COMP3667: Reinforcement learning practical 4

Dynamic Programming

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This is the version with answers!

1 Overview

named actions

LEFT, DOWN, RIGHT, UP = 0, 1, 2, 3

Welcome to the fourth reinforcement learning practical. In this practical, we will dive a bit more into the practical details of dynamic programming methods. In particular, we will:

- Use the OpenAI Gym "Frozen Lake" environment as a basic example.
- Implement policy evaluation and understand how this algorithm updates state values.
- Implement policy improvement to learn better policies.
- Implement *policy iteration* to learn optimal policies.

For this practical, you will need a basic Python environment with numpy, matplotlib, and OpenAI gym (version 0.20.0).

You can download the Jupyter notebook here or directly open the Colab notebook.

2 Frozen Lake Environment

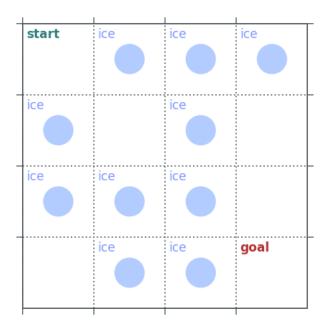
```
Install OpenAI gym if you have not done this already. You will need to use version 0.20.0:
                                                                                          </> copy code
   %%capture
    !pip install setuptools==65.5.0 "wheel<0.40.0"
    !apt update
   !pip install 'gym==0.20.0'
   Do the relevant imports
                                                                                          </> copy code
   import gym
    import numpy as np
    import matplotlib.pyplot as plt
    import matplotlib.font_manager
   and load the Frozen Lake environment
                                                                                          </> copy code
   name = 'FrozenLake-v1' # small version
    # name = 'FrozenLake8x8-v1' # larger version
    env = gym.make(name, is_slippery=False)
3
    env.seed(742)
    env.action_space.seed(742)
```

```
def plot(env, v=None, policy=None, col_ramp=1, dpi=175, draw_vals=False, mark_ice=True):
1
        # set up plot
2
        plt.rcParams['figure.dpi'] = dpi
        plt.rcParams.update({'axes.edgecolor': (0.32,0.36,0.38)})
        plt.rcParams.update({'font.size': 4 if env.env.nrow == 8 else 7})
5
        gray = np.array((0.32,0.36,0.38))
6
        plt.figure(figsize=(3, 3))
        ax = plt.gca()
        ax.set_xticks(np.arange(env.env.ncol)-.5)
9
        ax.set_yticks(np.arange(env.env.nrow)-.5)
10
        ax.set_xticklabels([])
        ax.set_yticklabels([])
12
        plt.grid(color=(0.42,0.46,0.48), linestyle=':')
13
        ax.set_axisbelow(True)
        ax.tick_params(color=(0.42,0.46,0.48),
15
                       which='both', top='off', left='off', right='off', bottom='off')
16
        # use zero value as dummy if not provided
17
        if v is None:
            v = np.zeros(env.nS)
19
        # plot values
20
        plt.imshow(1-v.reshape(env.env.nrow,env.env.ncol)**col_ramp,
21
                   cmap='gray', interpolation='none',
22
                   clim=(0,1), zorder=-1)
23
        # go through states
24
        for s in range(env.nS):
25
            x, y = s % env.env.nrow, s // env.env.ncol
            # print numeric values
27
            if draw_vals and v[s] > 0:
28
                vstr = '\{0:.1e\}'.format(v[s]) if env.env.nrow == 8 else '\{0:.6f\}'.format(v[s])
29
                plt.text(x - 0.45, y + 0.45, vstr, color=(0.2, 0.8, 0.2), fontname='Sans')
            # mark ice, start, goal
31
            if env.desc.tolist()[y][x] == b'F':
32
                plt.text(x-0.45,y-0.3, 'ice', color=(0.5, 0.6, 1), fontname='Sans')
                 if mark_ice:
34
                     ax.add_patch(plt.Circle((x, y), 0.2, color=(0.7, 0.8, 1), zorder=0))
35
            elif env.desc.tolist()[y][x] == b'S':
36
                plt.text(x-0.45,y-0.3, 'start',color=(0.2,0.5,0.5), fontname='Sans',
37
                          weight='bold')
38
            elif env.desc.tolist()[y][x] == b'G':
39
                plt.text(x-0.45,y-0.3, 'goal', color=(0.7,0.2,0.2), fontname='Sans',
40
                          weight='bold')
                 continue # don't plot policy for goal state
            else:
43
                 continue # don't plot policy for holes
44
            # plot policy
            def plot_arrow(x, y, dx, dy, v, scale=0.4):
46
                plt.arrow(x, y, scale * float(dx), scale * float(dy), color=gray+0.2*(1-v),
47
                           head_width=0.1, head_length=0.1, zorder=1)
            if policy is not None:
                 a = policy[s]
50
                 if a[0] > 0.0: plot_arrow(x, y, -a[0],
                                                            0., v[s]) # left
51
                if a[1] > 0.0: plot_arrow(x, y,
                                                  0., a[1], v[s]) # down
52
                 if a[2] > 0.0: plot_arrow(x, y, a[2],
                                                          0., v[s]) # right
                if a[3] > 0.0: plot_arrow(x, y,
                                                   0., -a[3], v[s]) # up
54
        plt.show()
55
```

Have a look

```
print('action space: ' + str(env.action_space))
print('reward range: ' + str(env.reward_range))
print('observation space: ' + str(env.observation_space))
```

```
plot(env=env)
--> action space: Discrete(4)
--> reward range: (0, 1)
--> observation space: Discrete(16)
```

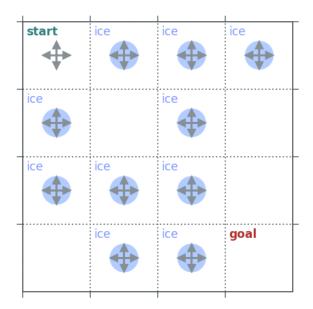


• Define a uniform policy and plot it in the environment. *Hint:* env.nS and env.nA give you the number of states and actions, respectively; you can provide the policy to the plotting function via the policy argument.

Answer:

</pre

```
def uniform_policy(env):
    return np.ones((env.nS, env.nA)) / env.nA
plot(env=env, policy=uniform_policy(env))
```



3 Policy Evaluation

Now, we would like to know how well a policy performs, that is, what the state value of a particular state is when following the policy in the future. This *policy evaluation* procedure can be efficiently done using dynamic

programming, which iteratively improves state value estimates and converges to the true state values.

• Write a policy_eval_step function that takes an initial state value estimate $v_{\pi}^{k}(s)$ and computes an improved estimate $v_{\pi}^{k+1}(s)$ as

$$v_{\pi}^{k+1}(s) = \sum_{a} \pi(a \mid s) \sum_{s'} p(s' \mid s, a) \left[\mathcal{R}_{ss'}^{a} + \gamma v_{\pi}^{k}(s') \right] . \tag{1}$$

Hint: env.env.P[s][a]: is giving you a list of tuples (p, s', r, done), one for each possible transition from state s when taking action a: p is the probability of this transition to happen, s' is the state the agent transitions to, r is the reward it receives, and done indicates whether the episode is over (which you will not need to use).

Answer: ⟨/> copy code

```
def policy_eval_step(env, policy, gamma, v_init=None):
    if v_init is None:
        v_init = np.zeros(env.nS)

v = np.zeros(env.nS)

for s_from in range(env.nS):
    for a in range(env.nA):
        pi = policy[s_from, a]

for p, s_to, r, done in env.env.P[s_from][a]:
        v[s_from] += pi * p * (r + gamma * v_init[s_to])

return v
```

• Initialise the state values with zero, take one policy_eval_step at a time and plot the result to observe how updates are being performed. *Hint:* You can provide the state values to the plot function using the v argument, specifying draw_vals=True additionally shows the numeric state values.

First, initialise the state values:

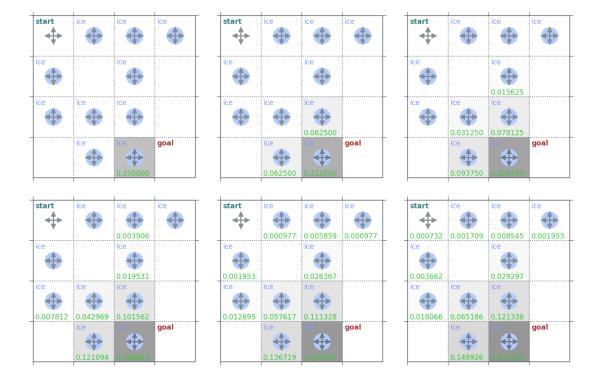
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```
v = np.zeros(env.nS)
```

Then repeatedly execute the update step:

</> copy code

```
v = policy_eval_step(env, uniform_policy(env), 1, v)
plot(env, v, uniform_policy(env), draw_vals=True)
```



• On an intuitive level, how would you describe the dynamics you see? What seems to be inefficient about the current implementation? How could this be improved?

The environment is "flooded" with values from the goal state (or, more generally, any reward transition); values "propagate" one step with each iteration. In the first iterations, most state values do not change (they remain zero) and would not need to be updated, but the current implementation always updates *all* state values, which seems inefficient. If we knew which state values change, we could only update those.

- Implement a modified policy_eval_step_inplace version of the function that performs state value updates in-place. That is, instead of clearly separating the "old" values $v_{\pi}^{k}(s)$ and the "new" values $v_{\pi}^{k+1}(s)$, you operate on a single state value estimate $v_{\pi}(s)$, which is updated as you go. *Hint:* Make sure the updates for a particular state are still "atomic" and you do not use the half-computed values (in case of transitions that stay in a particular state).
 - Think about a clever order in which to update state values in-place. *Hint:* States are ordered from top-left (start: 0) to bottom-right (goal: 15).
 - In the modified implementation, the order of states is reversed because this makes updates more efficient as newly computed values will directly be reused in the following updates (the update direction is the same as the "flow" of state values from goal state to start state; this is obviously specific to this environment and not a general recipe).
 - Again, observe the step-wise update of values from one iteration to the next. How does that compare to the original implementation?

Already in the first iteration, values propagate all the way to the start state.

Answer: ⟨/> copy code

```
def policy_eval_step_inplace(env, policy, gamma, v_init=None):
        if v_init is None:
            v_init = np.zeros(env.nS)
3
        v = v_init.copy() # opearate on copy in-place
        for s_from in reversed(range(env.nS)): # reverse order of states
            v_s_from = 0 # compute value for this state
            for a in range(env.nA):
                pi = policy[s_from, a]
                for p, s_to, r, done in env.env.P[s_from][a]:
                    v_sfrom += pi * p * (r + gamma * v[s_to]) # use the values we also update
10
            v[s_from] = v_s_from # update
11
        return v
12
```



• Write a policy_evaluation function that iteratively updates the state values using the policy_eval_step or policy_eval_step_inplace function and stops if they do not change (by some small tolerance value). Print the number of iterations needed to converge and compare for the policy_eval_step and policy_eval_step_inplace implementation.

Answer:

```
v = v_init
for i in range(1, max_iter + 1):
    new_v = policy_eval_step(env, policy, gamma, v)
    if np.allclose(v, new_v, atol=atol):
        break
    v = new_v
    if print_iter:
        print(f"{i} iterations")
    return v
```

• How many iterations do you need until state values have converged to their true value? To make it simpler, do the following though experiment: Take an environment that has only a single state and a single action (so nothing can really change and there is only one possible policy) and you get a reward of 1 upon every transition. Take the update equation for the state value from above, which now simplifies to

$$v^{k+1} = 1 + \gamma v^k \tag{2}$$

Can you write down v^n in a non-recursive form assuming you start with $v^0 = 0$? Can you write down v^{∞} in closed form? Is v^{∞} the exact state value? How long does it take to converge? What happens for $\gamma = 1$ as opposed to $\gamma < 1$?

Starting with $v^0 = 0$, v^n can be "unrolled" and rewritten as

$$v^n = 1 + \gamma v^{n-1} \tag{3}$$

$$=1+\gamma\left(1+\gamma\,v^{n-2}\right)\tag{4}$$

$$=\cdots$$
 (5)

$$=1+\gamma(1+\gamma\cdots(1+\gamma v^0)) \tag{6}$$

$$=1+\gamma+\gamma^2\cdots(1+\gamma^2v^0)\tag{7}$$

$$= 1 + \gamma + \gamma^{2} + \dots + \gamma^{n-1} + \gamma^{n} v^{0}$$
 (8)

$$= \sum_{k=0}^{n-1} \gamma^k \qquad \text{(non-recursive; geometric series)} \tag{9}$$

$$= \frac{1 - \gamma^n}{1 - \gamma} \qquad \text{(closed form formula for } \gamma < 1) \tag{10}$$

$$\Rightarrow \qquad v^{\infty} = \frac{1}{1 - \gamma} \tag{11}$$

Inserting v^{∞} into the update equation gives

$$\frac{1}{1-\gamma} = 1 + \gamma \frac{1}{1-\gamma} \tag{12}$$

$$\Leftrightarrow 1 = 1 - \gamma + \gamma \tag{13}$$

$$\Leftrightarrow 1 = 1 \tag{14}$$

(10)

(16)

(15)

so v^{∞} is indeed the exact state value. The values converge exponentially fast (the γ^n term decays at an exponential rate), but will never be exact if there are paths of infinite length in the environment (typically loops; like the self-loop in this minimal though experiment). If only finite-length paths exist, the series can be truncated and values will be exact after a finite number of iterations.

4 Policy Improvement

• Implement a function that computes state-action values $q_{\pi}(s, a)$ (for all actions in a given state) from the state values $v_{\pi}(s)$ using their known relation

$$q_{\pi}(s, a) = \sum_{s' \in \mathcal{S}} \mathcal{P}_{ss'}^{a} \left[\mathcal{R}_{ss'}^{a} + \gamma \, v_{\pi}(s') \right] . \tag{17}$$

Answer:

• Implement a policy_improvement function that takes a state value estimate and defines a policy (by first computing state-action values) that achieves maximal return (i.e. always chooses an action with maximal state-action value). Optionally, either choose actions deterministically, or choose all actions with maximum value with the same probability (if there is more than one such action).

Answer: ⟨/> copy code

```
def policy_improvement(env, v, gamma, deterministic=False):
    policy = np.zeros([env.nS, env.nA]) / env.nA
    for s in range(env.nS):
        q = q_from_v(env, v, s, gamma)
        if deterministic:
        # deterministic policy
        policy[s][np.argmax(q)] = 1
    else:
        # stochastic policy with equal probability on maximizing actions
        best_a = np.argwhere(q==np.max(q)).flatten()
        policy[s, best_a] = 1 / len(best_a)
    return policy
```

• Load the larger 'FrozenLake8x8-v1' environment (again with is_slippery=False for now), compute state values by evaluating the uniform policy, plot the result. Then compute an improved policy and plot the result again.

Load and plot the larger environment with uniform policy:

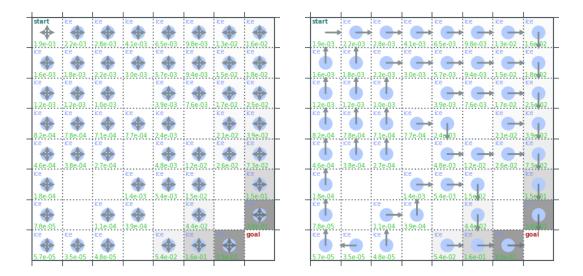
⟨⟩ copy code

```
env = gym.make('FrozenLake8x8-v1', is_slippery=False)
env.seed(742)
env.action_space.seed(742)
policy = uniform_policy(env)
v = policy_evaluation(env, policy, gamma=1)
plot(env, v, policy, draw_vals=True)
```

Compute improved policy and plot again:

⟨⟩ copy code

```
new_policy = policy_improvement(env, v, gamma=1)
plot(env, v, new_policy, draw_vals=True)
```



5 Policy Iteration

- Starting from the improved policy from above, perform two updates by doing
 - policy evaluation (plot the results)
 - policy improvement (plot the results).

Use a stochastic policy improvement (i.e. choose optimal action with equal probability) and a discount value of $\gamma = 1$. What do you observe? To you see anything that could be problematic? Can you explain what you observe (you may need to print the raw state values)? Would a deterministic policy improvement help? *Hint:* Remember what we learned above about value convergence and paths of finite/infinite length.

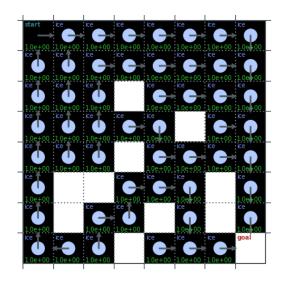
First update: The improved policy is deterministic and during evaluation state values of 1 propagate through the entire state space. In the policy improvement step we get an almost uniform policy (except it does not jump into holes) because state values are equal everywhere.

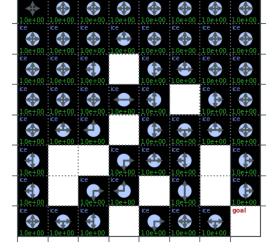
Second update: The state values from the (almost) uniform policy do not converge to their exact value of 1, because we have infinite-length paths (loops). Therefore, in the policy improvement step, we again get a deterministic policy (moving towards the goal state, because these state values have "seen" the reward in earlier iterations and are therefore converged more closely to the exact value of 1).

Potential problems: We do not get a stable policy, but instead alternate between the deterministic and stochastic version with state values being exactly 1 in the deterministic case and only approximately 1 in the stochastic case. This can be problematic when trying to detect convergence (as both policy and values keep on changing).

Deterministic policy: A deterministic policy improvement step **cannot** solve this problem (we may still get loops if we are unlucky). In the contrary, since all state values are the same, the deterministic policy may choose arbitrary actions (as long as they do not jump into a hole), which in total may lead to an arbitrarily bad policy.

```
gamma = 1
v = policy_evaluation(env, new_policy, gamma=gamma)
plot(env, v=v, policy=new_policy, draw_vals=True)
print(v)
new_policy = policy_improvement(env, v, gamma=gamma)
plot(env, v=v, policy=new_policy, draw_vals=True)
```

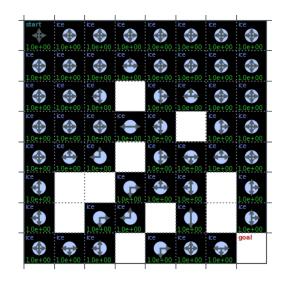


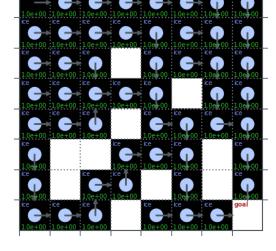


Policy Evaluation #1

Policy Improvement #1

State values:





Policy Evaluation #2

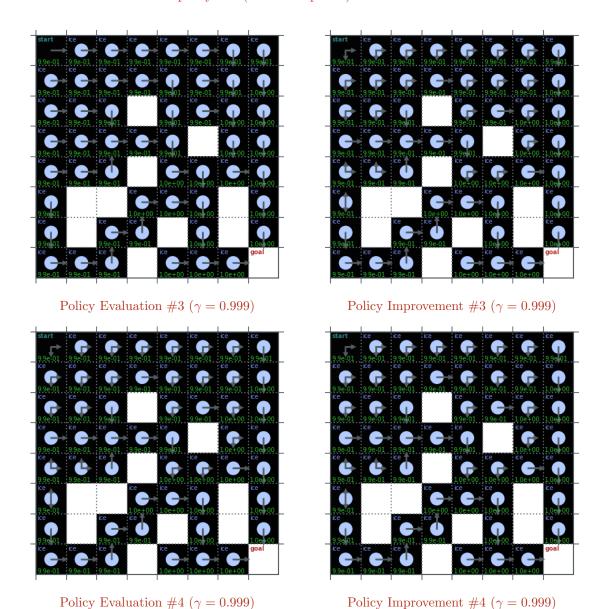
Policy Improvement #2

State values:

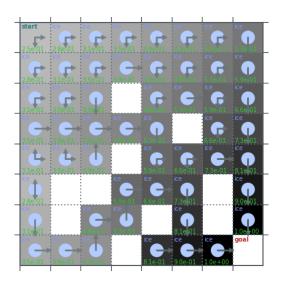
```
 \begin{bmatrix} 0.99584585 & 0.99588645 & 0.99596896 & 0.99609428 & 0.99624201 & 0.996371 \\ 0.99647678 & 0.99653576 & 0.99584504 & 0.99588397 & 0.99596476 & 0.99610928 \\ 0.99629677 & 0.99642897 & 0.99655735 & 0.99662794 & 0.99584513 & 0.99587906 \\ 0.99593551 & 0. & 0.9964423 & 0.99652499 & 0.99672868 & 0.99682301 \\ 0.99585109 & 0.99589112 & 0.99599191 & 0.99625233 & 0.9965307 & 0. \\ 0.9970407 & 0.99714287 & 0.99585676 & 0.9958818 & 0.9959271 & 0. \\ 0.9969224 & 0.99720212 & 0.99727183 & 0.99759226 & 0.99587711 & 0. \\ 0. & 0.99683596 & 0.99705648 & 0.99743225 & 0. & 0.99838516 \\ 0.99592708 & 0. & 0.99644137 & 0.99663059 & 0. & 0.9980566 \\ 0. & 0.99918967 & 0.99600632 & 0.99612383 & 0.9962692 & 0. \\ 0.99868395 & 0.99869025 & 0.99934277 & 0. & \end{bmatrix}
```

• Change the value to $\gamma = 0.999$ and do another two updates. What is different? *Hint:* You can use $\gamma = 0.9$ to see the effect more clearly.

State values are almost the same, but a value of $\gamma < 1$ penalises moving in circles "for no reason". Therefore, we get a policy the is moving towards the goal state as quickly as possible without creating any loop (even though there are sometimes multiple equally-valued paths, so the policy is not deterministic). As a result we obtain a stable policy and (as a consequence) stable state values.



10



Converged Policy/Values ($\gamma = 0.9$)