3     | 2     |
| a = 1 | 7     | 4     | 5     |
| a=2   | 1     | 0     | 0     |

#### Updates are very well defined

|       | s = 0 | s = 1 | s = 2 |
|-------|-------|-------|-------|
| a = 0 | 5     | 3     | 2     |
| a = 1 | 7     | 4     | 5     |
| a=2   | 1     | 0     | 0     |

Traditional Q learning used a table

|       | s = 0 | s = 1 | s = 2 |
|-------|-------|-------|-------|
| a = 0 | 1     | 3     | 2     |
| a = 1 | 7     | 4     | 5     |
| a=2   | 1     | 0     | 0     |

Updates are very well defined

|       | s = 0 | s = 1 | s = 2 |
|-------|-------|-------|-------|
| a = 0 | 5     | 3     | 2     |
| a = 1 | 7     | 4     | 5     |
| a=2   | 1     | 0     | 0     |

Neural networks are different. Increasing a single (s = 0, a = 0)entry will often **perturb** the Q value for all states and actions.

|       | s = 0 | s = 1 | s = 2 |
|-------|-------|-------|-------|
| a = 0 | 5     | 4     | 4     |
| a = 1 | 9     | 5     | 5     |
| a=2   | 4     | 0     | 0     |

These perturbations  $\varepsilon$  ripple through the Q recursion, with the max operator resulting in overestimation

$$Q(s, a, \theta') \leftarrow r + \gamma \max_{a' \in A} [Q(s', a', \theta) + \varepsilon_{\theta}]$$

These perturbations  $\varepsilon$  ripple through the Q recursion, with the max operator resulting in overestimation

$$Q(s, a, \theta') \leftarrow r + \gamma \max_{a' \in A} [Q(s', a', \theta) + \varepsilon_{\theta}]$$

$$Q(s, a, \theta'') \leftarrow r + \gamma \Bigl( \max_{a' \in A} [Q(s', a', \theta') + \varepsilon_{\theta} + \varepsilon_{\theta'}] \Bigr)$$

These perturbations  $\varepsilon$  ripple through the Q recursion, with the max operator resulting in overestimation

$$Q(s, a, \theta') \leftarrow r + \gamma \max_{a' \in A} [Q(s', a', \theta) + \varepsilon_{\theta}]$$

$$Q(s, a, \theta'') \leftarrow r + \gamma \Big( \max_{a' \in A} [Q(s', a', \theta') + \varepsilon_{\theta} + \varepsilon_{\theta'}] \Big)$$

$$Q(s, a, \theta''') \leftarrow r + \gamma \left( \max_{a' \in A} [Q(s', a', \theta'') + \varepsilon_{\theta} + \varepsilon_{\theta'} + \varepsilon_{\theta''}] \right)$$

These perturbations  $\varepsilon$  ripple through the Q recursion, with the max operator resulting in overestimation

$$Q(s, a, \theta') \leftarrow r + \gamma \max_{a' \in A} [Q(s', a', \theta) + \varepsilon_{\theta}]$$

$$Q(s, a, \theta'') \leftarrow r + \gamma \bigg( \max_{a' \in A} [Q(s', a', \theta') + \varepsilon_{\theta} + \varepsilon_{\theta'}] \bigg)$$

$$Q(s, a, \theta''') \leftarrow r + \gamma \left( \max_{a' \in A} [Q(s', a', \theta'') + \varepsilon_{\theta} + \varepsilon_{\theta'} + \varepsilon_{\theta''}] \right)$$

Compounding pertubations combined with the max operator result in exploding Q values (i.e.,  $Q(\cdot, \cdot) = \infty)$ 

Compounding pertubations combined with the max operator result in exploding Q values (i.e.,  $Q(\cdot, \cdot) = \infty)$ 

Compounding pertubations combined with the max operator result in exploding Q values (i.e.,  $Q(\cdot, \cdot) = \infty)$ 

**Solution 1:** Constrained optimization of neural networks (hard)

Compounding pertubations combined with the max operator result in exploding Q values (i.e.,  $Q(\cdot, \cdot) = \infty)$ 

Solution 1: Constrained optimization of neural networks (hard)Solution 2: Very large batch sizes that cover all (*s*, *a*) (intractable)

Compounding pertubations combined with the max operator result in exploding Q values (i.e.,  $Q(\cdot, \cdot) = \infty)$ 

Solution 1: Constrained optimization of neural networks (hard)
Solution 2: Very large batch sizes that cover all (*s*, *a*) (intractable)
Solution 3: Surrogate target network to break recurrence (easy)

**Solution 3:** Surrogate **target network** to break recurrence (easy)

Initialize target parameters  $\psi = \theta$ 

$$Q(s, a, \theta') \leftarrow r + \gamma \max_{a' \in A} \left[ Q(s', a', \psi) + \varepsilon_{\psi} \right]$$

**Solution 3:** Surrogate **target network** to break recurrence (easy) Initialize target parameters  $\psi = \theta$ 

$$Q(s, a, \theta') \leftarrow r + \gamma \max_{a' \in A} \left[ Q(s', a', \psi) + \varepsilon_{\psi} \right]$$

$$Q(s, a, \theta'') \leftarrow r + \gamma \Big( \max_{a' \in A} \big[ Q(s', a', \psi) + \varepsilon_\psi \big] \Big)$$

**Solution 3:** Surrogate **target network** to break recurrence (easy) Initialize target parameters  $\psi = \theta$ 

$$Q(s, a, \theta') \leftarrow r + \gamma \max_{a' \in A} \left[ Q(s', a', \psi) + \varepsilon_{\psi} \right]$$

$$Q(s, a, \theta'') \leftarrow r + \gamma \Big( \max_{a' \in A} \big[ Q(s', a', \psi) + \varepsilon_{\psi} \big] \Big)$$

$$Q(s, a, \theta''') \leftarrow r + \gamma \Big( \max_{a' \in A} \big[ Q(s', a', \psi) + \varepsilon_{\psi} \big] \Big)$$

**Solution 3:** Surrogate **target network** to break recurrence (easy) Initialize target parameters  $\psi = \theta$ 

$$Q(s, a, \theta') \leftarrow r + \gamma \max_{a' \in A} \left[ Q(s', a', \psi) + \varepsilon_{\psi} \right]$$

$$Q(s, a, \theta'') \leftarrow r + \gamma \Bigl( \max_{a' \in A} \bigl[ Q(s', a', \psi) + \varepsilon_{\psi} \bigr] \Bigr)$$

$$Q(s, a, \theta''') \leftarrow r + \gamma \Big( \max_{a' \in A} \big[ Q(s', a', \psi) + \varepsilon_{\psi} \big] \Big)$$

After a while, set  $\psi = \theta$  again

#### Behold, the combination of early termination and target networks

$$Q(s, a, \theta) = r + \neg \boldsymbol{d} \cdot \boldsymbol{\gamma} \cdot \max_{a' \in A} Q(s', a', \boldsymbol{\psi})$$

Behold, the combination of early termination and target networks

$$Q(s, a, \theta) = r + \neg d \cdot \gamma \cdot \max_{a' \in A} Q(s', a', \psi)$$

$$Q(s, a, \theta) - \left(r + \neg \boldsymbol{d} \cdot \boldsymbol{\gamma} \cdot \max_{a' \in A} Q(s', a', \boldsymbol{\psi})\right) = 0$$

Behold, the combination of early termination and target networks

$$Q(s, a, \theta) = r + \neg \boldsymbol{d} \cdot \gamma \cdot \max_{a' \in A} Q(s', a', \psi)$$

$$Q(s, a, \theta) - \left(r + \neg \boldsymbol{d} \cdot \boldsymbol{\gamma} \cdot \max_{a' \in A} Q(s', a', \boldsymbol{\psi})\right) = 0$$

The standard objective for DQN is

$$\min_{\boldsymbol{\theta}} \left( Q(s, a, \boldsymbol{\theta}) - \left( r + \neg \boldsymbol{d} \cdot \boldsymbol{\gamma} \cdot \max_{a' \in A} Q(s', a', \boldsymbol{\psi}) \right) \right)^2$$

We need to make a few small updates given our new objective

```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta = Q.init(seed=0)
pi, pi_e = max_q(Q, theta), e_greedy(Q, theta)
```

for update in range(num\_updates):
 collected\_data = collect\_training\_data(env, pi\_e)
 dataset += collected\_data
 train\_data = dataset.sample()
 theta = train(Q, theta, train\_data)
 metrics = evaluate(env, pi)

Initialize target parameters, and use target params in loss function

```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
==> theta, psi = Q.init(seed=0), Q.init(seed=0)
pi, pi_e = max_q(Q, theta), e_greedy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi_e)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, psi, train_data)
    metrics = evaluate(env, pi)
```

## **Evaluation**

Same as collect\_training\_data but use  $\pi$  not  $\pi_E$ 

```
env = LunarLander()
Q = nn.Module(env.state_space, env.action_space)
theta, psi = Q.init(seed=0), Q.init(seed=0)
pi, pi_e = max_q(Q, theta), e_greedy(Q, theta)
```

```
for update in range(num_updates):
    collected_data = collect_training_data(env, pi_e)
    dataset += collected_data
    train_data = dataset.sample()
    theta = train(Q, theta, psi, train_data)
    metrics = evaluate(env, pi)
```

## **Summary**

- Review
- State of the field
- Implement Deep Q Networks (DQN) (Mnih et al.)

## **Next Time**

- Zach, Riccardo, Grace, Dylan, and Saksham will be lecturing
  - 1. Give us a hint on the topic!
  - 2. Turn in reports (email is best)
  - 3. 10 min presentation + 5 min questions and discussion
    - Next lecture is not recorded (reduce anxiety)
- Miniproject handout
  - Moodle says due 22 March 16:00