CS 6300: Artificial Intelligence

Midterm Review

Midterm Logistics

- In our classroom during normal class time
 - Wednesday from 3-4:20pm
- I sheet of notes (front and back)
- Simple calculator allowed but not needed (all math will be simple)
- Lots of extra-credit.
 - Choose your own adventure.
 - Focus on solving the easiest problems first and then move to the harder ones.

Topics you'll need to know

- A* and consistent/admissible heuristics
- Alpha-Beta pruning for min-max search
- Expectimax search
- Probability
 - conditional prob, (cond.) independence, Bayes' rule, chain rule
- MDPs
 - Value Iteration
 - Policy Iteration (iterative version, not the closed form solution)
 - Temporal difference learning

Topics you'll need to know

- Q-Learning
- Linear value function approximation
- Policy Gradients
 - Be able to follow math for policy gradient derivation in slides.
 - I won't ask you to rederive full policy gradient
- AlphaGo
 - Understand high-level pieces and how they connect

Search Problems

A search problem consists of:

A state space



 A successor function (with actions, costs)



- A start state and a goal test
- A solution is a sequence of actions (a plan) which transforms the start state to a goal state

Graph Search Pseudo-Code

```
function GRAPH-SEARCH(problem, fringe) return a solution, or failure
    closed \leftarrow an empty set
    fringe \leftarrow \text{INSERT}(\text{MAKE-NODE}(\text{INITIAL-STATE}[problem]), fringe)
    loop do
       if fringe is empty then return failure
       node \leftarrow \text{REMOVE-FRONT}(fringe)
       if GOAL-TEST(problem, STATE[node]) then return node
       if STATE[node] is not in closed then
           add STATE[node] to closed
           for child-node in EXPAND(STATE[node], problem) do
               if STATE[child-node] is not in closed then fringe \leftarrow \text{INSERT}(child-node, fringe)
           \mathbf{end}
```

 \mathbf{end}

A-star: Combining UCS and Greedy

- Uniform-cost orders by path cost, or backward cost g(n)
- Greedy orders by goal proximity, or *forward cost* h(n)



A* Search orders by the sum: f(n) = g(n) + h(n)

Example: Teg Grenager

Admissible Heuristics

A heuristic h is admissible (optimistic) if:

 $0 \leq h(n) \leq h^*(n)$

where $h^*(n)$ is the true cost to a nearest goal

Examples:





 Coming up with admissible heuristics is most of what's involved in using A* in practice.

Consistency of Heuristics



- Main idea: estimated heuristic costs ≤ actual costs
 - Admissibility: heuristic cost ≤ actual cost to goal

$h(A) \leq actual cost from A to G$

- Consistency: heuristic "arc" cost ≤ actual cost for each arc
 h(A) h(C) ≤ cost(A to C)
- Consequences of consistency:
 - The f value along a path never decreases

 $h(A) \leq cost(A to C) + h(C)$

A* graph search is optimal

Adversarial Search

Minimax Values

C



Minimax Implementation



def min-value(state):

$$V(s') = \min_{s \in \text{successors}(s')} V(s)$$

$$V(s) = \max_{s' \in \text{successors}(s)} V(s')$$

Minimax Implementation (Dispatch)

def value(state):

if the state is a terminal state: return the state's utility if the next agent is MAX: return max-value(state) if the next agent is MIN: return min-value(state)





def min-value(state):
 initialize v = +∞
 for each successor of state:
 v = min(v, value(successor))
 return v

Minimax Example



Alpha-Beta Implementation

 α : MAX's best option on path to root β : MIN's best option on path to root

At root you should initialize $\alpha = -\infty$ and $\beta = +\infty$

```
\begin{array}{l} \mbox{def max-value(state, } \alpha, \beta): \\ \mbox{initialize } v = -\infty \\ \mbox{for each successor of state:} \\ v = max(v, value(successor, \alpha, \beta)) \\ \mbox{if } v \geq \beta \mbox{ return } v \\ \alpha = max(\alpha, v) \\ \mbox{return } v \end{array}
```

 $\begin{array}{l} \mbox{def min-value(state , \alpha, \beta):} \\ \mbox{initialize } v = +\infty \\ \mbox{for each successor of state:} \\ v = min(v, value(successor, \alpha, \beta)) \\ \mbox{if } v \leq \alpha \ return \ v \\ \beta = min(\beta, v) \\ \ return \ v \end{array}$

Alpha-Beta Quiz

 α : MAX's best option on path to root β : MIN's best option on path to root

def max-value(state, α , β): initialize $v = -\infty$ for each successor of state: $v = max(v, value(successor, \alpha, \beta))$ if $v \ge \beta$ return v $\alpha = \max(\alpha, v)$ return v

def min-value(state , α , β): initialize $v = +\infty$ for each successor of state: $v = min(v, value(successor, \alpha, \beta))$ if $v \leq \alpha$ return $v = \sqrt{-1}$ $\beta = \min(\beta, v)$ return v



Alpha-Beta Example 2



Uncertain Search

Expectimax Pseudocode

def value(state):

if the state is a terminal state: return the state's utility if the next agent is MAX: return max-value(state) if the next agent is EXP: return exp-value(state)

```
def max-value(state):
    initialize v = -∞
    for each successor of state:
        v = max(v, value(successor))
    return v
```

def exp-value(state):
 initialize v = 0
 for each successor of state:
 p = probability(successor)
 v += p * value(successor)
 return v

Expectimax Pseudocode





v = (1/2) (8) + (1/3) (24) + (1/6) (-12) = 10

Mixed Layer Types

- E.g. Backgammon
- Expectiminimax
 - Environment is an extra "random agent" player that moves after each min/max agent
 - Each node computes the appropriate combination of its children



Probability

Probability Distributions

Unobserved random variables have distributions





- A distribution is a TABLE of probabilities of values
- A probability (lower case value) is a single number

P(W = rain) = 0.1

• Must have: $\forall x \ P(X = x) \ge 0$ a

and
$$\sum_{x} P(X = x)$$

Shorthand notation: P(hot) = P(T = hot), P(cold) = P(T = cold), P(rain) = P(W = rain),....

OK *if* all domain entries are unique

= 1

Joint Distributions

• A *joint distribution* over a set of random variables: $X_1, X_2, \ldots X_n$ specifies a real number for each assignment (or *outcome*):

$$P(X_1 = x_1, X_2 = x_2, \dots, X_n = x_n)$$

 $P(x_1, x_2, \dots, x_n)$

• Must obey: $P(x_1, x_2, \dots x_n) \ge 0$

$$\sum_{(x_1, x_2, \dots, x_n)} P(x_1, x_2, \dots, x_n) = 1$$

\mathcal{P}	(T	7	W)
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Т	W	Р
hot	sun	0.4
hot	rain	0.1
cold	sun	0.2
cold	rain	0.3

- Size of distribution if n variables with domain sizes d?
 - For all but the smallest distributions, impractical to write out!

Quiz: Events

P(+x, +y) ?

•
$$P(+x)$$
? = $\sum_{y} P(x = +x, Y = y)$

P(-y OR +x) ?

P(X,Y)

Х	Υ	Р
+x	+y	0.2
+x	-у	0.3
-X	+у	0.4
-X	-у	0.1

Marginal Distributions

- Marginal distributions are sub-tables which eliminate variables
- Marginalization (summing out): Combine collapsed rows by adding



$$P(X_1 = x_1) = \sum_{x_2} P(X_1 = x_1, X_2 = x_2)$$



Quiz: Marginal Distributions



Conditional Probabilities

- A simple relation between joint and conditional probabilities
 - In fact, this is taken as the *definition* of a conditional probability

$$P(a|b) = \frac{P(a,b)}{P(b)}$$

$$P(T,W)$$

$$T W P$$

$$hot sun 0.4$$

$$hot rain 0.1$$

$$cold sun 0.2$$

$$cold rain 0.3$$



$$P(W = s | T = c) = \frac{P(W = s, T = c)}{P(T = c)} = \frac{0.2}{0.5} = 0.4$$
$$= P(W = s, T = c) + P(W = r, T = c)$$
$$= 0.2 + 0.3 = 0.5$$

Quiz: Conditional Probabilities

P(+x | +y) ?



X	Y	Р
+x	+y	0.2
+x	-у	0.3
-X	+у	0.4
-X	-у	0.1

P(-x | +y) ?

P(-y | +x) ?

The Product Rule

Sometimes have conditional distributions but want the joint

$$P(y)P(x|y) = P(x,y)$$
 $(x|y) = \frac{P(x,y)}{P(y)}$

 \mathbf{D}



The Product Rule

$$P(y)P(x|y) = P(x,y)$$

• Example:

P(W)

Ρ

0.8

0.2

R

sun

rain

<i>P</i> ((D W))	_
D	W	Р	
wet	sun	0.1	
dry	sun	0.9	
wet	rain	0.7	
dry	rain	0.3	

P	(T)	W)
1	$\langle \mathcal{L} \rangle$	' '	* *)

D	W	Р
wet	sun	
dry	sun	
wet	rain	
dry	rain	

The Chain Rule

- More generally, can always write any joint distribution as an incremental product of conditional distributions $P(x_1, x_2, x_3) = P(x_1)P(x_2|x_1)P(x_3|x_1, x_2)$ $P(x_1, x_2, \dots, x_n) = \prod_i P(x_i|x_1 \dots x_{i-1})$ $P(x_1, x_2, \dots, x_n) = \prod_i P(x_i|x_1 \dots x_{i-1})$
- You can pick any order.
- Why is the Chain Rule always true?

Bayes' Rule

That's my rule!

Two ways to factor a joint distribution over two variables:

$$P(x,y) = P(x|y)P(y) = P(y|x)P(x)$$

Dividing, we get:

$$P(x|y) = \frac{P(y|x)}{P(y)}P(x)$$

- Why is this at all helpful?
 - Lets us build one conditional from its reverse
 - Often one conditional is tricky but the other one is simple
 - Foundation of many systems (e.g. ASR, MT, IRL)
- In the running for most important AI equation!

Independence

 $X \perp \!\!\!\perp Y$

Two variables are *independent* in a joint distribution if:

P(X,Y) = P(X)P(Y) $\forall x, y P(x,y) = P(x)P(y)$

- Says the joint distribution *factors* into a product of two simple ones
- Usually variables aren't independent!
- Can use independence as a *modeling assumption*
 - Independence can be a simplifying assumption
 - *Empirical* joint distributions: at best "close" to independent
 - What could we assume for {Weather, Traffic, Cavity}?
- Independence is like something from CSPs: what?



Example: Independence?



0.4

rain

Conditional Independence

- Unconditional (absolute) independence very rare (why?)
- Conditional independence is our most basic and robust form of knowledge about uncertain environments.
- X is conditionally independent of Y given Z

 $X \perp \!\!\!\perp Y | Z$

if and only if:

 $\forall x, y, z : P(x, y|z) = P(x|z)P(y|z)$

or, equivalently, if and only if

$$\forall x, y, z : P(x|z, y) = P(x|z)$$

Probability Recap

Conditional probability

$$P(x|y) = \frac{P(x,y)}{P(y)}$$

• Product rule P(x,y) = P(x|y)P(y)

• Chain rule
$$P(X_1, X_2, \dots, X_n) = P(X_1)P(X_2|X_1)P(X_3|X_1, X_2)\dots$$

 $= \prod_{i=1}^n P(X_i|X_1, \dots, X_{i-1})$

- X, Y independent if and only if: $\forall x, y : P(x, y) = P(x)P(y)$
- X and Y are conditionally independent given Z if and only if: $X \perp\!\!\!\perp Y | Z$ $\forall x, y, z : P(x, y | z) = P(x | z) P(y | z)$

Markov Decision Processes

- An MDP is defined by:
 - A set of states s ∈ S
 - A set of actions $a \in A$
 - A transition function T(s, a, s')
 - Probability that a from s leads to s', i.e., P(s' | s, a)
 - Also called the model or the dynamics
 - A reward function R(s, a, s')
 - Sometimes just R(s) or R(s')
 - A start state
 - Maybe a terminal state
 - · dissount factor y
- MDPs are non-deterministic search problems
 - One way to solve them is with expectimax search
 - We'll have a new tool soon



[Demo – gridworld manual intro (L8D1)]

What is Markov about MDPs?

- "Markov" generally means that given the present state, the future and the past are independent
- For Markov decision processes, "Markov" means action outcomes depend only on the current state

$$P(S_{t+1} = s' | S_t = s_t, A_t = a_t, S_{t-1} = s_{t-1}, A_{t-1}, \dots, S_0 = s_0)$$

= $P(S_{t+1} = s' | S_t = s_t, A_t = a_t)$



Andrey Markov (1856-1922)

 This is just like search, where the successor function could only depend on the current state (not the history)

Important Quantities

- The value (utility) of a state s:
 - V^{*}(s) = expected utility starting in s and acting optimally
- The value (utility) of a q-state (s,a):
 - Q^{*}(s,a) = expected utility starting out having taken action a from state s and (thereafter) acting optimally
- The optimal policy:
 π^{*}(s) = optimal action from state s



Bellman Equations

- Fundamental operation: compute the (expectimax) value of a state
 - Expected utility under optimal action
 - Average sum of (discounted) rewards
 - This is just what expectimax computed!
- Recursive definition of value:

$$V^{*}(s) = \max_{a} Q^{*}(s, a)$$
$$Q^{*}(s, a) = \sum_{s'} T(s, a, s') \left[R(s, a, s') + \gamma V^{*}(s') \right]$$
$$V^{*}(s) = \max_{a} \sum_{s'} T(s, a, s') \left[R(s, a, s') + \gamma V^{*}(s') \right]$$



Value Iteration

- Start with V₀(s) = 0: no time steps left means an expected reward sum of zero
- Given vector of V_k(s) values, do one ply of expectimax from each state:

$$V_{k+1}(s) \leftarrow \max_{a} \sum_{s'} T(s, a, s') \left[R(s, a, s') + \gamma V_k(s') \right]$$

Bellman Update Equation

- Repeat until convergence
- Complexity of each iteration: O(S²A)
- Theorem: will converge to unique optimal values
 - Basic idea: approximations get refined towards optimal values
 - Policy may converge long before values do



Policy Iteration

- Alternative approach for optimal values:
 - Step 1: Policy evaluation: calculate utilities for some fixed policy (not optimal utilities!) until convergence
 - Step 2: Policy improvement: update policy using one-step look-ahead with resulting converged (but not optimal!) utilities as future values
 - Repeat steps until policy converges
- This is policy iteration
 - It's still optimal!
 - Can converge (much) faster under some conditions

Policy Iteration

- Evaluation: For fixed current policy π , find values with policy evaluation:
 - Iterate until values converge:

$$V_{k+1}^{\pi_i}(s) \leftarrow \sum_{s'} T(s, \pi_i(s), s') \left[R(s, \pi_i(s), s') + \gamma V_k^{\pi_i}(s') \right]$$

- Improvement: For fixed values, get a better policy using policy extraction
 - One-step look-ahead:

$$\pi_{i+1}(s) = \arg\max_{a} \sum_{s'} T(s, a, s') \left[R(s, a, s') + \gamma V^{\pi_i}(s') \right]$$

Temporal Difference Learning

- Big idea: learn from every experience!
 - Update V(s) each time we experience a transition (s, a, s', r)
 - Likely outcomes s' will contribute updates more often
- - Move values toward value of whatever successor occurs: running average

Sample of V(s):
$$sample = R(s, \pi(s), s') + \gamma V^{\pi}(s')$$
Update to V(s): $V^{\pi}(s) \leftarrow (1 - \alpha)V^{\pi}(s) + (\alpha)sample$ Same update: $V^{\pi}(s) \leftarrow V^{\pi}(s) + \alpha(sample - V^{\pi}(s))$

Q-Learning

(So aoro S, A, r, Sz. ... uptatel uptatel

 $Q(\varsigma_{o_1}a_{o_2})$

Q-Learning: sample-based Q-value iteration

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

- Learn Q(s,a) values as you go
 - Receive a sample (s,a,s',r)
 - Consider your old estimate: Q(s, a)
 - Consider your new sample estimate:

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

Incorporate the new estimate into a running average:

$$Q(s,a) \leftarrow (1-\alpha)Q(s,a) + (\alpha) [sample]$$

$$Q(S,a) \leftarrow Q(S,a) + \ll (Sample - Q(S,a))$$



Linear Value Functions

 Using a feature representation, we can write a q function (or value function) for any state using a few weights:

$$V(s) = w_1 f_1(s) + w_2 f_2(s) + \ldots + w_n f_n(s)$$

$$Q(s,a) = w_1 f_1(s,a) + w_2 f_2(s,a) + \ldots + w_n f_n(s,a)$$

- Advantage: our experience is summed up in a few powerful numbers
- Disadvantage: states may share features but actually be very different in value!

Approximate Q-Learning

$$Q(s,a) = w_1 f_1(s,a) + w_2 f_2(s,a) + \ldots + w_n f_n(s,a)$$

Q-learning with linear Q-functions:

transition =
$$(s, a, r, s')$$

difference = $\left[r + \gamma \max_{a'} Q(s', a')\right] - Q(s, a)$
 $Q(s, a) \leftarrow Q(s, a) + \alpha$ [difference] Exact
 $w_i \leftarrow w_i + \alpha$ [difference] $f_i(s, a)$ Approx

Approximate Q's

- Intuitive interpretation:
 - Adjust weights of active features
 - E.g., if something unexpectedly bad happens, blame the features that were on: disprefer all states with that state's features
- Formal justification: online least squares



DQN

- Approximate Q-Learning at scale.
- Uses Neural Network for Q-value function approximation.

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Two approaches to model-free RL

Learn Q-values

- Trains Q-values to be consistent. Not directly optimizing for performance.
- Use an objective based on the Bellman Equation

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

- Learn Policy Directly
 - Have a parameterized policy π_{θ}
 - Update the parameters θ to optimize performance of policy.

Policy Gradient RL

Find a policy that maximizes expected utility (discounted cumulative rewards)

$$\pi^* = \arg\max_{\pi} E_{\pi} \left[\sum_{t=0}^{\infty} \gamma^t R(s, \pi(s), s') \right]$$

Notation

- Trajectory (rollout, episode) $\tau = (s_0, a_0, s_1, a_1, ...)$
 - $s_0 \sim \rho_0(\cdot)$ (initial state distribution)
 - $s_{t+1} \sim P(\cdot | s_t, a_t)$ (transition probabilities)
- Rewards $r_t = R(s_t, a_t, s_{t+1})$
- Finite-horizon undiscounted return of a trajectory

$$R(\tau) = \sum_{t=0}^{T} r_t$$

• Actions are sampled from a stochastic parameterized policy π_{θ} $a_t \sim \pi_{\theta}(\cdot | s_t)$

Notation

- Probability of a trajectory (rollout, episode) $\tau = (s_0, a_0, s_1, a_1, ...)$ $P(\tau|\pi) = \rho_0(s_0) \prod_{t=0}^{T-1} P(s_{t+1}|s_t, a_t) \pi_{\theta}(a_t|s_t)$
- Expected Return of a policy $J(\pi)$

$$J(\pi) = \sum_{\tau} P(\tau|\pi) R(\tau) = E_{\tau \sim \pi}[R(\tau)]$$

Goal of RL: Solve the following optimization problem

$$\pi^* = \operatorname*{argmax}_{\pi} J(\pi)$$

The Policy Gradient

We can now perform gradient ascent to improve our policy!

$$\theta_{k+1} \leftarrow \theta_k + \alpha \nabla_{\theta} J(\pi_{\theta}) \Big|_{\theta_k}$$

$$\nabla_{\theta} J(\pi_{\theta}) = E_{\tau \sim \pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_t | s_t) R(\tau) \right]$$

Estimate with a sample mean over a set D of policy rollouts given current parameters

$$\approx \frac{1}{|D|} \sum_{\tau \in D} \sum_{t=0}^{T} (\nabla_{\theta} \log \pi_{\theta}(a_t | s_t) R(\tau))$$

Alpha Go



There will be one short answer question about AlphaGo. Review high-level ideas from slides. Don't worry about nitty-gritty details.