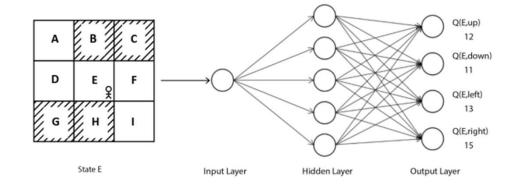
Policy Gradient Methods



Value-based RL

- In the previous chapters, we learned various methods to find the optimal policy in a given environment.
- The methods involved (iteratively) calculating Q-values of actions in each state.
- From the Q-values, we extract the optimal policy.
 - In each state, we choose an action with the highest Q value.
- These methods using Q-values are called "value-based" RL methods.

State	Action	Value
S ₀	0	10
S ₀	1	5
S ₁	0	10
S ₁	1	11





Disadvantages of Value-based RL Methods

- It is suitable for discrete environments
 - environments with discrete action space
- It cannot be directly applied to continuous environments
 - environments with continuous action space
- Suppose we are training an agent to drive a car and say we have one continuous action in our action space: speed.
- The speed ranges from 0km/h to 150km/h.
- In this case, how can we compute the Q value of all possible state-action pairs?



Disadvantages of Value-based RL Methods

- One way to use a Q table or a Q network is to discretize the continuous actions.
 - action 1: 0km/h to 10km/h
 - action 2: 10km/h to 20km/h
 - and so on.
- Discretization can work, but is not desirable depending on the task.
 - Granularity of action is degraded
 - may end up with huge number of actions
- Many real-world problems have continuous action space
 - self-driving cars, robots, etc.



Policy-based RL Methods

- There is another approach of reinforcement learning called policy-based RL methods.
- In policy-based RL, we do not need to compute the Q function.
- Instead, the optimal policy is computed directly.
- Policy-based methods have several advantages over valuebased methods
 - One is that they support continuous action spaces.



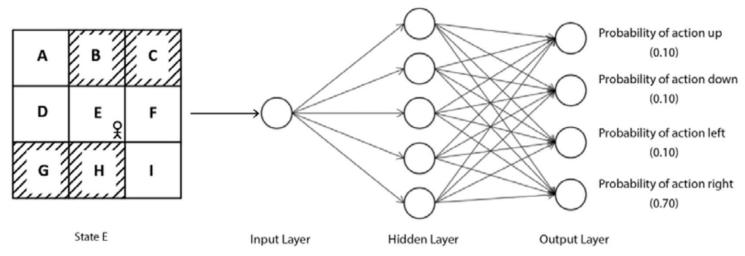
Policy Gradient: Intuition

- A policy-based method: does not compute Q function
- Directly parameterizes the policy using some parameter θ
- In a policy-based method, we use a stochastic policy.
 - In a stochastic policy, actions in a state is chosen based on a probability distribution.
 - Using a stochastic policy is the means for exploration in a policybased method.
 - According to a stochastic policy π , the probability of choosing action a in a state s is $\pi(a|s)$.
 - Furthermore, we parameterized our policy, and denote the policy as $\pi_{\theta}(a|s)$.
 - The policy is parameterized by θ .



Policy Gradient: Intuition

- What does it mean to parameterize the policy?
- It means we assume there is some function that maps the state to the probability of actions in that state.
- As for the function, we can use a neural network called policy network as a function approximator.
- Parameters of the policy network is θ , which is trained using the gradient descent algorithm.





DQN vs. Policy Gradient

- In DQN, we feed the state as an input to the network, and it returns the Q values of all possible actions in that state.
- We select an action that has the maximum Q value.
- In the Policy Gradient method, we feed the state as an input to the network, and it returns the probability distribution over an action space.
- We select an action using the probability distribution.



Learning in the Policy Gradient Method

- We play an episode and store the states, actions, and rewards.
 This becomes our training data.
- If we get a positive return, we increase the probability of all the actions that we took in each state until the end of the episode.
- If we get a negative return, we decrease the probability of all the actions that we took in each state until the end of the episode.



Learning in the Policy Gradient Method: Example

- We have eight states s_1 to s_8 , and our goal is to reach state s_8 .
- The action space only consists of two actions: *left* and *right*.
- We have a policy network, which returns the probability of each actions when we feed the state.
- Suppose we select an action in each state based on the probability distribution returned by the policy network.

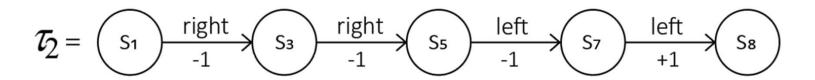
$$T_{l} = S_{1} \xrightarrow{left} S_{2} \xrightarrow{right} S_{4} \xrightarrow{left} S_{6} \xrightarrow{right} S_{8}$$

- The return of this trajectory is $R(\tau_1) = 1 + 1 + 1 + 1 = 4$.
- Since we got a positive return, we increase the probabilities of all the actions that we took in each state until the end of the episode.
- The probabilities of $(s_1, left)$, $(s_2, right)$, $(s_4, left)$, $(s_6, right)$ are increased.



Learning in the Policy Gradient Method: Example

• We generate another trajectory τ_2 , where we also select actions based on the probability distribution given by the policy network.



- The return of this trajectory is $R(\tau_2) = -1 1 1 + 1 = -2$.
- Since we got a negative return, we decrease the probabilities of all the actions that we took in each state until the end of the episode.
- The probabilities of $(s_1, right)$, $(s_3, right)$, $(s_5, left)$, $(s_7, left)$ are decreased.



Learning in the Policy Gradient Method

- How exactly do we increase and decrease probabilities?
- This is done using backpropagation, an algorithm used to update the parameters of the policy network.
- The updates are done in such a way that the actions yielding high return will get high probabilities, and actions yielding low return will get low probabilities.



- The goal of the policy gradient method is to find the optimal parameter θ of the neural network, so that the network returns the "correct" probability distribution over the action space.
- The objective of our network is to assign high probabilities to actions that maximize the expected return of the trajectory.
- The objective function can be written as:
 - $-J(\theta) = \mathbb{E}_{\tau \sim \pi_{\theta}(\tau)}[R(\tau)]$
 - τ : trajectory
 - $\tau \sim \pi_{\theta}(\tau)$: we are sampling the trajectory based on the policy π given by the network parameterized by θ .
 - $R(\tau)$: return of the trajectory τ .
- We want to maximize the objective function.



- To minimize the objective function, we calculate the gradients of the objective function and update the parameter using gradient descent.
- This time, we need to maximize the objective function. So we use gradient ascent.
- $\theta = \theta + \alpha \nabla_{\theta} J(\theta)$
 - $\nabla_{\theta} J(\theta)$: gradients of the the objective function



• The gradient $\nabla_{\theta} J(\theta)$ can be derived as:

$$- \nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim \pi_{\theta}(\tau)} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} \left(a_{t} | s_{t} \right) R(\tau) \right]$$

- So the gradient ascent becomes:
 - $-\theta = \theta + \alpha \sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} (a_t | s_t) R(\tau)$
 - $\log \pi_{\theta}(a_t|s_t)$: log probability of taking an action a given the state s at time t.
 - $R(\tau)$: return of the trajectory.
- We can see from the equation that when $R(\tau)$ is positive, we are increasing the probability of action a_t in state s_t . When $R(\tau)$ is negative, we are decreasing the probability of action a_t in state s_t .



 For the expectation, we follow the Monte Carlo approach and change the expectation to average over N trajectories.

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim \pi_{\theta}(\tau)} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} (a_t | s_t) R(\tau) \right]$$

$$\nabla_{\theta} J(\theta) = \frac{1}{N} \sum_{i=1}^{N} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} (a_t | s_t) R(\tau) \right]$$

Update rule

$$\theta = \theta + \alpha \frac{1}{N} \sum_{i=1}^{N} \sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} (a_t | s_t) R(\tau)$$

We repeat this step for many iterations to find the optimal policy.



Derivation of the following equation

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim \pi_{\theta}(\tau)} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) R(\tau) \right]$$

• We start from the definition of $J(\theta)$,

$$J(\theta) = \mathbb{E}_{\tau \sim \pi_{\theta}(\tau)}[R(\tau)]$$

• When X is a continuous random variable whose probability density function (pdf) is given as p(x), the expection of f(X) can be defined as:

$$\mathbb{E}_{x \sim p(x)}[f(X)] = \int_{\mathcal{X}} p(x)f(x)dx$$

• Thus, from the definition of $J(\theta)$,

$$J(\theta) = \int \pi_{\theta}(\tau) R(\tau) d\tau$$



Now we apply the derivative,

$$\nabla_{\theta} J(\theta) = \int \nabla_{\theta} \pi_{\theta}(\tau) R(\tau) d\tau$$

• multiply and divide by $\pi_{\theta}(\tau)$,

$$\nabla_{\theta} J(\theta) = \int \nabla_{\theta} \pi_{\theta}(\tau) \frac{\pi_{\theta}(\tau)}{\pi_{\theta}(\tau)} R(\tau) d\tau$$

and rearrange the equation.

$$\nabla_{\theta} J(\theta) = \int \pi_{\theta}(\tau) \frac{\nabla_{\theta} \pi_{\theta}(\tau)}{\pi_{\theta}(\tau)} R(\tau) d\tau$$



We use the log derivative trick which is:

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

• Using the trick, we get

$$\nabla_{\theta} J(\theta) = \int \pi_{\theta}(\tau) \nabla_{\theta} \log \pi_{\theta}(\tau) R(\tau) d\tau$$

• If we write this in the expection form:

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim \pi_{\theta}(\tau)} [\nabla_{\theta} \log \pi_{\theta}(\tau) R(\tau)]$$



- How to compute $\nabla_{\theta} \log \pi_{\theta}(\tau)$
- The probability distribution of trajectory can be given as:

$$\pi_{\theta}(\tau) = p(s_0) \prod_{t=0}^{T-1} \pi_{\theta}(a_t|s_t) p(s_{t+1}|s_t, a_t)$$

- $p(s_0)$ is the initial state distribution.
- Taking the log on both sides, we can write:

$$\log \pi_{\theta}(\tau) = \log \left[p(s_0) \prod_{t=0}^{T-1} \pi_{\theta}(a_t|s_t) p(s_{t+1}|s_t, a_t) \right]$$



• Log of a product is equal to sum of the logs, thus:

$$\log \pi_{\theta}(\tau) = \log p(s_0) + \log \prod_{t=0}^{T-1} \pi_{\theta}(a_t|s_t) p(s_{t+1}|s_t, a_t)$$

• Using the same rule, we can further write

$$\log \pi_{\theta}(\tau) = \log p(s_0) + \sum_{t=0}^{T-1} \left[\log \pi_{\theta}(a_t|s_t) + \log p(s_{t+1}|s_t, a_t) \right]$$

Now we apply the derivative to the equation.

$$\nabla_{\theta} \log \pi_{\theta}(\tau) = \nabla_{\theta} \left[\log p(s_0) + \sum_{t=0}^{T-1} \left[\log \pi_{\theta}(a_t|s_t) + \log p(s_{t+1}|s_t, a_t) \right] \right]$$



• When taking derivative, the terms $\log p(s_0)$ and $\log p(s_{t+1}|s_t,a_t)$ are not related to θ , so they become zero.

$$\nabla_{\theta} \log \pi_{\theta}(\tau) = \sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta}(a_t|s_t)$$

• Using this equation, we get the following:

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\tau \sim \pi_{\theta}(\tau)} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) R(\tau) \right]$$



The Policy Gradient Algorithm

- This algorithm is called REINFORCE, or Monte Carlo policy gradient.
- 1. Initialize the network parameter θ with random values
- 2. Generate N trajectories $\{\tau^i\}_{i=1}^N$ following the policy π_{θ}
- 3. Compute the return of the trajectory $R(\tau)$
- 4. Compute the gradients

$$\nabla_{\theta} J(\theta) = \frac{1}{N} \sum_{i=1}^{N} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) R(\tau) \right]$$

- 5. Update the network parameter as $\theta = \theta + \alpha \nabla_{\theta} J(\theta)$
- 6. Repeat *steps* 2 to 5 for several iterations



Policy Gradient on Continuous Action Space

- We saw that in the policy gradient method, the policy network returns a probability distribution of actions in a state.
- Then, we select an action based on the probability distribution.
- However, this applies to a discrete action space.
- If the action space is continuous, we can have the policy network return the mean and the variance of the action as output.
- We can generate a Gaussian distribution using the mean and the variance, and select an action by sampling from the distribution.



Variance Reduction Methods

- The REINFORCE method is an on-policy method, because we improve the same policy with which we are generating episodes in every iteration.
- One of the major issue in the method is the the gradient, $\nabla_{\theta} J(\theta)$, will have high variance in each update.
 - The return varies greatly in each episode.
- When the variance of gradient is high, it takes a lot of time to converge.



Variance Reduction Methods

- Methods to reduce the variance
 - policy gradient with reward-to-go (causality)
 - policy gradient with baseline



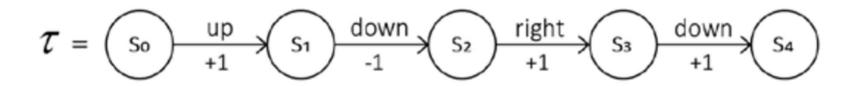
In the original method, the policy gradient is computed as:

$$\nabla_{\theta} J(\theta) = \frac{1}{N} \sum_{i=1}^{N} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} (a_t | s_t) R(\tau) \right]$$

- We make a small change in the equation.
- Instead of using the full return of the trajectory $R(\tau)$, we use the reward-to-go R_t , which is the return of the trajectory starting from state s_t .



Let us consider one episode, in which the trajectory was as follows.



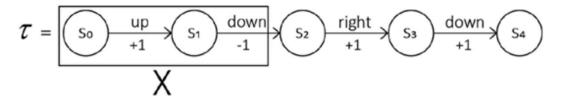
- The return of the trajectory is $R(\tau) = +1 1 + 1 + 1 = 2$.
- Now, we compute the gradient as :
 - $\nabla_{\theta} J(\theta) = \nabla_{\theta} \log \pi_{\theta}(\text{up}|s_{0})R(\tau) + \nabla_{\theta} \log \pi_{\theta}(\text{down}|s_{1})R(\tau) + \nabla_{\theta} \log \pi_{\theta}(\text{right}|s_{2})R(\tau) + \nabla_{\theta} \log \pi_{\theta}(\text{down}|s_{3})R(\tau)$
- In every step of the episode, we are multiplying the log probability of the action by the return of the full trajectory $R(\tau)$.



• Suppose we want to know how good the action right is in the state s_2 .

$$T = S_0 \xrightarrow{\text{up}} S_1 \xrightarrow{\text{down}} S_2 \xrightarrow{\text{right}} S_3 \xrightarrow{\text{down}} S_4$$

- If the action *right* in the state s_2 is a good action, we increase its probability. If not, then we decrease its probability.
- To find out whether the action right in the state s_2 is good, we do not need to consider the states and actions that came before s_2 .



- Thus, we use the return starting from s_2 , which we call reward-to-go.
 - $\nabla_{\theta} J(\theta) = \nabla_{\theta} \log \pi_{\theta}(\text{up}|s_{0}) R_{0} + \nabla_{\theta} \log \pi_{\theta}(\text{down}|s_{1}) R_{1} + \nabla_{\theta} \log \pi_{\theta}(\text{right}|s_{2}) R_{2} + \nabla_{\theta} \log \pi_{\theta}(\text{down}|s_{3}) R_{3}$

$$R_t = \sum_{t'=t}^{T-1} r(s_{t'}, a_{t'})$$



 In policy gradient with reward-to-go, the gradient is expressed as:

$$\nabla_{\theta} J(\theta) = \frac{1}{N} \sum_{i=1}^{N} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} (a_t | s_t) R_t \right]$$

$$R_t = \sum_{t'=t}^{T-1} r(s_{t'}, a_{t'})$$



- The algorithm
- 1. Initialize the network parameter θ with random values
- 2. Generate *N* number of trajectories $\{\tau^i\}_{i=1}^N$ following the policy π_θ
- 3. Compute the return (reward-to-go) R_t
- 4. Compute the gradients:

$$\nabla_{\theta} J(\theta) = \frac{1}{N} \sum_{i=1}^{N} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) R_t \right]$$

- 5. Update the network parameter as $\theta = \theta + \alpha \nabla_{\theta} J(\theta)$
- 6. Repeat *steps* 2 to 5 for several iterations



We start from the gradient in policy gradient with reward-to-go.

$$\nabla_{\theta} J(\theta) = \frac{1}{N} \sum_{i=1}^{N} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} (a_t | s_t) R_t \right]$$

 We introduce a new function b called a baseline function, and subtract the baseline from the reward.

$$\nabla_{\theta} J(\theta) = \frac{1}{N} \sum_{i=1}^{N} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} (a_t | s_t) (R_t - b) \right]$$



- The purpose of the baseline is to reduce the variance in the return.
- Thus, if the baseline b is a value that can give us the expected return from the state the agent is in, then subtracing b in every step will reduce the variance in return.
- A choice for the baseline function is:

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

• We can use other functions as the baseline function as long as it does not depend on the network parameter θ .



- One of the most popular function for baseline is the value function.
- The value function is the expected return an agent would obtain starting from the state following the policy π .
- Using the value function as the baseline, we can write:

$$\nabla_{\theta} J(\theta) = \frac{1}{N} \sum_{i=1}^{N} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta} (a_t | s_t) (R_t - V(s_t)) \right]$$



- Say we use the value function as our baseline function. How can we get the optimal value function?
- We approximate the value function using another neural network parameterized by ϕ . We call this the value network.
- To train the value network, we define the loss function as follows:

$$J(\phi) = \frac{1}{N} \sum_{i=1}^{N} \sum_{t=0}^{T} (R_t - V_{\phi}(s_t))^2$$

• Then, we minimize the error using gradient descent and update the network parameter as:

$$\phi = \phi - \alpha \nabla_{\phi} J(\phi)$$



- The algorithm
- 1. Initialize the policy network parameter θ and value network parameter ϕ
- 2. Generate *N* number of trajectories $\{\tau^i\}_{i=1}^N$ following the policy π_θ
- 3. Compute the return (reward-to-go) R_t
- 4. Compute the policy gradient:

$$\nabla_{\theta} J(\theta) = \frac{1}{N} \sum_{i=1}^{N} \left[\sum_{t=0}^{T-1} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) \left(R_t - V_{\phi}(s_t) \right) \right]$$

- 5. Update the policy network parameter θ using gradient ascent as $\theta = \theta + \alpha \nabla_{\theta} J(\theta)$
- 6. Compute the MSE of the value network:

$$J(\phi) = \frac{1}{N} \sum_{i=1}^{N} \sum_{t=0}^{T-1} \left(R_t - V_{\phi}(s_t) \right)^2$$

- 7. Compute gradients $\nabla_{\phi}J(\phi)$ and update the value network parameter ϕ using gradient descent as $\phi = \phi \alpha \nabla_{\phi}J(\phi)$
- 8. Repeat *steps* 2 to 7 for several iterations



- We implement the policy gradient algorithm for the cart pole balancing task.
- The code includes all three versions:
 - policy gradient
 - policy gradient with reward-to-go
 - policy gradient with baseline
 - baseline: value function



library imports

- namedtuple is data structure where we can record a value with its attribute name. It is used to record the (state, action, reward, logp, done) for all steps in a trajectory.
- matplotlib.pyplot is used to draw a graph.

```
import torch
import gym
from collections import namedtuple
import matplotlib.pyplot as plt
```



PolicyNet: neural network used for policy gradient

```
class PolicyNet(torch.nn.Module):
    def init (self, input size, output size, hidden layer size=64):
        super(PolicyNet, self). init ()
        self.fc1 = torch.nn.Linear(input size, hidden layer size)
        self.fc2 = torch.nn.Linear(hidden layer size, output size)
        self.softmax = torch.nn.Softmax(dim=0)
    def forward(self, x):
       x = torch.from_numpy(x).float()
        return self.softmax(self.fc2(torch.nn.functional.relu(self.fc1(x))))
    def get_action_and_logp(self, x):
        action prob = self.forward(x)
       m = torch.distributions.Categorical(action prob)
        action = m.sample()
        logp = m.log prob(action)
        return action.item(), logp
    def act(self, x):
        action, = self.get action and logp(x)
        return action
```



- ValueNet: neural network used to calculate value function (baseline)
 - The output layer has one neuron, which indicates the value of a state given as input.

```
class ValueNet(torch.nn.Module):
    def __init__(self, input_size, hidden_layer_size=64):
        super(ValueNet, self).__init__()
        self.fc1 = torch.nn.Linear(input_size, hidden_layer_size)
        self.fc2 = torch.nn.Linear(hidden_layer_size, 1)

    def forward(self, x):
        x = torch.from_numpy(x).float()
        return self.fc2(torch.nn.functional.relu(self.fc1(x)))
```



- policyGradient: runs policy gradient algorithm while running episodes.
 - collect_trajectory runs a single episode and collects (state, action, reward, logp, done) in each step. logp is the log probability of selecting the action in the state.

```
def policyGradient(env, max_num_steps=1000, gamma=0.98, lr=0.01,
                   num traj=10, num iter=200):
    input size = env.observation_space.shape[0]
   output_size = env.action_space.n
    Trajectory = namedtuple('Trajectory', 'states actions rewards dones logp')
    def collect_trajectory():
        state list = []
        action list = []
        reward_list = []
        dones list = []
        logp list = []
        state = env.reset()
        done = False
        steps = 0
        while not done and steps <= max_num_steps:
            action, logp = policy.get_action_and_logp(state)
            newstate, reward, done, _ = env.step(action)
            state list.append(state)
            action list.append(action)
            reward list.append(reward)
            dones list.append(done)
            logp_list.append(logp)
            steps += 1
            state = newstate
        traj = Trajectory(states=state list, actions=action list,
                          rewards=reward list, logp=logp_list, dones=dones_list)
        return traj
```



- policyGradient: runs policy gradient algorithm while running episodes.
 - calc_returns calculates (discounted) R_t for each step in the episode

```
def calc_returns(rewards):
    dis_rewards = [gamma**i * r for i, r in enumerate(rewards)]
    return [sum(dis_rewards[i:]) for i in range(len(dis_rewards))]
```

- We prepare a policy network and a value network. Both needs training, so we use the Adam optimizer to train the networks.
 - Value network is only needed when we use the baseline

```
policy = PolicyNet(input_size, output_size)
policy_optimizer = torch.optim.Adam(policy.parameters(), lr=lr)

value = ValueNet(input_size)
value_optimizer = torch.optim.Adam(value.parameters(), lr=lr)
```



loss terms for the three versions of policy gradient

```
mean return list = []
for it in range(num_iter):
   traj_list = [collect_trajectory() for _ in range(num_traj)]
   returns = [calc_returns(traj.rewards) for traj in traj list]
   #-----#
   # policy gradient with base function #
   #========#
   #policy loss terms = [-1. * traj.logp[i] * (returns[i][i] - value(traj.states[i]))
                    for i, traj in enumerate(traj_list) for j in range(len(traj.actions))]
   #=======#
   # policy gradient with reward-to-go #
   #=======#
   #policy loss terms = [-1. * traj.logp[j] * (torch.Tensor([returns[i][j]]))
                    for i, traj in enumerate(traj list) for j in range(len(traj.actions))]
   # policy gradient
   #=======#
   policy loss terms = [-1. * traj.logp[j] * (torch.Tensor([returns[i][0]]))
                    for i, traj in enumerate(traj_list) for j in range(len(traj.actions))]
```



Updating parameters of neural networks through backpropagation

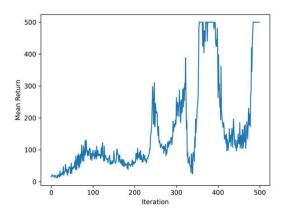
Saving mean returns and printing intermediate results



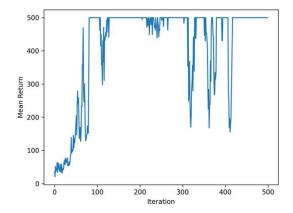
- Running the policy gradient algorithm to solve the Cart Pole task
 - _max_episode_steps is maximum steps the environment allows in an episode.

```
env = gym.make('CartPole-v1')
env._max_episode_steps=500
agent, mean_return_list = policyGradient(env, num_iter=500, max_num_steps=500, gamma=1.0, num_traj=5)

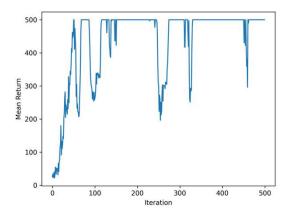
plt.plot(mean_return_list)
plt.xlabel('Iteration')
plt.ylabel('Mean Return')
plt.savefig('pg_returns.png', format='png', dpi=300)
```



policy gradient



policy gradient with reward-to-go



policy gradient with baseline



Summary

- While DQN is a value-based method, policy gradient is a policy-based RL method.
- In policy gradient, we do not calculate Q values. Instead, the neural network outpus probability distribution of actions in a state.
- To train the policy network, we run episodes with stochastic policy. If the agent obtains a positive reward in an episode, probability of the actions taken is increased. If the agent obtains a negative reward, probability of the actions is decreased.
- Since the original policy gradient method can lead to high variance of gradients, which can make the training take long time to converge.
- We can use two improved algorithms: policy gradient with reward-to-go and policy gradient with baseline.
- In policy gradient with baseline, we use another neural network to approximate value function, and use it as the baseline.



End of Class

Questions?

Email: jso1@sogang.ac.kr

