

ECE 1508: Reinforcement Learning

Chapter 2: Model-based RL

Ali Bereyhi

`ali.bereyhi@utoronto.ca`

Department of Electrical and Computer Engineering
University of Toronto

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Last Piece: *Dynamic Programming*

Right now, we know what to do when we know *MDP of environment*

- ① We can find optimal values from *Bellman optimality equations*
- ② We could then find the *optimal action*-values
- ③ We finally get the *optimal policy* from *optimal action*-values

The only remaining challenge is to find

an algorithmic approach to solve *Bellman optimality equations*

We complete this last piece using

Dynamic Programming \equiv *DP*

Dynamic Programming: *Basic Idea*

Assume, we want to solve the problem of

$$x = f(x)$$

for some function $f(x)$

We could solve it via *direct approach*:

- ① Rewrite is as $f(x) - x = 0$
- ② Solve it via classic algorithms
 - ↳ Reduce it to a *known form*, e.g., a *polynomial*
 - ↳ Solve it via an *iterative method*, e.g., *Newton-Raphson* or *method of intervals*

Dynamic Programming: *Basic Idea*

Assume, we want to solve the problem of

$$x = f(x)$$

for some function $f(x)$

We could also solve it by *recursion*:

- ① Start with an x^0 and set $x^1 = f(x^0)$
- ② Until $x^{k+1} \approx x^k$, we do
 - ↳ Update *recursively* as $x^{k+1} = f(x^k)$
 - ↳ Set $k \leftarrow k + 1$

Under *some conditions on $f(\cdot)$* , this approach can *converge*

Dynamic Programming: Example

We want to solve

$$x = \frac{-1}{2 + x}$$

- ① Start with an $x^0 = 0$
- ② We now get into the recursion loop

$$\hookrightarrow x^1 = f(x^0) = -\frac{1}{2}$$

$$\hookrightarrow x^2 = f(x^1) = -\frac{2}{3}$$

$$\hookrightarrow x^3 = f(x^2) = -\frac{3}{4}$$

$$\hookrightarrow \dots$$

$$\hookrightarrow x^k = f(x^{k-1}) = -\frac{k}{k+1}$$

We asymptotically **converge** to $x^\infty = -1$ which is the **solution**

\hookrightarrow Note that we **always** converge **no matter** which point we start

Dynamic Programming: Example

Now, let's write the *same equation* in a *different recursive form*

$$x = \frac{-1 - x^2}{2}$$

① Start with an $x^0 = 0$

② We get into recursion loop

$$\hookrightarrow x^1 = f(x^0) = -0.5$$

$$\hookrightarrow x^2 = f(x^1) = -0.625$$

$$\hookrightarrow \dots$$

$$\hookrightarrow x^\infty = -1$$

① Start with an $x^0 = 5$

② We get into recursion loop

$$\hookrightarrow x^1 = f(x^0) = -13$$

$$\hookrightarrow x^2 = f(x^1) = -85$$

$$\hookrightarrow \dots$$

$$\hookrightarrow x^\infty = -\infty$$

We can now diverge if we start with a wrong initial point!

Not all recursive forms are always converging!

Dynamic Programming: *Applications to Our Problem*

Our problem has a similar form: we need to solve *Bellman equations*
which are *recursive equations*

So, we could use **DP** to find the solution

There are two major **DP** approaches

- Policy Iteration that uses recursion to iterate between
 - ↳ Policy *Evaluation*
 - ↳ Policy *Improvement*
- Value Iteration which applies recursion on *optimal Bellman* equation

Let's look at these two approaches in detail

Policy Evaluation: Step 1

The first step is *policy evaluation*: we can formulate this problem as follows

Ultimate Goal of Policy Evaluation

Given a *policy* π , we intend to *evaluate* values of *all states* by *recursion*

Before we start, let's recap a few definitions: recall *expected policy reward*

$$\bar{\mathcal{R}}_{\pi}(s) = \sum_{m=1}^M \bar{\mathcal{R}}(a^m, s) \pi(a^m | s)$$

For sake of compactness, we use the following notation

$$\bar{\mathcal{R}}_{\pi}(s) = \mathbb{E}_{\pi} \{ \bar{\mathcal{R}}(A, s) | s \}$$

Policy Evaluation: Step I

Similarly, we define the notation

$$\mathbb{E}_{\pi} \{ v_{\pi} (\bar{S}) | s, a \} = \sum_{n=1}^N v_{\pi} (s^n) p (s^n | s, a)$$

and also denote its expected form over the **action set** by

$$\begin{aligned} \mathbb{E}_{\pi} \{ v_{\pi} (\bar{S}) | s \} &= \sum_{n=1}^N v_{\pi} (s^n) p_{\pi} (s^n | s) \\ &= \sum_{m=1}^M \underbrace{\sum_{n=1}^N v_{\pi} (s^n) p (s^n | s, a^m)}_{\mathbb{E}_{\pi} \{ v_{\pi} (\bar{S}) | s, a^m \}} \pi (a^m | s) \\ &= \sum_{m=1}^M \mathbb{E}_{\pi} \{ v_{\pi} (\bar{S}) | s, a^m \} \pi (a^m | s) \end{aligned}$$

Policy Evaluation: Step 1

We can then write the *Bellman equations* compactly as

$$v_{\pi}(s) = \bar{\mathcal{R}}_{\pi}(s) + \gamma \mathbb{E}_{\pi} \{ v_{\pi}(\bar{S}) | s \}$$

for *value function* and also as

$$q_{\pi}(s, a) = \bar{\mathcal{R}}(s, a) + \gamma \mathbb{E}_{\pi} \{ v_{\pi}(\bar{S}) | s, a \}$$

for *action-value function*

Now, we are ready to *evaluate a policy* by *recursion*

Policy Evaluation: *Value Computation via Recursion*

Recall our perspective on value computation:

values are N unknowns that we want to compute from Bellman equations

Now, if someone claims that *the values*

$$v_{\pi}(s^n) = v_n$$

for $n = 1 : N$ are values of policy π , can we confirm it?

- + Shouldn't we simply use *Bellman Equation*?!
 - Exactly!

Policy Evaluation: Value Computation via Recursion

We could confirm

$$v_{\pi}(s^n) = v_n$$

by writing first finding for every state s

$$\begin{aligned}
 \mathbb{E}_{\pi} \{v_{\pi}(\bar{S}) | s\} &= \sum_{n=1}^N v_{\pi}(s^n) p_{\pi}(s^n | s) \\
 &= \sum_{n=1}^N \sum_{m=1}^M v_{\pi}(s^n) p(s^n | s, a^m) \pi(a^m | s) \\
 &= \sum_{n=1}^N \sum_{m=1}^M \underbrace{v_n}_{\text{claimed value}} \underbrace{p(s^n | s, a^m)}_{\text{transition model}} \underbrace{\pi(a^m | s)}_{\text{policy}}
 \end{aligned}$$

Policy Evaluation: Value Computation via Recursion

We could confirm

$$v_{\pi}(s^n) = v_n$$

by writing first finding for every state s

$$\mathbb{E}_{\pi} \{v_{\pi}(\bar{S}) | s\} = \text{computed from } v_n\text{'s} := F(\{v_1, \dots, v_N\}, s)$$

and then checking if

$$\begin{aligned} v_{\pi}(s^n) &= v_n = \bar{\mathcal{R}}_{\pi}(s^n) + \gamma \mathbb{E}_{\pi} \{v_{\pi}(\bar{S}) | s^n\} \\ &= \bar{\mathcal{R}}_{\pi}(s^n) + \gamma F(\{v_1, \dots, v_N\}, s) \end{aligned}$$

holds for all $n = 1 : N$

Policy Evaluation: Value Computation via Recursion

If it happens that the claimed $v_\pi(\cdot)$ is **not** a **valid claim**; then, we get out of Bellman equation

$$\bar{v}_\pi(s^n) = \bar{v}_n = \bar{\mathcal{R}}_\pi(s^n) + \gamma \mathbb{E}_\pi \{v_\pi(\bar{S}) | s^n\}$$

which is different from the claimed $v_\pi(\cdot)$, i.e., $v_n \neq \bar{v}_n$

Policy Evaluation

We iterate this procedure until we can confirm, i.e., we

- 1 set $v_\pi(\cdot) \leftarrow \bar{v}_\pi(\cdot)$
- 2 repeat the same procedure and compute **new** $\bar{v}_\pi(\cdot)$

We **stop** when $v_\pi(\cdot) = \bar{v}_\pi(\cdot)$, or at least it happens approximately

Policy Evaluation

```

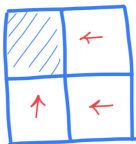
PolicyEval( $\pi, v_\pi^0$ ):
1: Initiate values with  $v_\pi^0$  and set  $k = 0$ 
2: Make sure that  $v_\pi^0(s) = 0$  for terminal states  $s$ 
3: Choose a small threshold  $\epsilon$  and initiate  $\Delta = +\infty$  # stopping criteria
4: for  $n = 1 : N$  do
5:   Compute  $\bar{\mathcal{R}}_\pi(s^n) = \mathbb{E}_\pi\{\bar{\mathcal{R}}(s^n, a)\}$  # average rewards
6: end for
7: while  $\Delta > \epsilon$  do
8:   for  $n = 1 : N$  do
9:     Update  $v_\pi^{k+1}(s^n) = \bar{\mathcal{R}}_\pi(s^n) + \gamma \mathbb{E}_\pi\{v_\pi^k(\bar{S}) | s^n\}$  # DP update
10:   end for
11:    $\Delta = \max_n |v_\pi^{k+1}(s^n) - v_\pi^k(s^n)|$  # check convergence
12:   Update  $k \leftarrow k + 1$ 
13: end while # Recursion Loop

```

Attention

We should make sure that **terminal states** are all initiated with **zero** value

Example: Dummy Grid World



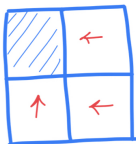
Let's try with our dummy grid world: we saw that

$$\bar{\mathcal{R}}_{\pi}(0) = 0 \quad \bar{\mathcal{R}}_{\pi}(1) = -1 \quad \bar{\mathcal{R}}_{\pi}(2) = -1 \quad \bar{\mathcal{R}}_{\pi}(3) = -1$$

Now let's *evaluate its values* by *recursion*: we first note that, if we have

$$\begin{aligned} \mathbb{E}_{\pi} \left\{ v_{\pi}^k(\bar{S}) \mid 0 \right\} &= v_{\pi}^k(0) & \mathbb{E}_{\pi} \left\{ v_{\pi}^k(\bar{S}) \mid 1 \right\} &= v_{\pi}^k(0) \\ \mathbb{E}_{\pi} \left\{ v_{\pi}^k(\bar{S}) \mid 2 \right\} &= v_{\pi}^k(0) & \mathbb{E}_{\pi} \left\{ v_{\pi}^k(\bar{S}) \mid 3 \right\} &= v_{\pi}^k(2) \end{aligned}$$

Example: Dummy Grid World



PolicyEval(π, v_π^0):

- 1: Initiate values with $v_\pi^0(1)$, $v_\pi^0(2)$ and $v_\pi^0(3)$ *at random* and *set* $v_\pi^0(0) = 0$
- 2: Set $\epsilon = 0.001$, and initiate $\Delta = 1000$ # stopping criteria
- 3: **while** $\Delta > \epsilon$ **do**
- 4: Update $v_\pi^{k+1}(1) = -1 + v_\pi^k(0)$ # DP update
- 5: Update $v_\pi^{k+1}(2) = -1 + v_\pi^k(0)$ # DP update
- 6: Update $v_\pi^{k+1}(3) = -1 + v_\pi^k(2)$ # DP update
- 7: $\Delta = \max_{s \in \{1,2,3\}} |v_\pi^{k+1}(s) - v_\pi^k(s)|$ # check convergence
- 8: Update $k \leftarrow k + 1$
- 9: **end while**

It converges after only *one* recursion!

Policy Improvement

Let us now recall **optimality constraint**: with optimal policy, we have

$$v_{\star}(s) = \max_m q_{\star}(s, a^m)$$

which can be achieved by policy

$$\pi^{\star}(a^m | s) = \begin{cases} 1 & m = \operatorname{argmax}_m q_{\star}(s, a^m) \\ 0 & m \neq \operatorname{argmax}_m q_{\star}(s, a^m) \end{cases}$$

This means that if π is **not** optimal, we would have

$$\pi(a^m | s) \neq \begin{cases} 1 & m = \operatorname{argmax}_m q_{\pi}(s, a^m) \\ 0 & m \neq \operatorname{argmax}_m q_{\pi}(s, a^m) \end{cases}$$

Policy Improvement

In other words, if we change our policy to

$$\bar{\pi}(a^m | s) = \begin{cases} 1 & m = \operatorname{argmax}_m q_{\pi}(s, a^m) \\ 0 & m \neq \operatorname{argmax}_m q_{\pi}(s, a^m) \end{cases}$$

Then, it should give us better values, i.e., $\bar{\pi} \geq \pi$!

- + *Are you sure?! I don't see it immediately*
- We can actually show it!

This is what we call *policy improvement theorem*

Policy Improvement

Policy Improvement

Given (deterministic) policy π^k , we can always design a *better* policy π^{k+1} by setting it to

$$\pi^{k+1}(a^m | s) = \begin{cases} 1 & m = \operatorname{argmax}_m q_{\pi^k}(s, a^m) \\ 0 & m \neq \operatorname{argmax}_m q_{\pi^k}(s, a^m) \end{cases}$$

Policy Improvement

PolicyImprov(v_π):

```

1: for  $n = 1 : N$  do
2:   for  $m = 1 : M$  do
3:     Compute  $\bar{R}(s^n, a^m)$ 
4:      $q_\pi(s^n, a^m) = \bar{R}(s^n, a^m) + \gamma \mathbb{E}_\pi \{v_\pi(\bar{S}) | s^n, a^m\}$    # action-values
5:   end for
6:   Compute an improved policy as                                     # policy improvement

```

$$\bar{\pi}(a^m | s^n) = \begin{cases} 1 & m = \underset{m}{\operatorname{argmax}} q_\pi(s^n, a^m) \\ 0 & m \neq \underset{m}{\operatorname{argmax}} q_\pi(s^n, a^m) \end{cases}$$

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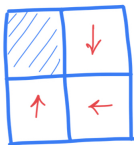
7: end for

```

Attention

Here, we do **no recursion**

Example: Dummy Grid World



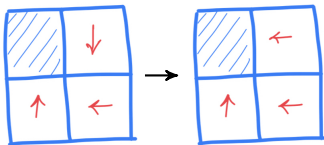
Let's try dummy grid world with above non-optimal policy: here, we have

$$v_{\pi}(0) = 0 \quad v_{\pi}(1) = -3 \quad v_{\pi}(2) = -1 \quad v_{\pi}(3) = -2$$

We now look at *action-values* at the *problematic state* $s = 1$

$$\begin{aligned} q_{\pi}(1, 0) &= -1 \\ q_{\pi}(1, 1) &= -3 \\ q_{\pi}(1, 2) &= -3.5 \\ q_{\pi}(1, 3) &= -3.5 \end{aligned} \quad \rightsquigarrow \quad -3 = v_{\pi}(1) \neq \max_a q_{\pi}(1, a) = -1$$

Example: *Dummy Grid World*



Now if we improve the policy, we get

$$\bar{\pi}(a|1) = \begin{cases} 1 & m = \operatorname{argmax}_m q_{\pi}(1, a) \\ 0 & m \neq \operatorname{argmax}_m q_{\pi}(1, a) \end{cases} = \begin{cases} 1 & a = 0 \\ 0 & a \neq 0 \end{cases}$$

which is actually *optimal*

Policy Iteration: Improving Policy by Recursion

Looking at the **policy improvement** theorem, we see

If we plug in $\pi^k = \pi^*$ into algorithm; then, after policy improvement

↳ we get $\pi^{k+1} = \pi^*$

↳ say we **evaluate** values for $\pi^{k+1} = \pi^*$ and plug back to algorithm

↳ we get $\pi^{k+2} = \pi^*$

↳ say we **evaluate** values for $\pi^{k+2} = \pi^*$ and plug back to algorithm

↳ ...

So, **optimal policy** is a **fixed-point** for this **recursion**

Policy Iteration

We can start with an **arbitrary policy** π^0 and keep doing the above recursion until we see that $\pi^{k+1} = \pi^k$ which indicates that we reached **optimal policy**

Policy Iteration

PolicyIter():

1: Initiate with **random** $v_\pi(s)$ for all **non-terminal** states s

2: Set $v_\pi(s) = 0$ for **terminal** states s

3: Initiate two random policies π and $\bar{\pi}$

4: **while** $\pi \neq \bar{\pi}$ **do**

5: $v_\pi = \text{PolicyEval}(\pi, v_\pi)$ and $\pi \leftarrow \bar{\pi}$ **Recursion**

6: $\bar{\pi} = \text{PolicyImprov}(v_\pi)$

7: **end while**

Recursion

Note that this is a **nested** recursive computation

- There is a loop for recursion inside the algorithm in which
 - ↳ at each iteration we **evaluate the policy recursively**
- But, we initiate each **policy evaluation** loop with the values of last iteration
 - ↳ this can improve the convergence speed

Back-Tracking by Recursion

- + *But wait a Moment! We already talked about back-tracking optimal policy from **Bellman optimality equation**! Don't we implement that?!*
- Sure! We can do the same thing by recursion

We follow the same idea but we use recursion

- 1 We can find optimal values from **Bellman optimality equations**
↳ This is where we use **recursion**
- 2 We could then find the **optimal action**-values
- 3 We finally get the **optimal policy** from **optimal action**-values

Recall: Back-Tracking from Optimal Values

OptimBackTrack():

```

1: Solve Bellman equations                                     # we use recursion
2: for  $n = 1 : N$  do
3:   for  $m = 1 : M$  do
4:     Set  $q_{\star}(s^n, a^m) = \bar{\mathcal{R}}(s^n, a^m) + \gamma \mathbb{E} \{v_{\star}(\bar{S}) | s^n, a^m\}$  # action-values
5:   end for
6:   Compute optimal policy via optimality constraint

```

$$\pi^{\star}(a^m | s) = \begin{cases} 1 & m = \underset{m}{\operatorname{argmax}} q_{\star}(s, a^m) \\ 0 & m \neq \underset{m}{\operatorname{argmax}} q_{\star}(s, a^m) \end{cases}$$

7: end for

Recursion with Bellman Optimality

Recall Bellman optimality equation

$$v_{\star}(s) = \max_m \left(\bar{\mathcal{R}}(s, a^m) + \gamma \mathbb{E} \{ v_{\star}(\bar{S}) | s, a^m \} \right)$$

We can again solve it by recursion: we start with some $v_{\star}^0(\cdot)$ and then for every state s and action a^m , we compute

$$\mathbb{E} \{ v_{\star}^k(\bar{S}) | s, a^m \} = \sum_{n=1}^N \underbrace{v_{\star}^k(s^n)}_{\text{last computed value}} \underbrace{p(s^n | s, a^m)}_{\text{transition model}}$$

We then update the optimal value function as

$$v_{\star}^{k+1}(s) = \max_m \left(\bar{\mathcal{R}}(s, a^m) + \gamma \mathbb{E} \{ v_{\star}^k(\bar{S}) | s, a^m \} \right)$$

Value Iteration vs Policy Iteration

Before we complete the value iteration algorithm: *it is interesting to put its recursion next to the one used for policy evaluation*

With optimality equation, we iterate as

$$v_{\star}^{k+1}(s) = \max_m \left[\bar{\mathcal{R}}(s, a^m) + \gamma \mathbb{E} \left\{ v_{\star}^k(\bar{S}) \mid s, a^m \right\} \right]$$

With Bellman equation for a given policy π , we iterate as

$$\begin{aligned} v_{\pi}^{k+1}(s) &= \bar{\mathcal{R}}_{\pi}(s) + \gamma \mathbb{E}_{\pi} \left\{ v_{\pi}^k(\bar{S}) \mid s \right\} \\ &= \sum_{m=1}^M \left(\bar{\mathcal{R}}(s, a^m) + \gamma \mathbb{E} \left\{ v_{\pi}^k(\bar{S}) \mid s, a^m \right\} \right) \pi(a^m | s) \end{aligned}$$

Value Iteration vs Policy Iteration

With optimality equation, we iterate as

$$v_{\star}^{k+1}(s) = \max_m \left[\bar{\mathcal{R}}(s, a^m) + \gamma \mathbb{E} \left\{ v_{\star}^k(\bar{S}) \mid s, a^m \right\} \right]$$

With Bellman equation for a given policy π , we iterate as

$$v_{\pi}^{k+1}(s) = \sum_{m=1}^M \left(\bar{\mathcal{R}}(s, a^m) + \gamma \mathbb{E} \left\{ v_{\pi}^k(\bar{S}) \mid s, a^m \right\} \right) \pi(a^m | s)$$

This indicates that for both recursive loops

- we compute M values *per iteration per state*
 - ↳ in policy iteration, we compute the *average* of these M via π
 - ↳ in value iteration, we take the *largest* among these M values

Value Iteration

ValueIter():

1: Initiate with **random** $v_{\star}^0(s)$ for all states, and set $v_{\star}^0(s) = 0$ for **terminal** states

2: Choose a **small** threshold ϵ , initiate $\Delta = +\infty$ and $k = 0$

3: **while** $\Delta > \epsilon$ **do**

4: **for** $n = 1 : N$ **do**

5: **for** $m = 1 : M$ **do**

6: Compute $q_{\star}(s^n, a^m) = \bar{\mathcal{R}}(s^n, a^m) + \gamma \mathbb{E} \{ v_{\star}^k(\bar{S}) | s^n, a^m \}$

7: **end for**

8: Update $v_{\pi}^{k+1}(s^n) = \max_m q_{\star}(s^n, a^m)$ # DP update

9: **end for**

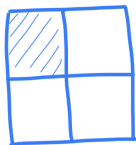
10: Set $\Delta = \max_n |v_{\pi}^{k+1}(s^n) - v_{\pi}^k(s^n)|$ and $k \leftarrow k + 1$

11: **end while** Recursion

12: Compute an optimal policy as

$$\bar{\pi}(a^m | s) = \begin{cases} 1 & m = \operatorname{argmax}_m q_{\star}(s, a^m) \\ 0 & m \neq \operatorname{argmax}_m q_{\star}(s, a^m) \end{cases}$$

Example: *Dummy Grid World*



You may try policy and value iteration for this problem at home!

Easy as Pie 😊

Example: A Bit Larger Grid World¹

0	1	2	3
4	5	6	7
8	9	10	11
12	13	14	15

Board \equiv states



moves \equiv actions

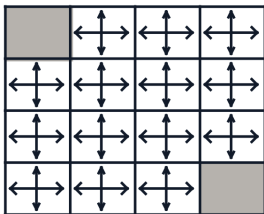
Let's do a bit of more serious example: we are now in a 4×4 grid world

- We have two **terminal states** shown in gray
- Each move we do gets a -1 reward
 - ↳ We also get -1 reward if we hit a corner
 - ↳ We get zero reward at terminal state

In simple words: we are looking for **shortest path** to the **corners**

¹This example is taken from Sutton and Barto's Book; Example 4.1 in Chapter 4

Example: A Bit Larger Grid World



initial policy

0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0

initial values

Let's first try *policy iteration*: we start with

- a uniform random policy π^0
- all values being zero, i.e., $v_{\pi^0}^0(s) = 0$ for all s

Example: A Bit Larger Grid World

Recall policy iteration:

PolicyIter():

1: Initiate with **random** $v_\pi(s)$ for all **non-terminal** states s

2: Set $v_\pi(s) = 0$ for **terminal** states s

3: Initiate two random policies π and $\bar{\pi}$

4: **while** $\pi \neq \bar{\pi}$ **do**

5: $v_\pi = \text{PolicyEval}(\pi, v_\pi)$ and $\pi \leftarrow \bar{\pi}$ **Recursion**

6: $\bar{\pi} = \text{PolicyImprov}(v_\pi)$

7: **end while**

Recursion

We should start with $v_{\pi^0}^0(\cdot)$ and do the **red recursion** first

- at the end of this recursion we have evaluated the random policy

Example: A Bit Larger Grid World

0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0

 $v_{\pi^0}^0$

0.0	-1.0	-1.0	-1.0
-1.0	-1.0	-1.0	-1.0
-1.0	-1.0	-1.0	-1.0
-1.0	-1.0	-1.0	0.0

 $v_{\pi^0}^1$

0.0	-1.7	-2.0	-2.0
-1.7	-2.0	-2.0	-2.0
-2.0	-2.0	-2.0	-1.7
-2.0	-2.0	-1.7	0.0

 $v_{\pi^0}^2$

0.0	-2.4	-2.9	-3.0
-2.4	-2.9	-3.0	-2.9
-2.9	-3.0	-2.9	-2.4
-3.0	-2.9	-2.4	0.0

 $v_{\pi^0}^3$

...

0.0	-14.	-20.	-22.
-14.	-18.	-20.	-20.
-20.	-20.	-18.	-14.
-22.	-20.	-14.	0.0

 $v_{\pi^0}^{\infty}$

We now have *evaluated* the value of random policy $v_{\pi^0} = v_{\pi^0}^{\infty}$

Example: A Bit Larger Grid World

Recall policy iteration:

PolicyIter():

1: Initiate with **random** $v_\pi(s)$ for all **non-terminal** states s

2: Set $v_\pi(s) = 0$ for **terminal** states s

3: Initiate two random policies π and $\bar{\pi}$

4: **while** $\pi \neq \bar{\pi}$ **do**

5: $v_\pi = \text{PolicyEval}(\pi, v_\pi)$ and $\pi \leftarrow \bar{\pi}$ **Recursion**

6: $\bar{\pi} = \text{PolicyImprov}(v_\pi)$

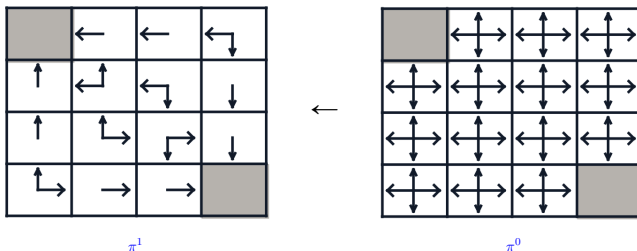
7: **end while**

Recursion

Next, we do the **outer recursion** recursion, i.e.,

- we **improve** the policy

Example: A Bit Larger Grid World



We improve policy by taking actions with **maximal action-values**

- if we have multiple **maximal action-values** we can behave **randomly**

Example: A Bit Larger Grid World

Recall policy iteration:

PolicyIter():

1: Initiate with **random** $v_\pi(s)$ for all **non-terminal** states s

2: Set $v_\pi(s) = 0$ for **terminal** states s

3: Initiate two random policies π and $\bar{\pi}$

4: **while** $\pi \neq \bar{\pi}$ **do**

5: $v_\pi = \text{PolicyEval}(\pi, v_\pi)$ and $\pi \leftarrow \bar{\pi}$ **Recursion**

6: $\bar{\pi} = \text{PolicyImprov}(v_\pi)$

7: **end while**

Recursion

We now $v_{\pi^1}^0 = v_{\pi^0} = v_{\pi^0}^\infty$ and do the **red recursion** again

- at the end of this recursion we have evaluated the **new policy** π^1

Example: A Bit Larger Grid World

0.0	-14.	-20.	-22.
-14.	-18.	-20.	-20.
-20.	-20.	-18.	-14.
-22.	-20.	-14.	0.0

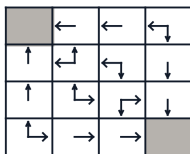
 $v_{\pi^1}^0$

...

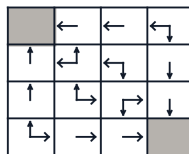
0.0	-1.0	-2.0	-3.0
-1.0	-2.0	-3.0	-2.0
-2.0	-3.0	-2.0	-1.0
-3.0	-2.0	-1.0	0.0

 $v_{\pi^1}^{+\infty}$

After *evaluating* policy π^1 as $v_{\pi^1} = v_{\pi^1}^{\infty}$, we do the next improvement

 $\pi^2 = \pi^1$

←

 π^1

Well $\pi^2 = \pi^1$ and we should *stop!*

Example: A Bit Larger Grid World

0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0

initial values

Now we try *value iteration*: for start, we only need an initial value, so we set

- all values being zero, i.e., $v_{\star}^0(s) = 0$ for all s

We keep recursion until we find the *optimal values*

Example: A Bit Larger Grid World

0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0
0.0	0.0	0.0	0.0

 v_{\star}^0

...

0.0	-1.0	-2.0	-3.0
-1.0	-2.0	-3.0	-2.0
-2.0	-3.0	-2.0	-1.0
-3.0	-2.0	-1.0	0.0

 $v_{\star}^{+\infty}$

Now, we back-track the optimal policy π^{\star}

0.0	-1.0	-2.0	-3.0
-1.0	-2.0	-3.0	-2.0
-2.0	-3.0	-2.0	-1.0
-3.0	-2.0	-1.0	0.0

 v_{\star} 

action-values



	←	←	↙
↑	↖	↙	↓
↑	↗	↘	↓
↖	→	→	

 π^{\star}

Complexity of Policy and Value Iteration

- + It seems that value iteration has *less complexity*!
- Well, it is not in order, but yes! It usually converge faster

In our example with *policy iteration*, we had to *evaluate two policies*

- once for π^0 and once for π^1
- say the first recursion took K_1 iterations and the second took K_2
 - ↳ the *total number* of iterations is then $K_1 + K_2$
 - ↳ in practice, it often happens that $K_2 \ll K_1$
 - ↳ because we already start from *good values* with $v_{\pi^1}^0 = v_{\pi^0}^{+\infty}$

With *value iteration*, we had to *only evaluate optimal policy*

- say it takes K_* iterations: there is *no reason* that K_* be same as K_1 or K_2
 - ↳ each evaluation has a *different initial* and *converging* point
- ↳ in practice, it often happens that $K_* > K_1$ and $K_* \gg K_2$
 - ↳ so it *might be* that $K_* \approx K_1 + K_2$
 - ↳ but usually with *multiple policy improvements*, we see $K_* < K_1 + K_2 + \dots$

Complexity of Policy and Value Iteration

- + If so, why should we use *policy iteration*?!
 - Well, not all problems are like a dummy grid world

In practice, it might be computationally *hard* to get *very close to optimal values*

- in this case, we take non-converged values
 - ↳ we consider them *estimates* of optimal values
- in value iteration we *approximate* optimal policy with on these *estimates*
 - ↳ this might be a *loose* estimate

If we do the same *approximative* computation with policy iteration

- we often end up with a *better policy*

Moral of Story

While *value iteration* typically show *faster convergence*, *policy iteration* can give *better policies* after convergence

Generalized Policy Iteration

In practice, we can terminate or change the order of computation in policy iteration to reduce its complexity: for instance, we could have

GenPolicyItr():

1: Initiate with **random** $v_\pi(s)$ for all **non-terminal** states s

2: Set $v_\pi(s) = 0$ for **terminal** states s

3: Initiate two random policies π and $\bar{\pi}$

4: **while** $\pi \neq \bar{\pi}$ **do**

5: $v_\pi = \text{TerminPolicyEval}(\pi, v_\pi)$ and $\pi \leftarrow$ **changed**

6: $\bar{\pi} = \text{PolicyImprov}(v_\pi)$

7: **end while**

where $\text{TerminPolicyEval}(\pi, v_\pi)$ evaluates **policy** π from starting value function v_π with a **terminating** recursion loop

Generalized Policy Iteration: *Terminating Evaluation*

TerminPolicyEval(π, v_π^0):

```

1: Initiate values with  $v_\pi^0$  and set  $k = 0$ 
2: Make sure that  $v_\pi^0(s) = 0$  for terminal states  $s$ 
3: Choose a small threshold  $\epsilon$  and initiate  $\Delta = +\infty$  # stopping criteria
4: for  $n = 1 : N$  do
5:   Compute  $\bar{\mathcal{R}}_\pi(s^n) = \mathbb{E}_\pi \{ \bar{\mathcal{R}}(s^n, a) \}$  # average response
6: end for
7: while  $\Delta > \epsilon$  and  $k < K$  do changed
8:   for  $n = 1 : N$  do
9:     Update  $v_\pi^{k+1}(s^n) = \bar{\mathcal{R}}_\pi(s^n) + \gamma \mathbb{E}_\pi \{ v_\pi^k(\bar{S}) | s^n \}$  # DP update
10:   end for
11:    $\Delta = \max_n |v_\pi^{k+1}(s^n) - v_\pi^k(s^n)|$  # check convergence
12:   Update  $k \leftarrow k + 1$ 
13: end while

```

Obviously, TerminPolicyEval(π, v_π) does **not** return the **exact values** of the policy π , but only an **estimate** of them

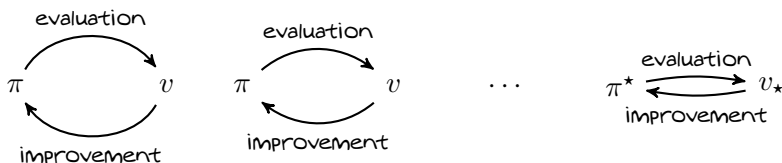
Generalized Policy Iteration

We can come up with various such ideas: *these variants are often called*

Generalized Policy Iteration \equiv GPI

These approaches all rely on

back-and-forth computation of policies and values



If designed properly, they all converge to optimal policy and optimal values

Some Final Remarks

- + We know the algorithms now, but how can we **guarantee** that they **converge**? You showed us an simple example that recursion could simply **diverge**!
- Well, we can show that what we discussed in this chapter converge: it comes from the nice properties of **Bellman equations**
 - ↳ There are several proofs; for instance see [a proof in Tom Mitchell's notes](#)

When it comes to practice, most known algorithms are proved to **converge to optimal policy and optimal values**; however, note that

- **Convergence guarantee** is different from the **speed of convergence**
 - ↳ An algorithm might **converge**, but **very slow**
- If you deal with an **unknown** algorithm; then, you should make sure that it **converges to optimal policy and optimal values**