Assignment 1: Markov Decision Processes

Due Date: October 15th (Fri), 11:59 pm.

Homework Instructions

- All your answers should be written in this notebook.
- You shouldn't need to write or modify any other files.
- Look for four instances between `BEGIN YOUR CODE` and `END YOUR CODE` --those are the only parts of the code you need to write. To grade your homework, we will check whether the printouts immediately following your code match up with the results we got. The portions used for grading are highlighted in yellow. (However, note that the yellow highlighting does not show up when github renders this file.)

To submit your homework, upload a PDF version of this file (To make the pdf, do `File - Print Preview`)

Score Breakdown

Question	Points
Question 1	6
Question 2	4
Question 3a	6
Question 3b	4
Total	20

Introduction

This assignment will review the two classic methods for solving Markov Decision Processes (MDPs) with finite state and action spaces. We will implement value iteration (VI) and policy iteration (PI) for a finite MDP, both of which find the optimal policy in a finite number of iterations.

The experiments here will use the Frozen Lake environment, a simple gridworld MDP that is taken from `gym` and slightly modified for this assignment. In this MDP, the agent must navigate from the start state to the goal state on a 4x4 grid, with stochastic transitions.

In [1]:
```
#!pip install gym  # uncomment if you haven't installed the gym library

from frozen_lake import FrozenLakeEnv
env = FrozenLakeEnv()
env.seed(0);
print(env.__doc__)

import numpy as np, numpy.random as nr, gym
np.set_printoptions(precision=3)
def begin_grading(): print("\x1b[43m")
def end_grading(): print("\x1b[0m")
```

```
    Winter is here. You and your friends were tossing around a frisbee at t
he park
    when you made a wild throw that left the frisbee out in the middle of t
he lake.
    The water is mostly frozen, but there are a few holes where the ice has
melted.
    If you step into one of those holes, you'll fall into the freezing wate
r.
    At this time, there's an international frisbee shortage, so it's absolu
tely imperative that
    you navigate across the lake and retrieve the disc.
    However, the ice is slippery, so you won't always move in the direction
you intend.
    The surface is described using a grid like the following

        SFFF
        FHFH
        FFFH
        HFFG

    S : starting point, safe
    F : frozen surface, safe
    H : hole, fall to your doom
    G : goal, where the frisbee is located

    The episode ends when you reach the goal or fall in a hole.
    You receive a reward of 1 if you reach the goal, and zero otherwise.
```

We extract the relevant information from the gym Env into the MDP class below. The `env` object won't be used any further, we'll just use the `mdp` object.

In [2]:

```
class MDP(object):
    def __init__(self, P, nS, nA, desc=None):
        self.P = P # state transition and reward probabilities, explained b
        self.nS = nS # number of states
        self.nA = nA # number of actions
        self.desc = desc # 2D array specifying what each grid cell means (u
mdp = MDP( {s : {a : [tup[:3] for tup in tups] for (a, tups) in a2d.items()


print("mdp.P is a two-level dict where the first key is the state and the s
print("The 2D grid cells are associated with indices [0, 1, 2, ..., 15] fro
print(np.arange(16).reshape(4,4))
print("mdp.P[state][action] is a list of tuples (probability, nextstate, re
print("For example, state 0 is the initial state, and the transition inform
print("As another example, state 5 corresponds to a hole in the ice, which
print("P[5][0] =", mdp.P[5][0], '\n')
```

```
mdp.P is a two-level dict where the first key is the state and the second k
ey is the action.
The 2D grid cells are associated with indices [0, 1, 2, ..., 15] from left
to right and top to down, as in
[[ 0  1  2  3]
 [ 4  5  6  7]
 [ 8  9 10 11]
 [12 13 14 15]]
mdp.P[state][action] is a list of tuples (probability, nextstate, reward).

For example, state 0 is the initial state, and the transition information f
or s=0, a=0 is
P[0][0] = [(0.1, 0, 0.0), (0.8, 0, 0.0), (0.1, 4, 0.0)]

As another example, state 5 corresponds to a hole in the ice, which transit
ions to itself with probability 1 and reward 0.
P[5][0] = [(1.0, 5, 0)]
```

Part 1: Value Iteration

Question 1: implement value iteration

In this problem, you'll implement value iteration, which has the following pseudocode:

Initialize $V^{(0)}(s) = 0$, for all s

For $i = 0, 1, 2, \ldots$

- $V^{(i+1)}(s) = \max_a \sum_{s'} P(s, a, s')[R(s, a, s') + \gamma V^{(i)}(s')]$, for all s

---

We additionally define the sequence of greedy policies $\pi^{(0)}, \pi^{(1)}, \ldots, \pi^{(n-1)}$, where

$$\pi^{(i)}(s) = \arg\max_a \sum_{s'} P(s, a, s')[R(s, a, s') + \gamma V^{(i)}(s')]$$

Your code will return two lists: $[V^{(0)}, V^{(1)}, \ldots, V^{(n)}]$ and $[\pi^{(0)}, \pi^{(1)}, \ldots, \pi^{(n-1)}]$

To ensure that you get the same policies as the reference solution, choose the lower-index action to break ties in $\arg\max_a$. This is done automatically by np.argmax. This will only affect the "# chg actions" printout below--it won't affect the values computed.

Warning: make a copy of your value function each iteration and use that copy for the update--don't update your value function in place. Updating in-place is also a valid algorithm, sometimes called Gauss-Seidel value iteration or asynchronous value iteration, but it will cause you to get different results than me.

In [3]:

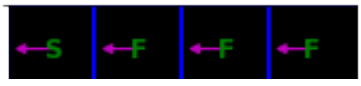
```
def value_iteration(mdp, gamma, nIt):
    """
    Inputs:
        mdp: MDP
        gamma: discount factor
        nIt: number of iterations, corresponding to n above
    Outputs:
        (value_functions, policies)

    len(value_functions) == nIt+1 and len(policies) == n
    """
    print("Iteration | max|V-Vprev| | # chg actions | V[0]")
    print("----------+--------------+---------------+---------")
    Vs = [np.zeros(mdp.nS)] # list of value functions contains the initial
    pis = []
    for it in range(nIt):
        Vprev = Vs[-1]
        oldpi = pis[-1] if len(pis) > 0 else None

        V = np.zeros(mdp.nS)  # Initialize V
        pi = np.zeros(mdp.nS)  # update states & policy

        # BEGIN YOUR CODE
        # ------------------------
        for state in range(mdp.nS):
            s_action=[]
            for a in range(mdp.nA):
                q=[]
                for next_state in range(len(mdp.P[state][a])):
                    q.append((mdp.P[state][a][next_state][0]*(mdp.P[state][
                s_action.append(sum(q))

            V[state]=(max(s_action))  # Value

            pi[state]=np.argmax(s_action)  # Policy
        # ------------------------
        # END YOUR CODE

        # Your code should define variables V: the bellman backup applied t
        # and pi: the greedy policy applied to Vprev
        max_diff = np.abs(V - Vprev).max()
        nChgActions="N/A" if oldpi is None else (pi != oldpi).sum()
        print("%4i      | %6.5f      | %4s          | %5.3f"%(it, max_diff,
        Vs.append(V)
        pis.append(pi)
    return Vs, pis

GAMMA=0.95 # we'll be using this same value in subsequent problems
begin_grading()
Vs_VI, pis_VI = value_iteration(mdp, gamma=GAMMA, nIt=20)
end_grading()
```

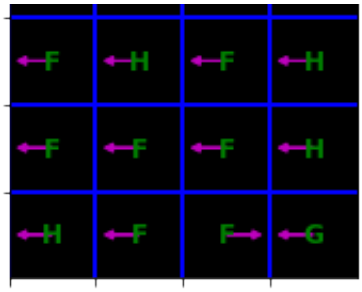
```
Iteration | max|V-Vprev| | # chg actions | V[0]
----------+--------------+---------------+---------
   0      | 0.80000      |  N/A          | 0.000
   1      | 0.60800      |    2          | 0.000
   2      | 0.51984      |    2          | 0.000
   3      | 0.39508      |    2          | 0.000
   4      | 0.30026      |    2          | 0.000
   5      | 0.25355      |    1          | 0.254
   6      | 0.10478      |    0          | 0.345
   7      | 0.09657      |    0          | 0.442
   8      | 0.03656      |    0          | 0.478
   9      | 0.02772      |    0          | 0.506
  10      | 0.01111      |    0          | 0.517
  11      | 0.00735      |    0          | 0.524
  12      | 0.00310      |    0          | 0.527
  13      | 0.00190      |    0          | 0.529
  14      | 0.00083      |    0          | 0.530
  15      | 0.00049      |    0          | 0.531
  16      | 0.00022      |    0          | 0.531
  17      | 0.00013      |    0          | 0.531
  18      | 0.00006      |    0          | 0.531
  19      | 0.00003      |    0          | 0.531
```

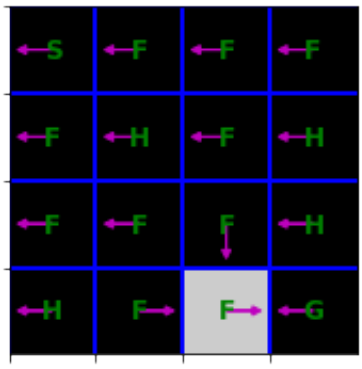
Below, we've illustrated the progress of value iteration. Your optimal actions are shown by arrows. At the bottom, the value of the different states are plotted.

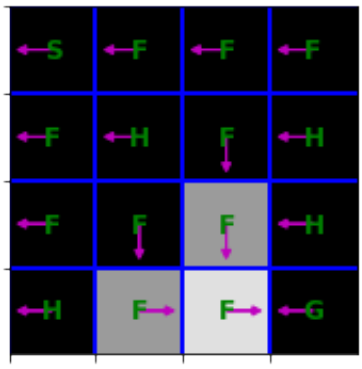
In [4]:

```
import matplotlib.pyplot as plt
%matplotlib inline
for (V, pi) in zip(Vs_VI[:10], pis_VI[:10]):
    plt.figure(figsize=(3,3))
    plt.imshow(V.reshape(4,4), cmap='gray', interpolation='none', clim=(0,1
    ax = plt.gca()
    ax.set_xticks(np.arange(4)-.5)
    ax.set_yticks(np.arange(4)-.5)
    ax.set_xticklabels([])
    ax.set_yticklabels([])
    Y, X = np.mgrid[0:4, 0:4]
    a2uv = {0: (-1, 0), 1:(0, -1), 2:(1,0), 3:(-1, 0)}
    Pi = pi.reshape(4,4)
    for y in range(4):
        for x in range(4):
            a = Pi[y, x]
            u, v = a2uv[a]
            plt.arrow(x, y,u*.3, -v*.3, color='m', head_width=0.1, head_len
            plt.text(x, y, str(env.desc[y,x].item().decode()),
                     color='g', size=12,  verticalalignment='center',
                     horizontalalignment='center', fontweight='bold')
    plt.grid(color='b', lw=2, ls='-')
plt.figure()
plt.plot(Vs_VI)
plt.title("Values of different states");
```




F H F H
F F F H
H F F G


S F F F
F H F H
F F F H
H F F G


S F F F
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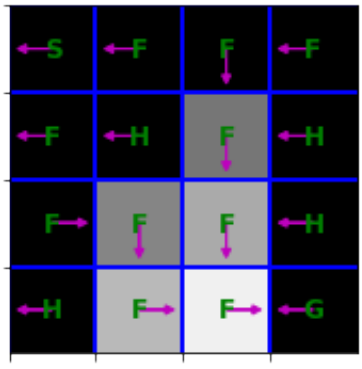
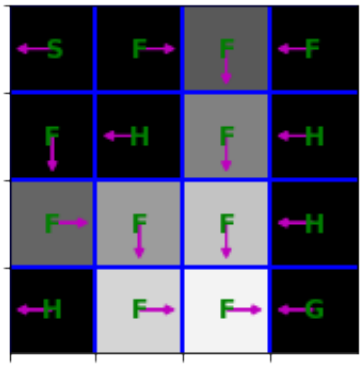
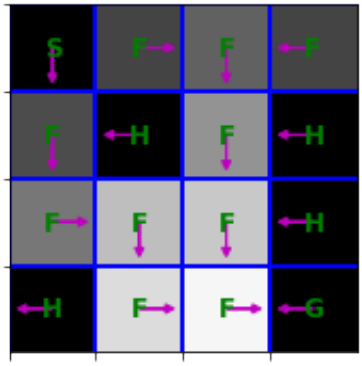
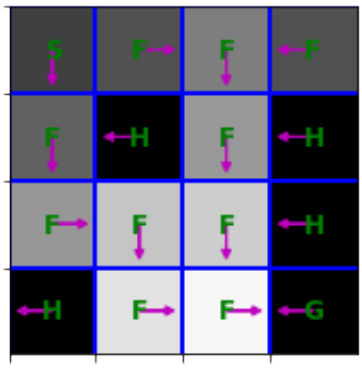
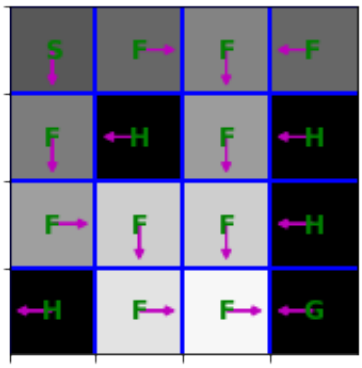
H F F G

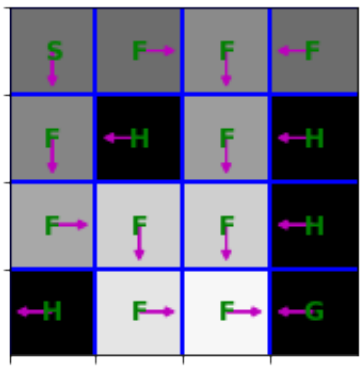

S F F F
F H F H
F F F H
H F F G

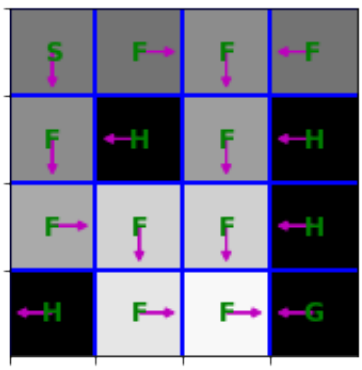

S F F F
F H F H
F F F H
H F F G

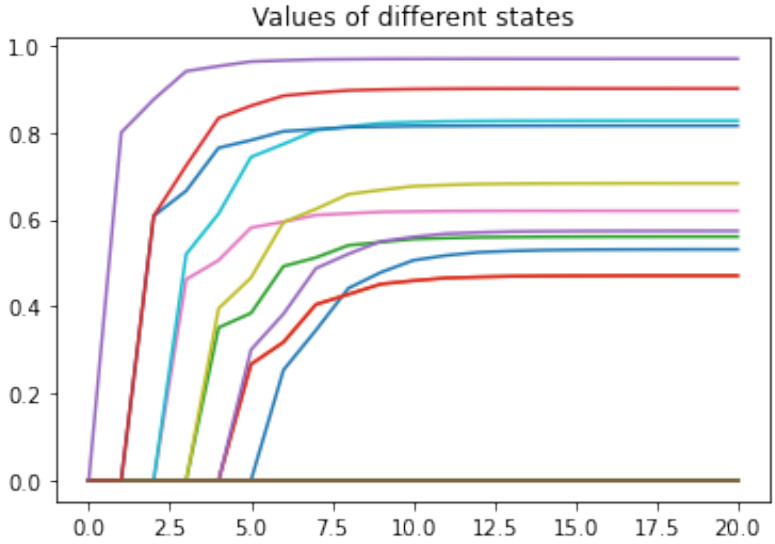

S F F F
F H F H
F F F H
H F F G


S F F F
F H F H
F F F H
H F F G


S F F F
F H F H
F F F H
H F F G


S F F F
F H F H
F F F H
H F F G


S F F F
F H F H
F F F H
H F F G



Question 2: construct an MDP where value iteration takes a long time to converge

When we ran value iteration on the frozen lake problem, the last iteration where an action changed was iteration 6--i.e., value iteration computed the optimal policy at iteration 6. Are there any guarantees regarding how many iterations it'll take value iteration to compute the optimal policy? There are no such guarantees without additional assumptions--we can construct the MDP in such a way that the greedy policy will change after arbitrarily many iterations.

Your task: define an MDP with at most 3 states and 2 actions, such that when you run value iteration, the optimal action changes at iteration >= 50. Use discount=0.95. (However, note that the discount doesn't matter here--you can construct an appropriate MDP with any discount.)

In [5]:

```
chg_iter = 50

# BEGIN YOUR CODE
# ----------------------
nS = 3
nA = 2

P = {}
P[0] = {}  # State 0
P[0][0] = []
P[0][0].append((1,1,1))
P[0][1] = []
P[0][1].append((1,0,0))

P[1] = {}  # State 1
P[1][0] = []
P[1][0].append((1,1,0))
P[1][1] = []
P[1][1].append((1,2,-17.5))#This reward affect the optimal action. So when

P[2] = {}  # State 2
P[2][0] = []
P[2][0].append((1,2,1))
P[2][1] = []
P[2][1].append((1,0,0))
# ----------------------
# END YOUR CODE

mymdp = MDP(P, nS, nA)

begin_grading()
Vs, pis = value_iteration(mymdp, gamma=GAMMA, nIt=chg_iter+1)
end_grading()
```

Iteration	max|V-Vprev|	# chg actions	V[0]
0	1.00000	N/A	1.000
1	0.95000	0	1.000
2	0.90250	0	1.000
3	0.85737	0	1.000
4	0.81451	0	1.000
5	0.77378	0	1.000
6	0.73509	0	1.000
7	0.69834	0	1.000
8	0.66342	0	1.000
9	0.63025	0	1.000
10	0.59874	0	1.000
11	0.56880	0	1.000
12	0.54036	0	1.000
13	0.51334	0	1.000
14	0.48767	0	1.000
15	0.46329	0	1.000
16	0.44013	0	1.000
17	0.41812	0	1.000
18	0.39721	0	1.000
19	0.37735	0	1.000
20	0.35849	0	1.000
21	0.34056	0	1.000
22	0.32353	0	1.000
23	0.30736	0	1.000
24	0.29199	0	1.000
25	0.27739	0	1.000
26	0.26352	0	1.000
27	0.25034	0	1.000
28	0.23783	0	1.000
29	0.22594	0	1.000
30	0.21464	0	1.000
31	0.20391	0	1.000
32	0.19371	0	1.000
33	0.18403	0	1.000
34	0.17482	0	1.000
35	0.16608	0	1.000
36	0.15778	0	1.000
37	0.14989	0	1.000
38	0.14240	0	1.000
39	0.13528	0	1.000
40	0.12851	0	1.000
41	0.12209	0	1.000
42	0.11598	0	1.000
43	0.11018	0	1.000
44	0.10467	0	1.000
45	0.09944	0	1.000
46	0.09447	0	1.000
47	0.08974	0	1.000
48	0.08526	0	1.000
49	0.08099	0	1.000
50	0.07694	1	1.000

Question 3: Policy Iteration

The next task is to implement exact policy iteration (PI), which has the following pseudocode:

Initialize π_0

For $n = 0, 1, 2, \ldots$

1. Compute the state-value function V^{π_n}.
2. Compute the state-action-value function Q^{π_n} using the state-value function.
3. Compute new policy $\pi_{n+1}(s) = \text{argmax}_a \, Q^{\pi_n}(s, a)$.

Below, you'll implement the first and second steps of the loop.

Question 3a: state value function

You'll write a function called `compute_vpi` that computes the state-value function V^π for an arbitrary policy π. Recall that V^π satisfies the following linear equation:

$$V^\pi(s) = \sum_{s'} P(s, \pi(s), s')[R(s, \pi(s), s') + \gamma V^\pi(s')]$$

You'll have to solve a linear system in your code.

In [6]:

```
def compute_vpi(pi, mdp, gamma):
    V = np.zeros(mdp.nS)

    # BEGIN YOUR CODE
    # ----------------------
    s_action=[]
    Vs = [np.zeros(mdp.nS)]
    V_prev=Vs[-1]

    for state in range(mdp.nS):

        q=[]
        for next_state in range(len(mdp.P[state][pi[state]])):
            q.append((mdp.P[state][pi[state]][next_state][0]*(mdp.P[state][

        s_action.append(sum(q))
        V[state]=s_action[state]
        Vs.append(V)
    #first iteration(To make V_prev!=V)

    while any(V_prev!=V):
        V=np.zeros(mdp.nS)
        V_prev=Vs[-1]
        s_action=[]
        for state in range(mdp.nS):
            q=[]
            for next_state in range(len(mdp.P[state][pi[state]])):
                q.append((mdp.P[state][pi[state]][next_state][0]*(mdp.P[sta
            s_action.append(sum(q))
            V[state]=s_action[state]

        Vs.append(V)

    # ----------------------
    # END YOUR CODE

    return V
```

Now let's compute the value of an arbitrarily-chosen policy.

In [7]:

```
begin_grading()
print(compute_vpi(np.ones(16), mdp, gamma=GAMMA))
end_grading()
```

```
[0.016 0.024 0.232 0.024 0.017 0.    0.299 0.    0.02  0.188 0.393 0.
 0.    0.196 0.494 0.   ]
```

As a sanity check, if we run `compute_vpi` on the solution from our previous value iteration run, we should get approximately (but not exactly) the same values produced by value iteration.

In [8]:
```
Vpi=compute_vpi(pis_VI[15], mdp, gamma=GAMMA)
V_vi = Vs_VI[15]
print("From compute_vpi", Vpi)
print("From value iteration", V_vi)
print("Difference", Vpi - V_vi)
```

```
From compute_vpi [0.531 0.471 0.56  0.471 0.574 0.    0.62  0.    0.683 0.8
27 0.815 0.
 0.    0.901 0.97  0.   ]
From value iteration [0.53  0.47  0.56  0.47  0.573 0.    0.62  0.    0.683
0.827 0.815 0.
 0.    0.901 0.97  0.   ]
Difference [9.580e-04 3.839e-04 2.254e-04 3.839e-04 4.495e-04 0.000e+00 4.5
22e-05
 0.000e+00 2.612e-04 1.071e-04 3.272e-05 0.000e+00 0.000e+00 3.977e-05
 7.051e-06 0.000e+00]
```

Question 3b: state-action value function

Next, you'll write a function to compute the state-action value function Q^π, defined as follows

$$Q^\pi(s,a) = \sum_{s'} P(s,a,s')[R(s,a,s') + \gamma V^\pi(s')]$$

In [9]:
```
def compute_qpi(vpi, pi, mdp,  gamma):
    Qpi = np.zeros([mdp.nS,mdp.nA])

    # BEGIN YOUR CODE
    # ------------------------
    for state in range(mdp.nS):
        s_action=[]
        for a in range(mdp.nA):
            q=[]
            for next_state in range(len(mdp.P[state][a])):
                q.append((mdp.P[state][a][next_state][0]*(mdp.P[state][a][n
            s_action.append(sum(q))
            Qpi[state,a]=s_action[-1]

    # ------------------------
    # END YOUR CODE

    return Qpi

begin_grading()
Qpi = compute_qpi(Vpi, pis_VI[-1], mdp, gamma=0.95)
end_grading()
print("Qpi:\n", Qpi)
```

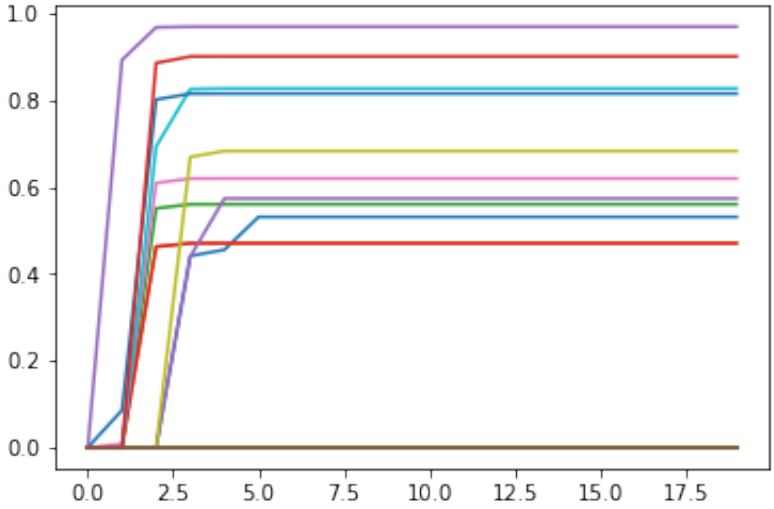
```
Qpi:
 [[0.509 0.531 0.463 0.499]
 [0.448 0.104 0.471 0.461]
 [0.47  0.56  0.47  0.515]
 [0.471 0.098 0.402 0.456]
 [0.551 0.574 0.115 0.458]
 [0.    0.    0.    0.   ]
 [0.131 0.62  0.131 0.426]
 [0.    0.    0.    0.   ]
 [0.574 0.143 0.683 0.579]
 [0.605 0.827 0.705 0.142]
 [0.78  0.815 0.151 0.55 ]
 [0.    0.    0.    0.   ]
 [0.    0.    0.    0.   ]
 [0.164 0.777 0.901 0.721]
 [0.854 0.922 0.97  0.805]
 [0.    0.    0.    0.   ]]
```

Now we're ready to run policy iteration!

In [10]:

```
def policy_iteration(mdp, gamma, nIt):
    Vs = []
    pis = []
    pi_prev = np.zeros(mdp.nS,dtype='int')
    pis.append(pi_prev)
    print("Iteration | # chg actions | V[0]")
    print("----------+---------------+---------")
    for it in range(nIt):
        vpi = compute_vpi(pi_prev, mdp, gamma)
        qpi = compute_qpi(vpi, pi_prev, mdp, gamma)
        pi = qpi.argmax(axis=1)
        print("%4i      | %6i        | %6.5f"%(it, (pi != pi_prev).sum(), v
        Vs.append(vpi)
        pis.append(pi)
        pi_prev = pi
    return Vs, pis
Vs_PI, pis_PI = policy_iteration(mdp, gamma=0.95, nIt=20)
plt.plot(Vs_PI);
```

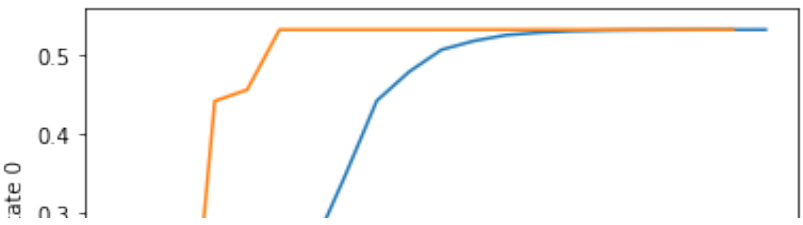
```
Iteration | # chg actions | V[0]
----------+---------------+---------
    0     |       1       | 0.00000
    1     |       6       | 0.00000
    2     |       3       | 0.00000
    3     |       1       | 0.44131
    4     |       1       | 0.45546
    5     |       0       | 0.53118
    6     |       0       | 0.53118
    7     |       0       | 0.53118
    8     |       0       | 0.53118
    9     |       0       | 0.53118
   10     |       0       | 0.53118
   11     |       0       | 0.53118
   12     |       0       | 0.53118
   13     |       0       | 0.53118
   14     |       0       | 0.53118
   15     |       0       | 0.53118
   16     |       0       | 0.53118
   17     |       0       | 0.53118
   18     |       0       | 0.53118
   19     |       0       | 0.53118
```

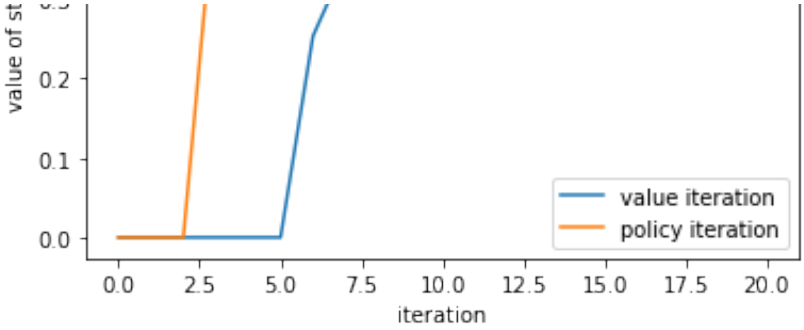


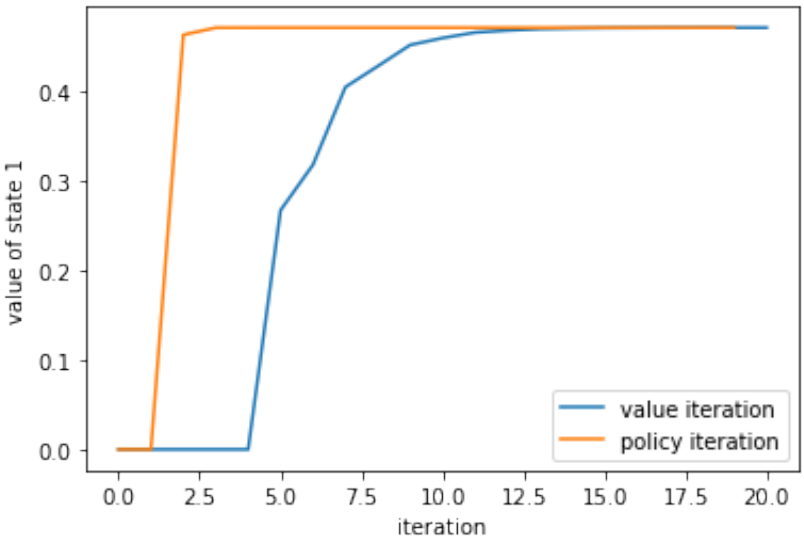
Now we can compare the convergence of value iteration and policy iteration on several states. For fun, you can try adding modified policy iteration.

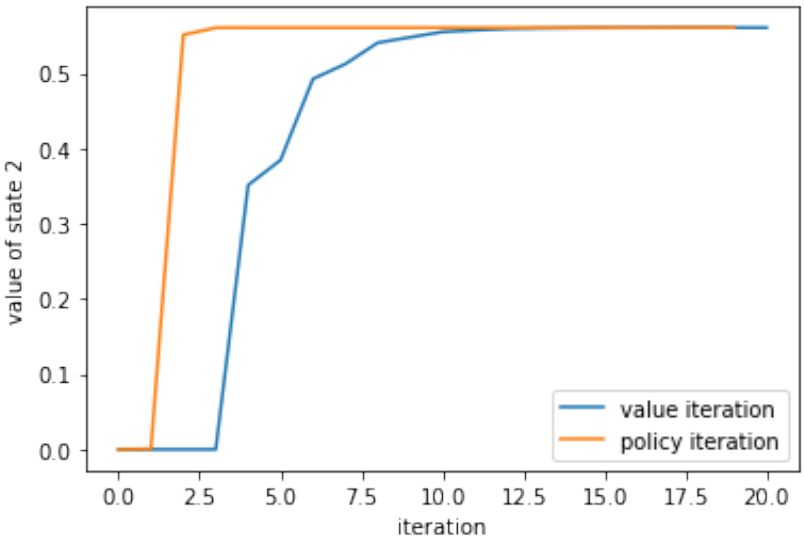
In [11]:

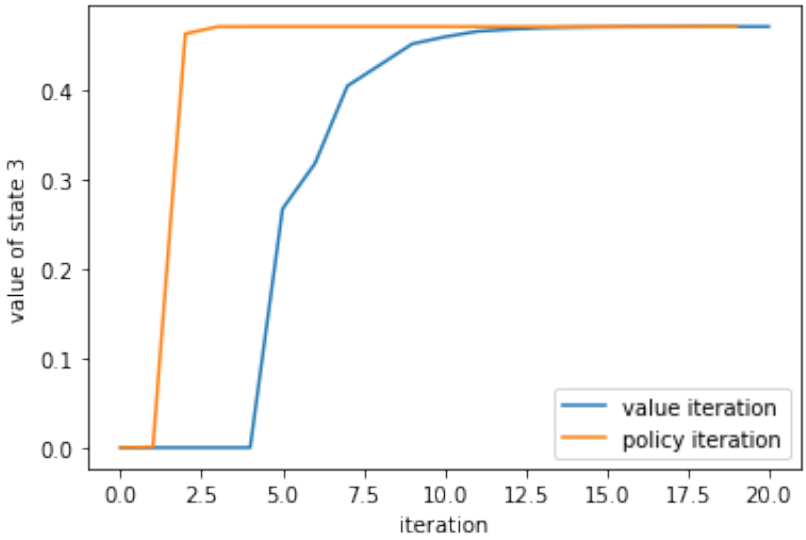
```
for s in range(5):
    plt.figure()
    plt.plot(np.array(Vs_VI)[:,s])
    plt.plot(np.array(Vs_PI)[:,s])
    plt.ylabel("value of state %i"%s)
    plt.xlabel("iteration")
    plt.legend(["value iteration", "policy iteration"], loc='best')
```

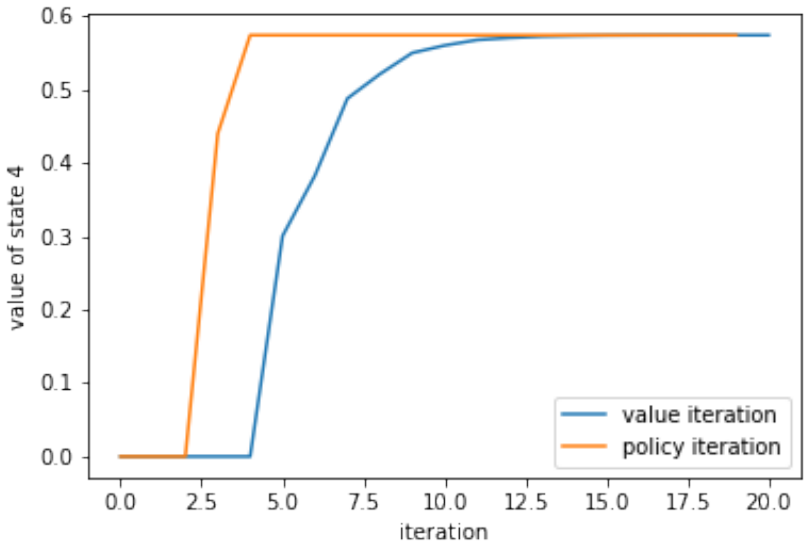



value of st
0.2
0.1
0.0
value iteration
policy iteration
0.0 2.5 5.0 7.5 10.0 12.5 15.0 17.5 20.0
iteration


value of state 1
0.4
0.3
0.2
0.1
0.0
value iteration
policy iteration
0.0 2.5 5.0 7.5 10.0 12.5 15.0 17.5 20.0
iteration


value of state 2
0.5
0.4
0.3
0.2
0.1
0.0
value iteration
policy iteration
0.0 2.5 5.0 7.5 10.0 12.5 15.0 17.5 20.0
iteration





Congratulations! You have completed HW1.

Make sure you have run all cells in your notebook in order before running the cell below, so that all images/graphs appear in the output.

Please generate pdf as follows and submit it to Gradescope.

File > Print Preview > Print > Save as pdf

Please save before submitting!

In [ ]: