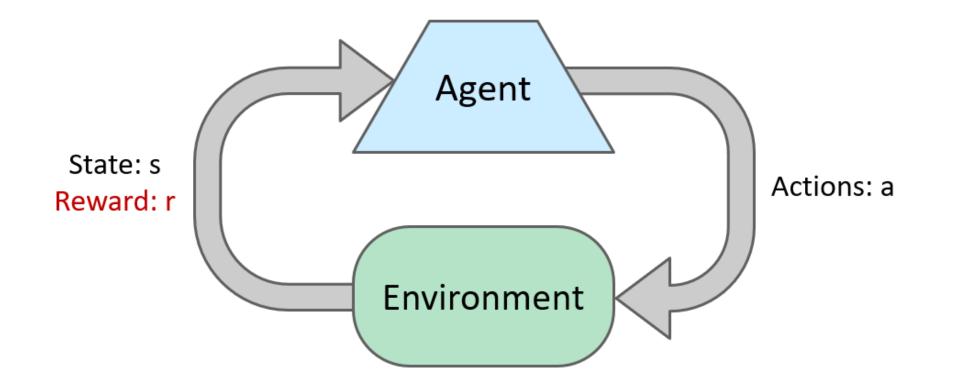
### **Reinforcement learning**

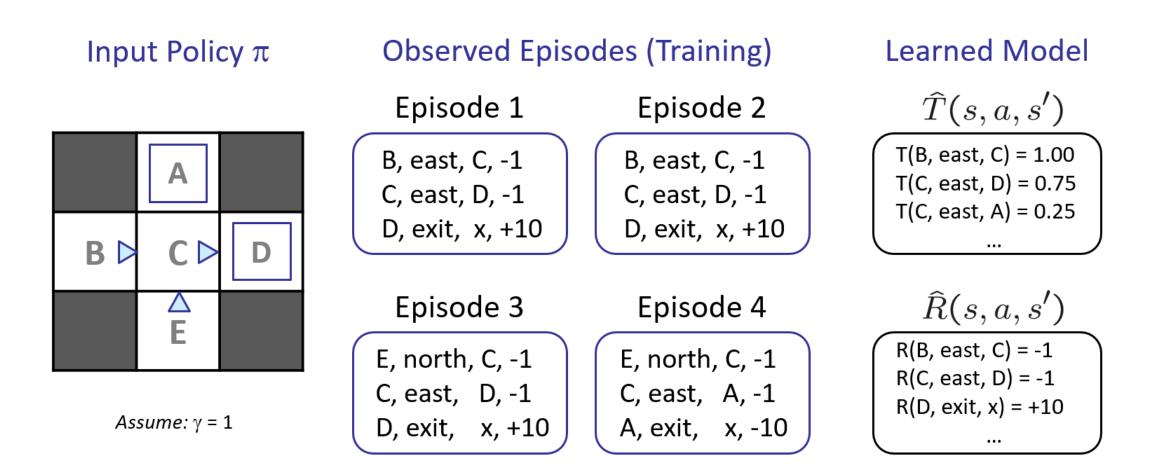


### **RL model**

- We assume a world governed by an MDP
  - $\circ\;$  States  $s\in S$
  - $\,\circ\,$  Actions A
  - $\,\circ\,$  System dynamics T(s,a,s')
  - $\circ\,$  Reward function R(s,a,s')
- We are looking for the policy  $\pi(s)$
- But we don't know T and/or R as functions
- But if we take an action *a*, in state *s*, we can observe that we landed in *s'* and received reward *r*.

#### **Model-based RL**

- General idea: if there is a MDP M in the environment, we can find a way to build an approximate model  $\hat{M}$  that approximates it.
  - $\,\circ\,$  Perform (or even better, observ) some actions a , count the outcomes s' and r
  - $\circ\,$  Normalize, to get the estimate of  $\hat{T}(s,a,s')$
  - $\circ~$  Calculate the associated average reward to get  $\hat{R}(s,a,s')$
- Then, we can solve  $\hat{M}$  using some technique (eg. value iteration or policy iteration)



#### **Model-based RL**

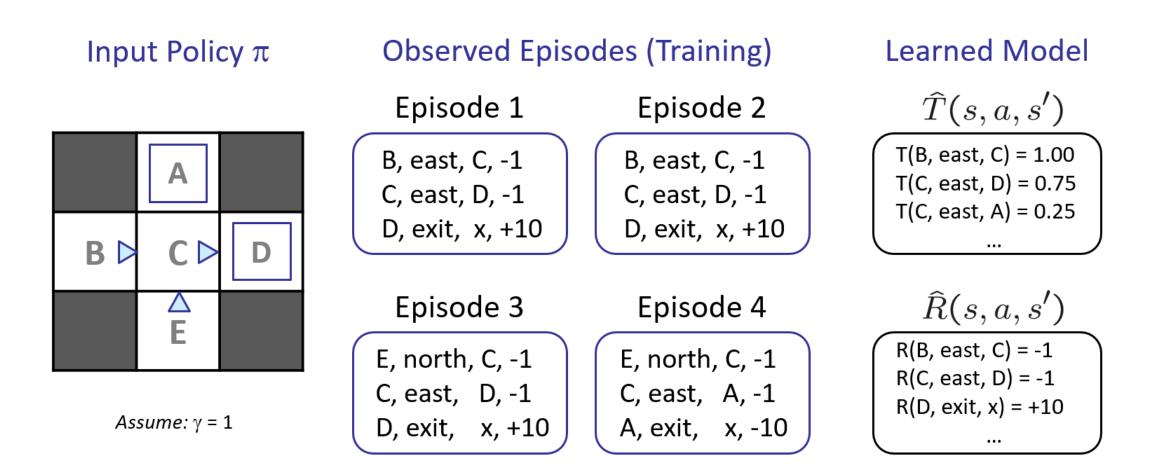
- Why it is good?
  - We didn't need to invent any new techniques
  - $\circ\,$  As an extra benefit, we get the T and R, which might come handy
  - $\circ~$  It is very convenient if we can get the data from somebody else's runs
  - $\circ~$  It can be convenient, if we have some prior information about the T or R
- Problems
  - $\circ~$  Often, we don't care about the T and R getting a decent policy should be much simpler than having a full model!

### **Model free RL**

- Idea: we can try to evaluate the Q value (or something of this sort) without bothering to find the T or R.
  - $\,\circ\,$  Then we can extract the  $\pi$  from the Q
  - This is going to get us variations of Q-learning
    - Another name: **critic only** algorithms
- Another idea: we can go directly to π without bothering to find the Q
   This is going to get as variations of **policy gradient** approaches
- Or we can try to do both simultaneously (**actor-critic algorithms**) [LAST TWO: NOT COVERED IN THIS CLASS]

### **Direct evaluation of V**

- Let us assume we have a policy  $\pi$
- We have a  $\operatorname{\mathsf{run}} s_1, a_1, r_1, s_2, a_2 \dots r_n, s_n$
- Let us say we are in  $s_k$ 
  - $\circ\,$  Calculate the discounted rewards to the end of the run $G=r_k+\gamma r_{k+1}+\gamma^2 r_{k+1}\ldots\gamma^{n-k-1}r_n$
- For all states s which have at least one run passing through them, average these g values, that will be the estimate  $\hat{V}(s)$
- This is a Monte Carlo method



### **Direct evaluation of V**

- The good
  - Simple to understand
  - $\circ~$  Doesn't require knowledge of T or R
  - $\circ\,$  It converges to the correct values of the  $V^{\pi}(s)$  given enough runs that pass through the state
- The bad
  - For rarely visited states, it can take a long while
  - It does not exploit knowledge about the connections between states: for all purposes, it learns each state separately.
  - In the previous figure: why are the values of E and B so different, when they both only go to C?

## **Temporal difference learning (TD-learning)**

- We are still doing policy evaluation, learning the  $V^{\pi}$  for a fixed  $\pi$
- Instead of waiting for a run to finish, learn from every transition (s,a,s',r)
  - $\circ\,$  The idea is that we **bootstrap** the V(s) with the value of V(s')
  - **Temporal difference**: we go back to our history and update those values
- A **sample** (notice the similarity to be Bellman equation...) or **TD-target**

$$sample = R(s, \pi(s), s') + \gamma V^{\pi}(s')$$

• TD-update:

$$V^{\pi}(s) \leftarrow (1-lpha) V^{\pi}(s) + lpha \cdot sample$$

or

$$V^{\pi}(s) \leftarrow V^{\pi}(s) + lpha(sample - V^{\pi}(s))$$

## Wait, will this thing converge?

- $\alpha$  is the learning rate: if  $\alpha = 1$ , we will keep replacing the value with a value based on the next state, whatever that turns out to be this time...
- We need to decrease the  $\alpha$  for it to converge
- A way to think about TD-learning is as follows:
  - $\circ$  we are actually already in s' but we are updating s (the previous one... this is the temporal difference)
  - $\circ\,$  we expect the sample to be  $V^{\pi}(s)$  ()
  - $\circ~$  if you got a very large reward, it makes the sample larger than expected
  - $\circ~$  this makes the sample-V large, and a part of it will be used to increase  $V^{\pi}$  such that next time you will not be surprised.
- Written like this, the reward is propagated only to the previous step (in one event) but you can develop variants that go back multiple steps

### **TD-learning limitations**

- Basically, we are mimicking the Bellman updates with running sample averages
- Notice that this is an **on-policy** algorithm: we are following  $\pi$  and learning  $V^{\pi}$
- The good: we are learning the  $V^{\pi}$  without a model!
- A problem
  - $\circ~$  We can do policy iteration starting from this, but we need T and R
  - So we still need a model, or an approximation of it

## Are we really falling into the fire-pit?

- The  $\pi$  value was fixed
- The good: with this approach we can learn from other people's experiences
- The bad: we don't control where we are going

#### **Active RL**

- Setup: we choose the actions now!
- T(s,a,s') and R(s,a,s') are unknown
- Goal: learn  $\pi^*$ ,  $V^*$ ,  $Q^*$
- This time, the learner actually falls into the firepit.
- The choices of actions made during learning matter!
  - For instance, exploration...

#### **Q-value iteration?**

• We learned V-value iteration

$$V_{k+1}(s) \leftarrow \max_a \sum_{s'} ig( R(s,a,s') + \gamma V_k(s') ig)$$

• But we can also do it for Q-values

$$Q_{k+1}(s,a) \leftarrow \sum_{s'} T(s,a,s') \left( R(s,a,s') + \gamma \max_{a'} Q_k(s',a') 
ight)$$

• Why don't we do this for a known MDP? It is slower!

## **Q-learning**

• Do the Q-value iteration using samples! Sample is like the one for TD-learning, only replace V with max over Q.

$$sample = R(s, a, s') + \gamma \max_{a'} Q(s', a')$$

Update Q(s,a) based on the sample, with a learning rate:

$$Q(s,a) \leftarrow (1-lpha)Q(s,a) + lpha \cdot sample$$

## **Q-learning properties**

- Does it converge?
  - $\circ\,$  Yes. In fact it converges to  $Q^*$ , regardless of what policy you use!!!
  - off-policy learning
- With some limitations
  - $\circ~$  If you never perform an action a, you never learn about it
  - $\circ\,$  If you have never been in a state s, you never learning about it
  - $\circ \,\, \alpha$  need to decrease for convergence, but if it is too small, convergence is slow

## **Exploration vs. exploitation**

- In Q-learning you only learn about
  - actions you have actually took in
  - states you have actually been
  - This is a bit simplistic, but *roughly* true for most RL
- Exploitation: take advantage of what you learned
  - Take action according to the q-values they currently have attached
  - This means that in each state we always take the same action
  - $\circ~$  The only way we might land in a novel state is if the probabilistic T lands us there.
  - When we deploy the (fixed) learned policy, we should do pure exploitation

### **Exploration**

- As we only learn from states we have been and actions we took we should explore by taking a variety of actions
- First approximation: act randomly!  $a \sim U(a)$ 
  - Problem: in many scenarios, we reach unfavorable terminal states very soon, and never see advanced states closer to the goal!

### **Exploration (cont'd)**

- Reset and act randomly! Go to state s and try all actions from there  $a \sim U(a)$ . Repeat for all states of interest.
  - This is called **reset** to a state *s*
  - $\circ~$  It is actually a very good strategy when it works
  - It works in: simulation, board games, etc.
    - Whenever we have full control of the environment
    - ... and the negative rewards are make-believe
  - It doesn't work in: robotics, live finance, etc. whenever the negative rewards are real!

## **Exploration (cont'd)**

- Second approximation: act according to the current best policy and let the transition function take us to random places
  - $\circ$  We would still not see new actions taken a
  - The transition function might make some transitions extremely unlikely...

### **Stochastic policies**

- Make the policy  $\pi \operatorname{stochastic}$ 
  - $\circ\;$  Deterministic policy:  $\pi(s) 
    ightarrow a$
  - $\,\circ\,$  Stochastic policy:  $\pi(s,a) \to [0,1]$  the likelihood that the action a will be taken in state s
- There are scenarios where the optimal policy is a stochastic one
  - Eg. rock, paper, scissors...
  - There are algorithms that learn that.
- Alternatively, we can use a stochastic policy during learning, for exploration, and reset to a deterministic one during deployment.

### $\epsilon$ -greedy

- Consider a probability number  $\epsilon$  (eg. 0.2)
- At every time step in state  $\boldsymbol{s}$ 
  - $\circ\,$  with a probability  $1-\epsilon$  choose  $a=argmax\,Q(s,a)$
  - $\circ$  with probability  $\epsilon$  choose random a
- $\epsilon$  determines the balance between exploration and exploitation
- Often, we reduce  $\epsilon$  during learning.

### **Exploration functions and novelty seeking**

- A disadvantage of  $\epsilon$ -greedy is that it is taking random actions even in well known states.
- We would like to seek out novel information.
- One way to do it is to keep count of how many times we have been in a state and taken action count(s, a) and calculate a **novelty** function similar to this:

$$f(u,a) = u + rac{v}{count(s,a)+1}$$

• The exact form can vary, but the idea is that the less we visited a state/action pair, the more attractive it seems

#### **Using the exploration functions**

• We replace

$$Q_{t+1}(s,a) \leftarrow (1-lpha) Q_t(s,a) + lpha \left( R(s,a,s') + \gamma \max_{a'} Q_t(s',a') 
ight)$$

with

$$Q_{t+1}(s,a) \leftarrow (1-lpha) Q_t(s,a) + lpha \left( R(s,a,s') + \gamma \max_{a'} f\left(Q_t(s',a'), count(s',a')
ight) 
ight)$$

- The idea is that we give an extra novelty reward to unknown  $\boldsymbol{s}, \boldsymbol{a}$  combinations
- This novelty will propagate into the Q! So we will seek those out from afar!
- But in the long run, it will decay, and the real rewards will dominate.

### Summary: exploration vs exploitation

- Very important for all RL algorithms
- Plenty of human and animal analogies exist
  - Of various levels of plausibility and relevance
- It is almost sure that you will need to do at least  $\epsilon$ -greedy
- Recent years several high profile works in curiosity-driven exploration had became highly cited, as they were claiming success with approaches that learned with rewards coming **exclusively** from novelty.

### Regret

- We need to find a measure of the efficiency of learning
- We define regret as the gap between the learned policy and the optimal policy
- Let us consider that we are repeatedly solving a problem  $1 \dots K$ 
  - $\circ\,$  At each step, we start from  $s_k$  (chosen by nature or an adversary)
  - $\circ$  After each run, we are using RL to improve our policy,  $\pi_1 \dots \pi_K$

$$Regret = \sum_{k=1}^K \left( V^*(s_k) - V^{\pi_k}(s_k) 
ight)$$

- Algorithms that have lower regret are less costly to train than those with high regret
  - Even if the final policy and the number of runs is the same!

## **Approximate Q-learning**

- In a realistic situation, we cannot keep a table of Q values
  - $\circ\,$  Also, we cannot visit all of them in training.
- We want to generalize from experience to similar states
  - $\circ~$  Falling into a firepit in the library bad  $\rightarrow$  falling into the firepit bad in most locations
- It is helpful to consider Q(s,a) a function:
  - Can be implemented as a lookup table (as we did until now)
  - Can be a linear expression of **features**
  - Can be a **neural network** (deep RL)

#### **Feature-based representation**

- Idea: engineer a series of **features** that capture aspects of the problem
  - $\circ\;$  They are often functions f(s) or f(s,a) from state to a [0,1] or 0,1 range.
  - Traditionally, they are engineered by experts and capture human expertise and experience.
  - We only need to ask the expert to give us "things that are important", not to put an exact weight on them.
- Advantage: can represent large spaces with a few numbers!
- Disadvantage: if the set of the features does not capture everything about the state, states similar in features might be very diffrenet in value

#### **Examples**

- Examples for f(s):
  - Distance to the monster (normalized to maximum)
  - Distance to the exit (normalized to maximum)
  - Are we in a tunnel? (yes/no)
- Examples for f(s, a):
  - Moving towards monster? (yes/no)
  - Bumping into wall? (yes/no)

#### **Linear value functions**

$$V(s) = \sum_i^n w_i f_i(s) 
onumber \ Q(s,a) = \sum_i^n w_i f_i(s,a)$$

### **Approximate Q-learning with linear Q-functions**

• Line in exact Q-learning, we consider a sample of the transition (s,a,r,s')

$$ext{difference} = r + \gamma \max_a' Q(s',a') - Q(s,a)$$

The exact update rule was:

$$Q(s,a) \leftarrow Q(s,a) + lpha \cdot ext{difference}$$

The update rule will be now:

$$w_i \leftarrow w_i + lpha \cdot ext{difference} \cdot f_i(s,a)$$

# Intuitive interpretation of approximate Qlearning

- Note that we multiply the update with the feature value
  - Intuition: if we got a big negative reward, we blame the features that were present, not the ones that are not!
- Formal justification of this model is from the theory of online least squares