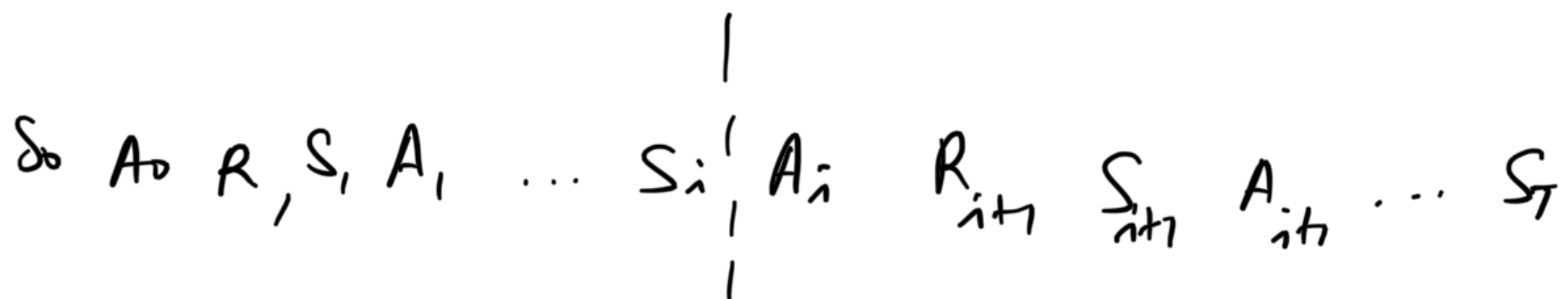
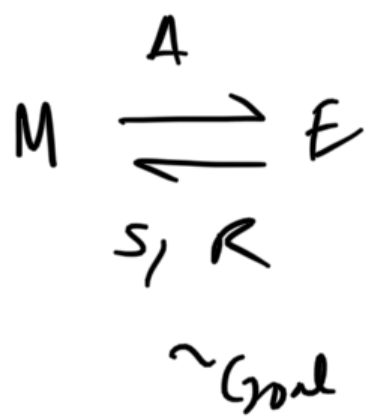


APS 1080 - LECTURE 3

MODEL BASED RL - DYNAMIC PROGRAMMING



What Action to take if you're in state i .

Return: Sum of subsequent rewards as a result of our action

$$G_i = R_{i+1} + \gamma R_{i+2} + \gamma^2 R_{i+3} + \dots$$

Goal: select A_i such that the expected value of the return is maximized.

\downarrow
 Action

$\max_{\pi} \mathbb{E} [G_i]$

VALUE FUNCTION

$$V_{\pi}(s) = \mathbb{E} [G_t \mid S_t = s]$$

$$q_{\pi}(s) = \dots$$

Given V_π

P : model of the environment

We can solve Action Selection Problem.

To solve Action Selection Problem, you need either:

① V and $P \Rightarrow \pi$

② $Q \Rightarrow \pi$



The Policy, which is the mechanism for action selection.

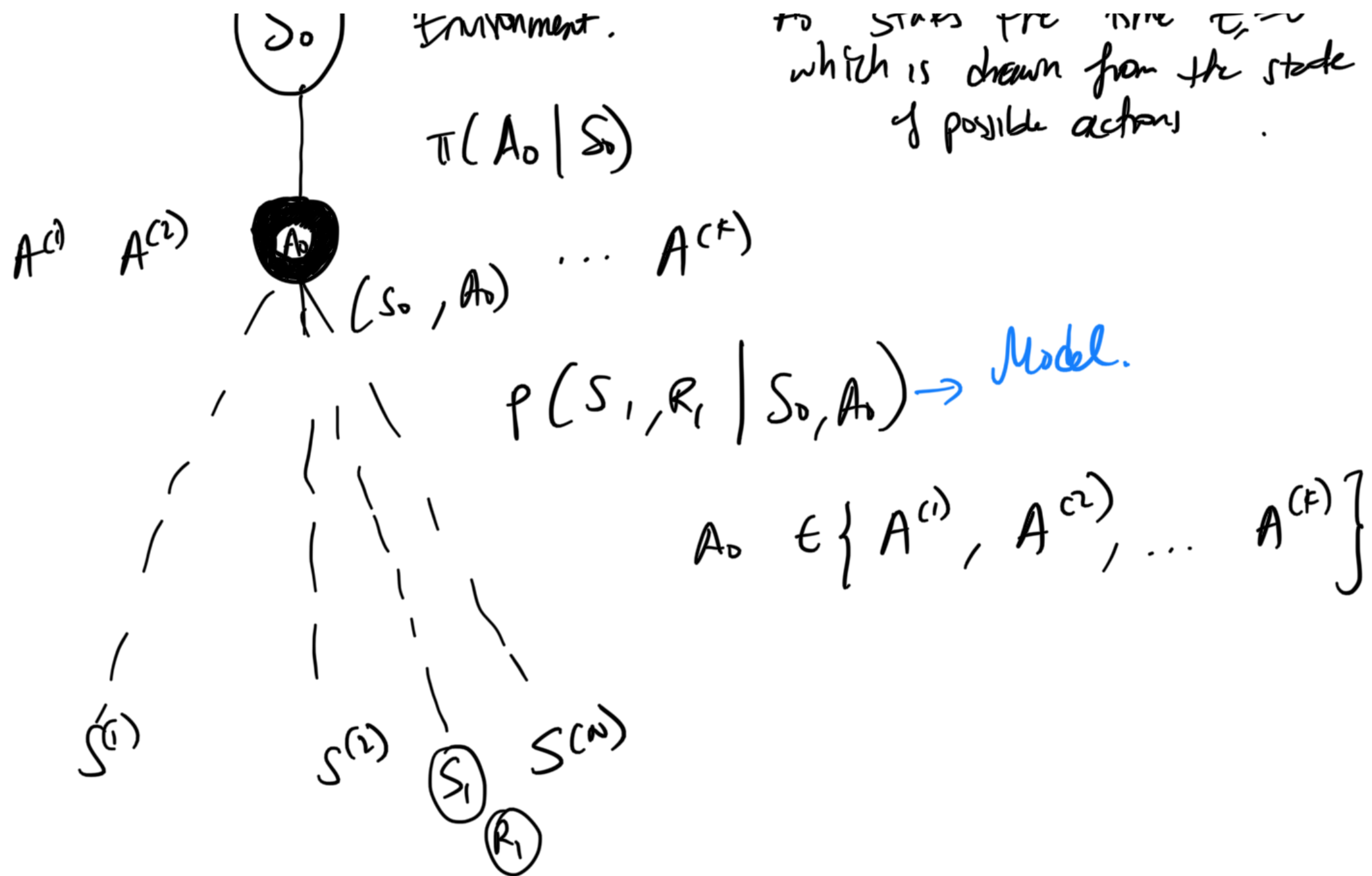
The states are drawn from $\mathcal{S} = \{S^{(1)}, S^{(2)}, \dots, S^{(N)}\}$

S	$V_\pi(S)$
$S^{(1)}$	$\mathbb{E}[G_t S_t = S^{(1)}]$
$S^{(2)}$	
\vdots	
$S^{(N)}$	

Target state:
which state to go to
then consult "P"
to see what action
to take to get
to that probability.

$$V_\pi(S) = \mathbb{E}[G_t | S_t = S]$$

Subscript: Time sequence
Superscript: Selection from the set.



This is how π and p interplay in continuous state/action rewards.

Action \rightarrow State (weighted by distribution p)

The model we may or may not have it.

The

with p we can calculate V_π , π and π^*
 \downarrow
 optimal policy.

Dynamic programming does NOT use experience

How we determine V_π and π when we don't have P ?

\rightarrow you need experience + learning to be surrogates of

$$V_{\pi}(s) = \sum_{a \in A} \pi(a|s) \sum_{s' \in S} \sum_{r \in R} [p(s', r | s, a) [r + \gamma V_{\pi}(s')]]$$

Probability of selecting an action given you're in state s .
 Indices of a for loop
 for loop
 consider all possible states
 over all possible rewards

Expected value \rightarrow Weighted Average over the entire tree.

You're building the Agent.

No knowledge about the problem $\rightarrow \pi$ would be an equiprobable distribution.

\rightarrow Eventually, you'd get to a more granular π , with a less naive distribution.

We have an equation that is recursively defined.

$$V_{\pi}(s) = \int V_{\pi}(\cdot)$$

\uparrow
gamma

We can calculate $V_{\pi}(\cdot)$ via linear system methods but it's not very scalable because you end up with too many variables and it's very computationally expensive. Also, it would only work if you have " p ".

Instead of doing that, we calculate $V_{\pi}(\cdot)$ via a successive approximation.

$$f(\cdot) = \mathcal{T}(f(\cdot))$$

↑
Contraction

$f(\cdot) \leftarrow$ initiate \leftarrow function $\forall x$

loop

loop $\forall x$

$$prev \leftarrow f(x)$$

$$next \leftarrow \mathcal{T}(prev)$$

$$f(x) \leftarrow next$$

$$\Delta(x) \leftarrow |next - prev|$$

$$\text{if } \|\Delta\| < \epsilon$$

\rightarrow break.

Initialize $V_{\pi}(x)$ for all x arbitrarily.

We set the terminal states to zero and all the other values at an arbitrary number.

States	V_π
$S^{(1)}$	
$S^{(2)}$	
\vdots	
$S^{(n)}$	

$$V_\pi(S_T) = 0$$

π	$a^{(1)}$ $a^{(2)}$... $a^{(k)}$	
$S^{(1)}$		$\Sigma = 1$
$S^{(2)}$		$\Sigma = 1$
\vdots		
$S^{(n)}$		

Input $\diamond \pi$

Init $V_\pi(s)$ arbitrarily $\forall s$

loop $\Delta \leftarrow 0$

| loop $\forall s$

| | prev $\leftarrow V_{\pi}(s)$

| | next $\leftarrow \Gamma[V_{\pi}(\cdot)]$

| | $\sum_{\pi} \dots \sum \sum [r + \gamma V_{\pi}(s)]$

| | $\Delta \leftarrow \max[\Delta, (prev - next)]$

| if $\Delta < \theta$:

Break

Iterate over the states, successively estimating the values of the value functions

When the deltas are low enough then you stop.

Given an arbitrary model and policy, we can determine a value function

$p, \pi \rightarrow V_{\pi}(\cdot)$

\rightarrow Evaluation of the policy.

Evaluation of π in $\mathcal{E} \sim$ "Prediction"

{ $\pi \xrightarrow{\text{eval}} V_{\pi}(\cdot)$
 $\pi \rightarrow \pi'$
 eval

$$\left\{ \begin{array}{l} \pi' \rightarrow V_{\pi'}(\cdot) \\ \pi' \xrightarrow[\text{Improvement}]{V_{\pi'}} \pi'' \end{array} \right.$$

$$\left\{ \begin{array}{l} \pi'' \rightarrow V_{\pi''}(\cdot) \\ \pi'' \rightarrow \pi''' \end{array} \right.$$

$$\left. \begin{array}{l} \text{until } V_{\pi}(i) \approx V_{\pi}(i+1) \\ \pi(i) \approx \pi(i+1) \end{array} \right\} \begin{array}{l} \text{Fixed Point} \\ \pi_{*} \end{array}$$

You modify the policy to help you improve the likelihood of going to the most desirable states.

Summarize Dynamic Programming

$$\textcircled{1} p, \pi_0 \rightarrow V_{\pi_0} \quad (\text{Prediction})$$

$$\textcircled{2} \pi_0, V_0 \rightarrow \pi_1 \quad (\pi_1 \geq \pi_0) \quad (\text{Improvement})$$

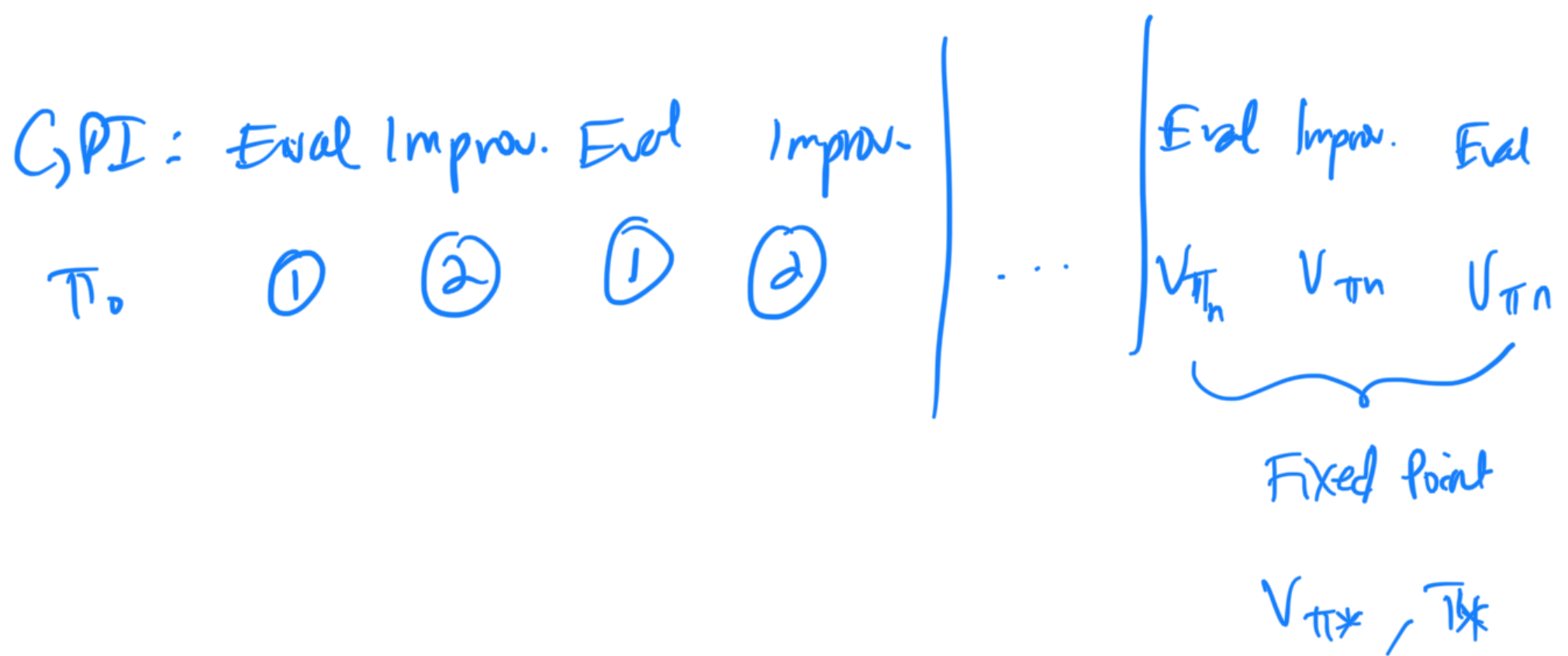
$\textcircled{3}$ Generalized Policy Iteration (GPI)

Prediction
Improvement

Until Fixed Point is reached $\rightarrow \pi_{*}$

$$\textcircled{4} \text{ GPI} \rightarrow \begin{array}{l} \text{Evaluation} \\ \text{Improvement} \end{array} \quad \begin{array}{l} \textcircled{1} \\ \textcircled{2} \end{array}$$

Evaluation
Improvement $\begin{array}{l} \textcircled{1} \\ \textcircled{2} \end{array}$



This is an asymptotic process

Other approach

Run it a few times and truncate the evaluation.

Select the best value amongst bag of truncated evaluations.

Step ④ helps us move away from the need of having a model.

What are other ways to approximate a value function that would not require the use of a "P"?

Next week:

How to approximate V_{π} without P by leveraging experience?

→ How can we approximate q_{π} and a policy from experience?

↳ π