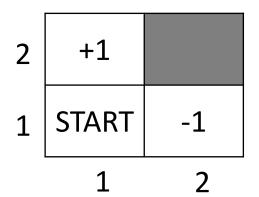
Markov Decision Processes Part 3: Computing State Utilities and Optimal Policies

CSE 4309 – Machine Learning
Vassilis Athitsos
Computer Science and Engineering Department
University of Texas at Arlington

Review: the Bellman Equation



$$U(s) = R(s) + \gamma \max_{a \in A(s)} \left\{ \sum_{s'} [p(s'|s,a)U(s')] \right\}$$

- For each state s, we get a Bellman equation.
- If our environment has *N* states, we need to solve a system of *N* Bellman equations.
- In this system of equations, there is a total of Nunknowns:
 - The N values U(s).
- There is an iterative algorithm for solving this system of equations, called the value iteration algorithm.

- The value iteration algorithm computes the utility of each state for a Markov Decision Process.
- The algorithm takes the following inputs:
 - The set of states $\mathbb{S} = \{s_1, \dots, s_N\}$.
 - The set A(s) of actions available at each state s.
 - The transition model p(s' | s, a).
 - The reward function R(s)
 - The discount factor γ .
 - $-\varepsilon$, which is the maximum error allowed in the utility of each state, in the result of the algorithm.

function ValueIteration(\mathbb{S} , A, p, R, γ , ε)

N = size of S.

U' = new array of doubles, of size N.

Initialize all values of U' to 0.

repeat:

U = copy of array U'

$$\delta = 0$$

for each state s in S:

$$U'[s] = R(s) + \gamma \max_{a \in A(s)} \{ \sum_{s'} [p(s'|s, a)U[s']] \}$$
if $|U'[s] - U[s]| > \delta$ then $\delta = |U'[s] - U[s]|$

if
$$|U'[s] - U[s]| > \delta$$
 then $\delta = |U'[s] - U[s]$

until $\delta < \varepsilon (1 - \gamma)/\gamma$

```
function ValueIteration(\mathbb{S}, A, p, R, \gamma, \varepsilon)
     N = size of S.
     U' = new array of doubles, of size N.
     Initialize all values of U' to 0.
     repeat:
          U = copy of array U'
          \delta = 0
          for each state s in S:
               U'[s] = R(s) + \gamma \max_{a \in A(s)} \{ \sum_{s'} [p(s'|s, a)U[s']] \}
if |U'[s] - U[s]| > \delta then \delta = |U'[s] - U[s]|
     until \delta < \varepsilon (1 - \gamma)/\gamma
     return U
```

- We will skip the proof, but it can be proven that this algorithm converges to the correct solutions of the Bellman equations.
 - Details can be found in S. Russell and P. Norvig, "Artificial Intelligence:
 A Modern Approach", third edition (2009), Prentice Hall.

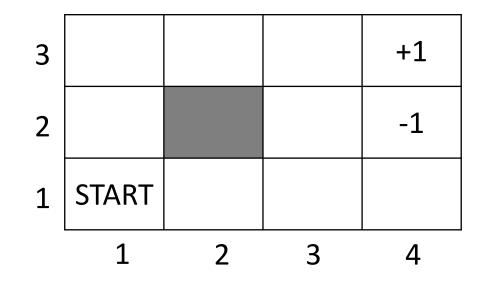
```
function ValueIteration(\mathbb{S}, A, p, R, \gamma, \varepsilon)
     N = \text{size of } S.
     U' = new array of doubles, of size N.
     Initialize all values of U' to 0.
     repeat:
          U = copy of array U'
          \delta = 0
          for each state s in S:
               U'[s] = R(s) + \gamma \max_{a \in A(s)} \{ \sum_{s'} [p(s'|s, a)U[s']] \}
if |U'[s] - U[s]| > \delta then \delta = |U'[s] - U[s]|
     until \delta < \varepsilon (1 - \gamma)/\gamma
     return U
```

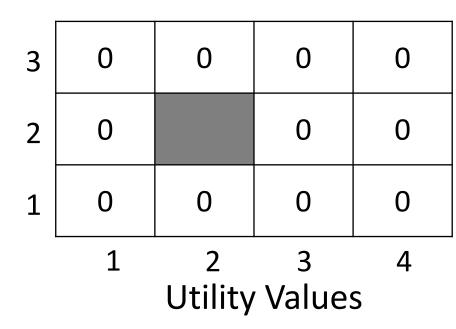
- The main operation of this algorithm is highlighted in red.
- We use the Bellman equation to update values U(s) using the previous estimates for those values.
 - This update step is called a Bellman update.

```
function ValueIteration(\mathbb{S}, A, p, R, \gamma, \varepsilon)
     N = \text{size of } S.
     U' = new array of doubles, of size N.
     Initialize all values of U' to 0.
     repeat:
          U = copy of array U'
          \delta = 0
          for each state s in S:
              U'[s] = R(s) + \gamma \max_{a \in A(s)} \{ \sum_{s'} [p(s'|s, a)U[s']] \}
if |U'[s] - U[s]| > \delta then \delta = |U'[s] - U[s]|
     until \delta < \varepsilon (1 - \gamma)/\gamma
     return U
```

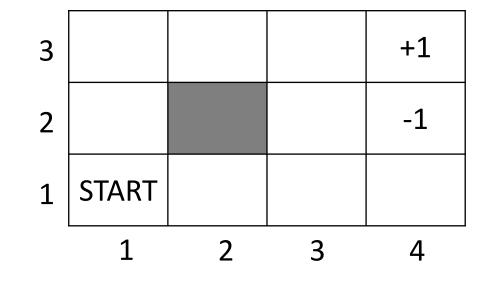
- So, the value iteration algorithm can be summarized as follows:
 - Initialize utilities of states to zero values.
 - Repeat updating utilities of states using Bellman updates, until the estimated values converge.

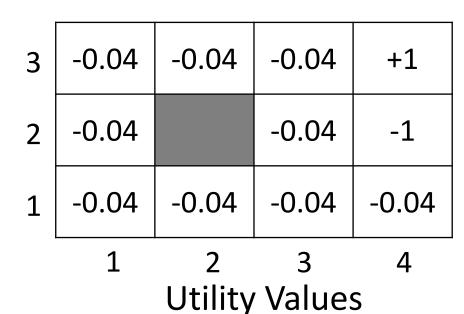
- Let's see how the value iteration algorithm works on our example.
- Assume:
 - -R(s) = -0.04 if s is a non-terminal state.
 - $-\gamma = 0.9$
- We initialize all utility values to 0.





- Let's see how the value iteration algorithm works on our example.
- Assume:
 - -R(s) = -0.04 if s is a non-terminal state.
 - $-\gamma = 0.9$
- This is the result after one round of updates:
 - The current estimate for each state s is R(s).





 Let's see how the value iteration algorithm works on our example.

• Assume:

- -R(s) = -0.04 if s is a non-terminal state.
- $\gamma = 0.9$
- This is the result after two rounds of updates:
 - Information about the +1 reward reached state (3,3).

3				+1
2				-1
1	START			
	1	2	3	4

3	-0.08	-0.08	0.67	+1
2	-0.08		-0.08	-1
1	-0.08	-0.08	-0.08	-0.08
	1	2	3	4
Utility Values				

 Let's see how the value iteration algorithm works on our example.

Assume:

- -R(s) = -0.04 if s is a non-terminal state.
- $-\gamma = 0.9$
- This is the result after three rounds of updates:
 - Information about the +1 reward reached more states.

3				+1
2				-1
1	START			
	1	2	3	4

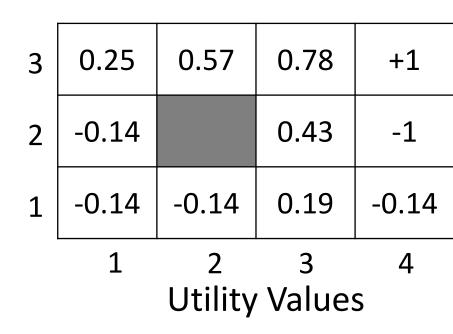
3	-0.11	0.43	0.73	+1	
2	-0.11		0.35	-1	
1	-0.11	-0.11	-0.11	-0.11	
	1	2	3	4	
	Utility Values				

 Let's see how the value iteration algorithm works on our example.

Assume:

- -R(s) = -0.04 if s is a non-terminal state.
- $\gamma = 0.9$
- This is the result after four rounds of updates:
 - Information about the +1 reward reached more states.

3				+1
2				-1
1	START			
,	1	2	3	4



 Let's see how the value iteration algorithm works on our example.

• Assume:

- -R(s) = -0.04 if s is a non-terminal state.
- $-\gamma = 0.9$
- This is the result after five rounds of updates:
 - Information about the +1 reward reached more states.

3				+1
2				-1
1	START			
	1	2	3	4

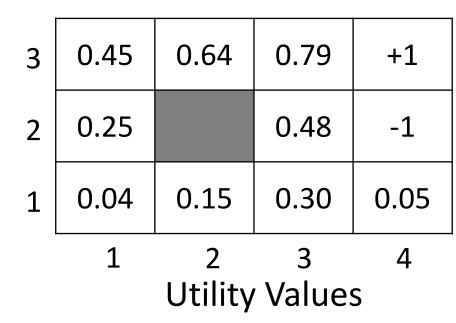
3	0.38	0.62	0.79	+1
2	0.12		0.47	-1
1	-0.16	0.07	0.24	-0.01
	1	2	3	4
Utility Values				

 Let's see how the value iteration algorithm works on our example.

Assume:

- -R(s) = -0.04 if s is a non-terminal state.
- $-\gamma = 0.9$
- This is the result after six rounds of updates:
 - Information about the +1 reward has reached all states.

3				+1
2				-1
1	START			
·	1	2	3	4



 Let's see how the value iteration algorithm works on our example.

Assume:

- -R(s) = -0.04 if s is a non-terminal state.
- $-\gamma = 0.9$
- This is the result after seven rounds of updates:
 - Values keep getting updated.

3				+1
2				-1
1	START			
'	1	2	3	4

3	0.48	0.65	0.79	+1
2	0.33		0.48	-1
1	0.16	0.21	0.32	0.09
•	1		2	4

Utility Values

 Let's see how the value iteration algorithm works on our example.

Assume:

- -R(s) = -0.04 if s is a non-terminal state.
- $\gamma = 0.9$
- This is the result after eight rounds of updates:
 - Values continue changing.

3				+1
2				-1
1	START			
·	1	2	3	4

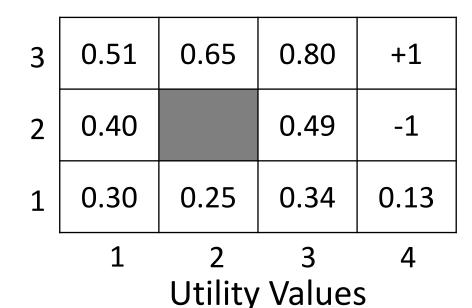
3	0.50	0.65	0.80	+1	
2	0.37		0.49	-1	
1	0.23	0.23	0.34	0.11	
,	1	2	3	4	
	Utility Values				

 Let's see how the value iteration algorithm works on our example.

Assume:

- -R(s) = -0.04 if s is a non-terminal state.
- $-\gamma = 0.9$
- This is the result after 13 rounds of updates:
 - Values don't change much anymore after this round.

3				+1
2				-1
1	START			
	1	2	3	4

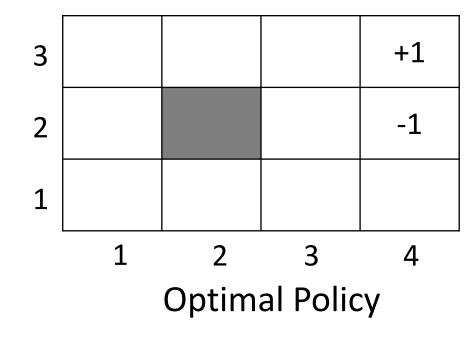


- The value iteration algorithm computes U(s) for every state s.
- Once we have computed all values U(s), we can get the optimal policy π^* using this equation:

$$\pi^*(s) = \underset{a \in A(s)}{\operatorname{argmax}} \left\{ \sum_{s'} [p(s'|s,a)U(s')] \right\}$$

- Thus, $\pi^*(s)$ identifies the action that leads to the highest expected utility for the next state, as measured over all possible outcomes of that action.
- This approach is called **one-step look-ahead.**

- At the bottom we see the result of the value iteration algorithm for:
 - -R(s) = -0.02 if s is a non-terminal state.
 - $-\gamma = 1$
- How can we figure out the optimal policy based on that output?

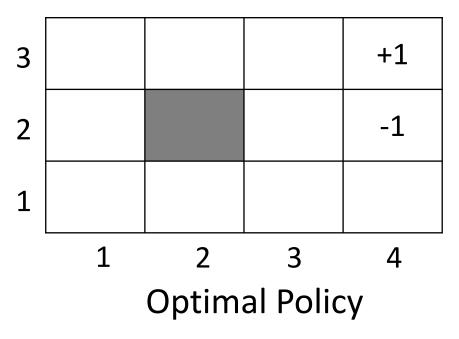


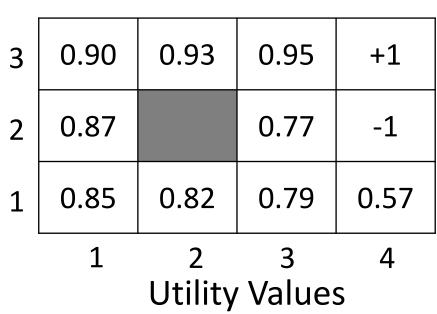
0.90	0.93	0.95	+1
0.87		0.77	-1
0.85	0.82	0.79	0.57
1	2	3	4
Utility Values			
	0.87	0.87 0.85 0.82	0.87 0.77 0.85 0.82 0.79 1 2 3

- Consider state (2,3).
- What is the optimal action for that state?
- We must consider each action.
- If the action is "left", these are the possible next states:

Probability	Next State	Utility
0.8	(2,3)	0.77
0.1	(3,3)	0.95
0.1	(1,3)	0.79

The weighted average is 0.79

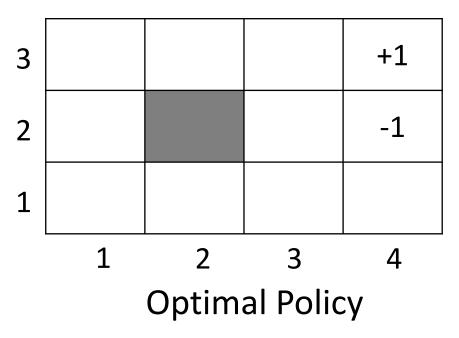


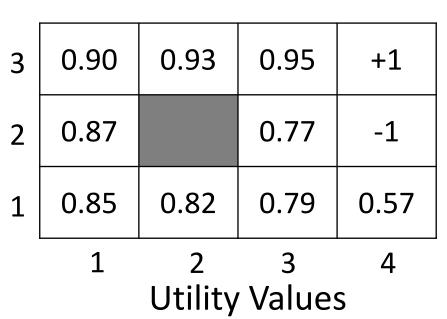


- Consider state (2,3).
- What is the optimal action for that state?
- We must consider each action.
- If the action is "right", these are the possible next states:

Probability	Next State	Utility
0.8	(2,3)	-1
0.1	(3,3)	0.95
0.1	(1,3)	0.79

• The weighted average is -0.63

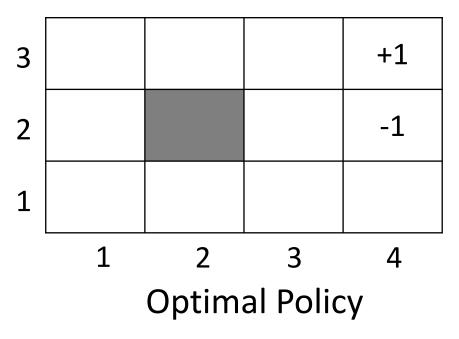


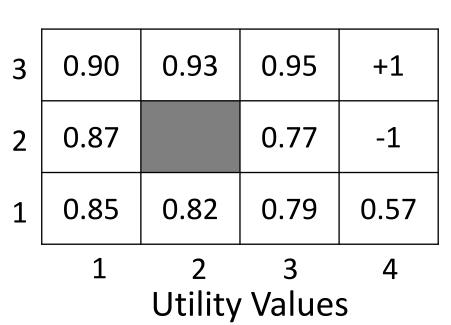


- Consider state (2,3).
- What is the optimal action for that state?
- We must consider each action.
- If the action is "up", these are the possible next states:

Probability	Next State	Utility
0.8	(3,3)	0.95
0.1	(2,4)	-1.00
0.1	(2,3)	0.77

The weighted average is 0.74

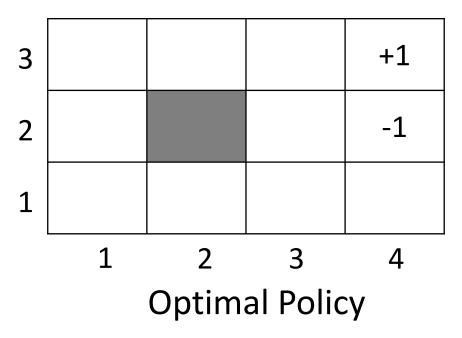


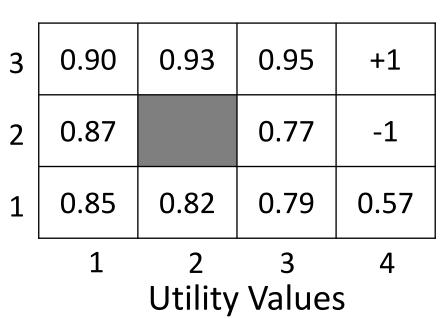


- Consider state (2,3).
- What is the optimal action for that state?
- We must consider each action.
- If the action is "down", these are the possible next states:

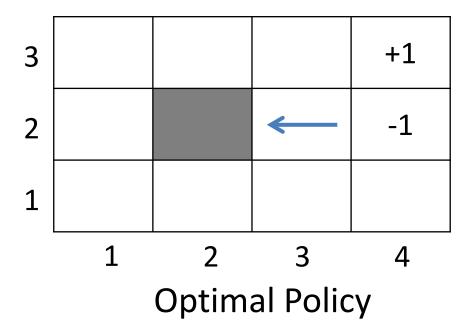
Probability	Next State	Utility
0.8	(1,3)	0.79
0.1	(2,4)	-1.00
0.1	(2,3)	0.77

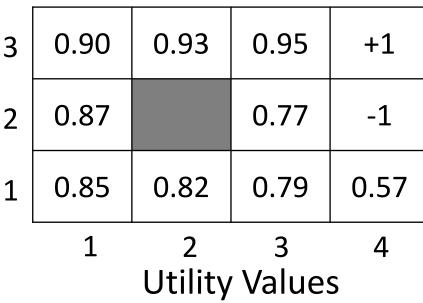
The weighted average is 0.61



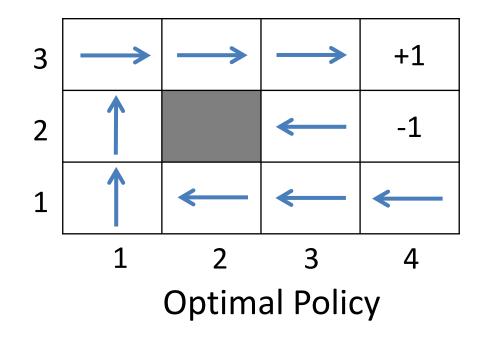


- For state (2,3), action "left" led to the highest expected utility for the next state.
- Thus, action "left" is the best action for state (2,3).
- Note that choosing the best action is not always to try to move towards the best state.
 - At state (2,3) the best action is towards the blocked square, to play it safe.
 - Going up is risky, it has a 10% chance to lead to the -1 state.





- Here is the optimal policy for:
 - -R(s) = -0.02 if s is a non-terminal state.
 - $-\gamma = 1$
- Note that choosing the best policy is more complicated than simply pointing to the direction of highest reward.
 - At state (2,3) the best action is towards the blocked square, to play it safe.
 - Going up is risky, it has a 10% chance to lead to the -1 state.



0.90	0.93	0.95	+1
0.87		0.77	-1
0.85	0.82	0.79	0.57
1	2 Utility	3 Value	4
		0.85 0.82 1 2	0.85 0.82 0.79

- There is an alternative algorithm for computing optimal policies, that is more efficient.
- Remember that, if we know the utility of each state, we can compute the optimal policy π^* using:

$$\pi^*(s) = \underset{a \in A(s)}{\operatorname{argmax}} \left\{ \sum_{s'} [p(s'|s,a)U(s')] \right\}$$

- However, to get the right $\pi^*(s)$, we don't need to know the utilities very accurately.
 - We just need to know the utilities accurately enough, so that, for each state s, argmax chooses the right action.

- This alternative algorithm for computing optimal policies is called the **policy iteration algorithm**.
- It is an iterative algorithm.
- Initialization:
 - Initiate some policy π_0 with random choices for the best action at each state.
- Main loop:
 - <u>Policy evaluation</u>: given the current policy π_i , calculate utility values $U^{\pi_i}(s)$, corresponding to the utility of each state s **if the agent follows policy** π_i .
 - <u>Policy improvement</u>: Given current utility values $U^{\pi_i}(s)$, use one-step look-ahead to compute new policy π_{i+1} .

- To be able to implement the policy iteration algorithm, we need to specify how to carry out each of the two steps of the main loop:
 - Policy evaluation.
 - Policy improvement.

- Task: calculate utility values $U^{\pi_i}(s)$, corresponding to the assumption that **the agent follows policy** π_i .
- When the policy was not known, we used the Bellman equation:

$$U(s) = R(s) + \gamma \max_{a \in A(s)} \left\{ \sum_{s'} [p(s'|s,a)U(s')] \right\}$$

• Now that the policy π_i is specified, we can instead use a simplified version of the **Bellman equation**:

$$U^{\pi_i}(s) = R(s) + \gamma \sum_{s'} [p(s'|s, \pi_i(s))U^{\pi_i}(s')]$$

• Key difference: now $\pi_i(s)$ specifies the action for each state s, so we do not need to look for the max over all possible actions.

$$U^{\pi_i}(s) = R(s) + \gamma \sum_{s'} [p(s'|s, \pi_i(s))U^{\pi_i}(s')]$$

- This is a linear equation.
 - The original Bellman equation, taking the max out of all possible actions, is **not** a linear equation.
- If we have *N* states, we get *N* linear equations of this form, with *N* unknowns.
- We can solve those N linear equations in $O(N^3)$ time, using standard linear algebra methods.

- For large state spaces, $O(N^3)$ is prohibitive.
- Alternative: do some rounds of iterations.

```
function PolicyEvaluation(\mathbb{S}, \mathbb{p}, \mathbb{R}, \gamma, \pi_i, K, \mathbb{U})
 \mathbb{U}_0 = \text{copy of } \mathbb{U} 
for k = 1 to K:
  for each state \mathbb{S} in \mathbb{S}:
 \mathbb{U}_k(s) = \mathbb{R}(s) + \gamma \sum_{s'} [p(s'|s,\pi_i(s)) \mathbb{U}_{k-1}(s')] 
return \mathbb{U}_k
```

- Obviously, doing K iterations does not guarantee that the utilities are computed correctly.
- Parameter K allows us to trade speed for accuracy. Larger values lead to slower runtimes and higher accuracy.

- For large state spaces, $O(N^3)$ is prohibitive.
- Alternative: do some rounds of iterations.

```
function PolicyEvaluation(\mathbb{S}, p, R, \gamma, \pi_i, K, U)
U_0 = \text{copy of } U
\text{for } k = 1 \text{ to } K:
\text{for each state } s \text{ in } \mathbb{S}:
U_k(s) = R(s) + \gamma \sum_{s'} [p(s'|s, \pi_i(s)) U_{k-1}(s')]
\text{return } U_k
```

- The PolicyEvaluation function takes as argument a current estimate U.
 - See later how the PolicyEvaluation function is called from the PolicyIteration function.

function PolicyIteration(\mathbb{S} , A, p, R, γ , K)

```
N= size of \mathbb{S}. U = new array of size N, all values initialized to 0 \pi= new array of actions, of size N Initialize all values of \pi to random (but legal) actions repeat:
```

 $U = PolicyEvaluation(S, p, R, \gamma, \pi, K, U)$ unchanged = **true for each** state s in S:

$$\begin{aligned} &\inf\max_{a\in A(s)} \left\{ \sum_{s'} \left[p(s'|s,a) \mathsf{U}[s'] \right] \right\} > \sum_{s'} \left[p(s'|s,\pi[s]) \mathsf{U}[s'] \right] \\ &\pi[s] = \underset{a\in A(s)}{\operatorname{argmax}} \left\{ \sum_{s'} \left[p(s'|s,a) \mathsf{U}[s'] \right] \right\} \\ &\operatorname{unchanged} = \mathbf{false} \end{aligned}$$

until unchanged == true

return π

function PolicyIteration(\mathbb{S} , A, p, R, γ , K)

N= size of \mathbb{S} . U = new array of size N, all values initialized to 0 $\pi=$ new array of actions, of size N Initialize all values of π to random (but legal) actions repeat:

 $U = PolicyEvaluation(S, p, R, \gamma, \pi, K, U)$ unchanged = **true for each** state s in S: The main loop alternates between:

- Updating the utilities given the policy.
- Updating the policy given the utilities.

The main loop exits when the policy stops changing.

$$\begin{aligned} &\inf\max_{a\in A(s)} \left\{ \sum_{s'} \left[p(s'|s,a) \mathsf{U}[s'] \right] \right\} > \sum_{s'} \left[p(s'|s,\pi[s]) \mathsf{U}[s'] \right] \\ &\pi[s] = \underset{a\in A(s)}{\operatorname{argmax}} \left\{ \sum_{s'} \left[p(s'|s,a) \mathsf{U}[s'] \right] \right\} \\ &\operatorname{unchanged} = \mathbf{false} \end{aligned}$$

until unchanged == true

return π

function PolicyIteration(\mathbb{S} , A, p, R, γ , K)

```
N= size of \mathbb{S}. U= new array of size N, all values initialized to 0 \pi= new array of actions, of size N Initialize all values of \pi to random (but legal) actions repeat:
```

```
U = PolicyEvaluation(S, p, R, \gamma, \pi, K, U)
unchanged = true
for each state s in S:
```

The main loop alternates between:

- Updating the utilities given the policy.
- Updating the policy given the utilities.

The main loop exits when the policy stops changing.

```
\begin{aligned} &\inf\max_{a\in A(s)} \left\{ \sum_{s'} \left[ p(s'|s,a) \mathsf{U}[s'] \right] \right\} > \sum_{s'} \left[ p(s'|s,\pi[s]) \mathsf{U}[s'] \right] \\ &\pi[s] = \underset{a\in A(s)}{\operatorname{argmax}} \left\{ \sum_{s'} \left[ p(s'|s,a) \mathsf{U}[s'] \right] \right\} \\ &\operatorname{unchanged} = \mathbf{false} \end{aligned}
```

until unchanged == true

function PolicyIteration(\mathbb{S} , A, p, R, γ , K)

N= size of \mathbb{S} . U= new array of size N, all values initialized to 0 $\pi=$ new array of actions, of size N Initialize all values of π to random (but legal) actions **repeat:**

 $U = PolicyEvaluation(S, p, R, \gamma, \pi, K, U)$ unchanged = **true**

for each state s in S:

 $\inf_{a \in A(s)} \{ \sum_{s'} [p(s'|s,a) \mathbf{U}[s']] \} > \sum_{s'} [p(s'|s,\pi[s]) \mathbf{U}[s']]$ $\pi[s] = \underset{a \in A(s)}{\operatorname{argmax}} \{ \sum_{s'} [p(s'|s,a) \mathbf{U}[s']] \}$

unchanged = false

until unchanged == true

The main loop alternates between:

- Updating the utilities given the policy.
- Updating the policy given the utilities.

The main loop exits when the policy stops changing.

Markov Decision Processes: Recap

In Markov Decision Processes:

- Each state has a reward R(s).
- Each state sequence $(s_0, s_1, ..., s_T)$ has a utility U_h which is computed by adding the discounted rewards of all states in the sequence.
- An action can lead to multiple outcomes. The probability of each outcome given the state and the action is known.
- A policy is a function mapping states to actions.
- The utility of a state s_0 is the expected utility measured over all state sequences that can lead from s_0 to a terminal state, under the assumption that the agent follows the optimal policy.

Markov Decision Processes: Recap

- The value iteration algorithm computes the utility of each state using an iterative approach.
 - Once the utilities of all states have been computed, the optimal policy is defined by identifying, for each state, the action leading to the highest expected utility.
- The policy iteration algorithm is a more efficient alternative, at the cost of possibly losing some accuracy.
 - It computes the optimal policy directly, without computing exact values for the utilities.
 - Utility values are updated for a few rounds only, and not until convergence.