CS50: Introduction to Artificial Inteligence Notes

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# Abstract CS50: Introduction to Artificial Inteligence notes. Note template by Pingbang Hu.

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# Chapter 1

# Knowledge

# 1.1 Propositional Logic

We use standard logic notation:

- ¬p
- $p \vee q$
- p ∧ q
- $p \Rightarrow q$ :

r · T				
p	q	$p \Rightarrow q$		
false	false	true		
false	true	true		
true	false	$_{ m false}$		
true	true	true		

p ⇔ q:

p	q	$p \Leftrightarrow q$
false	false	true
false	true	true
${ m true}$	false	$_{ m false}$
$\operatorname{true}$	true	true

Now we must establish *what* is considered to be "true" in our world by defining a **model**. We need to represent that knowledge. We do so by defining it via a **knowledge base**.

**Definition 1.1.1** (Model). Assignment of a truth value to every propositional symbol.

Definition 1.1.2 (Knowledge Base). A set of sentences known by a knowledge-based agent.

Definition 1.1.3 (Enatilment).

$$\alpha \models \beta$$
 " $\alpha$  entails  $\beta$ "

In every model in which sentence  $\alpha$  is true, sentence  $\beta$  is also true.

# 1.2 Inference

Our aim is to see if our knowledge base, KB, entails some query about the world,  $\alpha$ :

$$KB \models \alpha$$
?

We first define a **model checking algorithm** to determine if  $KB \models \alpha$ . We can determine this by doing the following:

- enumerate all possible models
- if in every model where KB is true,  $\alpha$  is also true, then  $KB \models \alpha$

# 1.3 Inference By Resolution

To determine if  $KB \models \alpha$  via knowledge resolution:

- Check if  $KB \wedge \neg \alpha$  is a contradiction
  - Convert  $KB \wedge \neg \alpha$  to Conjunctive Normal Form
  - Keep checking to see if we can use resolution to produce new clause
  - If we ever produce the empty clause (equivalent to False), we have a contradiction and so  $KB \models \alpha$
- If so, then  $KB \models \alpha$
- Otherwise, no entailment

# **Problem 1.3.1.** Does $(A \vee B) \wedge (\neg B \vee C) \wedge (\neg C)$ entail A?

**Answer.** First, we convert to CNF:

$$(A \lor B) \land (\neg B \lor C) \land (\neg C) \land (\neg A)$$

We can resolve  $(\neg B \lor C)$  and  $(\neg C)$  by concluding that  $\neg B$ . With the knowledge of  $\neg B$  we now see that, considering  $A \lor B$ , we can conclude A. We see that

$$A \land \neg A \Rightarrow False$$

and so we can conclude that the clause entails A.

# Chapter 2

# Probability

# 2.1 Introduction to Probability

We can represent a **possible world** using  $\omega$  where all possible worlds is the set  $\Omega$ . Thus, we can define total probability as:

$$\sum_{\omega\in\Omega}p(\omega).$$

We build machine models to predict outcomes based on data using **conditional probability**. In other words, we have probability based on some evidence. We can calculate conditional probability using the following formula:

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

# 2.2 Random Variables

In probability we define a variable with a set of possible values as a random variable:

**Definition 2.2.1** (Random Variable). A variable possessing a distribution of probabilities for various "states".

# 2.3 Baye's Rule

From the above equation we have:

$$P(a \wedge b) = P(b)P(a|b) = P(a)P(b|a)$$

**Definition 2.3.1** (Independence). The knowledge that one event occurs does not affect the probability of the other event. So we have

$$P(a \wedge b) = P(a)P(b)$$

since P(b) = P(b|a) if a and b are independent.

Thus, we can derive Baye's rule, which relates the probability of one event on the condition of another event, to the reverse relationship.

Definition 2.3.2 (Baye's Rule).

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

What Baye's rule allows us to do is that given:

 $P(\text{visible effect} \mid \text{unknown cause})$ 

# 2.4 Bayesian Networks

There are a number of different probabilistic models. The first we discuss are **Bayesian Networks**.

**Definition 2.4.1** (Bayesian Network). A data structure that represents the dependencies among random variables. They have the following characteristics:

- directed graph
- each node represents a random variable
- arrow from X to Y means X is a parent of Y
- each node X has probability distribution P(X|Parents(X))

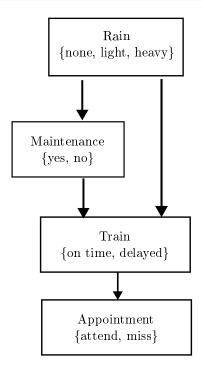


Figure 2.1: A basic example of a Bayesian Network.

We now aim to make an inference using the Bayesian Network. Given the following:

- Query X: variable for which to compute distribution
- $\bullet$  Evidence variables E: observed variables for event e
- $\bullet$  Hidden variables Y: non-evidence, non-query variable

our goal is to calculate P(X|e).

### **Problem 2.4.1.** Calculate P(Appointment|light, no)

**Answer**. Here, the *evidence* is that there is light rain and no maintenance. The *query* is the status of Appointment. The *hidden layer* is the status of the train, since you are not given the train's status and you are not querying it; it's just a confounding variable.

We note that

$$P(Appointment|light, no) = \alpha P(Appointment, light, no)$$

and by the marginalization technique:

 $= \alpha [P(\text{Appointment, light, no, on time}) + P(\text{Appointment, light, no, delayed})].$ 

Remark. The marginalization technique essentially states that

$$\alpha P(Q, E) = \alpha \left[ P(Q, E, Y) + P(Q, E, \neg Y) \right]$$

# 2.5 Inference by Enumeration

The above is an example of **inference by enumeration**. More formally, for the following:

- X: the query variable
- e: the evidence
- y: ranges over values of hidden variables
- $\alpha$ : normalizes the result

we have:

$$P(X|e) = \alpha P(X,e) = \alpha \sum_{y} P(X,e,y)$$

# 2.6 Sampling

Rather than attempting to calculate an exact probability, we can approximate a probability instead via **sampling**. By randomly generating samples for n = 1000 or n = 10000, we can get fairly useful results.

**Definition 2.6.1** (Rejection Sampling). **Rejection sampling** is the process of simulating numerous examples from a distribution while considering only the samples that possess the attributes of the desired query.

**Remark.** Rejection sampling is not particularly effective when the evidence you are looking for is fairly unlikely, since you are rejecting a lot of samples. This is inefficient since you are throwing away a large portion of your samples.

**Definition 2.6.2** (Likelihood Weighting). Rather than sampling everything, in **likelihood weighting**, we start by *fixing* the values for evidence variables. We then sample the non-evidence variables using conditional probabilities in the Bayesian Network. Finally we weight each sample by its *likelihood*.

### 2.7 Markov Models

As opposed to assigning one random variable to a value, but rather an array of random variables to a value over a timescale. This leads to a lot more data, and so we must make some assumptions.

**Definition 2.7.1** (Markov Assumption). The assumption that the current state depends only on a finite fixed number of previous states.

\*

**Definition 2.7.2** (Markov Chain). A sequence of random variables where the distribution of each variable follows the Markov assumption.

For example, consider the basic Markov Chain

$$X_t \to X_{t+1}$$

where  $X_t$  represents the weather today, and  $X_{t+1}$  represents the weather tomorrow. We can construct a **transition model** to describe the relationship as follows:

	$X_{t+1}$ : Sunny	$X_{t+1}$ : Rainy
$X_t$ : Sunny	0.8	0.2
$X_t$ : Rainy	0.3	0.7

The first line, for example, translates to "given today was sunny, tomorrow will be sunny with a probability of 0.8 and rainy with probability 0.2". We can refine our definition of the Markov Model further, by introducing *hidden* states.

 $\textbf{Definition 2.7.3} \ (\textbf{Hidden Markov Model}). \ A \ Markov \ model \ for \ a \ system \ with \ hidden \ states \ that \ generate some observed \ event.$ 

**Definition 2.7.4** (Filtering). Given observations from start until now, calculate distribution for current state.

**Definition 2.7.5** (Prediction). Given observations from start until now, calculate the distribution for a future state.

**Definition 2.7.6** (Smoothing). Given observations from start until now, calculate the distribution for a past state.

**Definition 2.7.7** (Most Likely Explanation). Given observations from start until now, calculate most likely sequence of states.

# Chapter 3

# Optimization

# 3.1 Hill Climbing

The most basic form of optimization we explore is that of the local search.

**Definition 3.1.1** (Local Search). Search algorithms that maintain a single node and searches by moving to a neighboring node.

We can implement a local searching using hill climbing in which we do the following:

- Check values of neighbors of current best value
- If a neighbor has a value closer to the desired value, set that value as current best

# Algorithm 3.1: HillClimb

The problem with this naive approach (**steepest ascent**) is that we may get stuck at a local extremum rather than the global extremum. Further, we may encounter a "shoulder", where several neighboring values are the same, and get stuck.

# 3.2 Simulated Annealing

Although there are a variety of hill climbing algorithms (stochastic, first-choice, beam, etc.). They have their flaws but the theme is that we never go from a good value to a worse value. Thus, we must overcome local extrema. We can tackle this with **simulated annealing**.

**Definition 3.2.1** (Simulated Annealing). Simulated annealing is akin to a "cooling" physical process.

- Early on, higher "temperature": more likely to accept neighbors that are worse than current state
- Later on, lower "temperature": less likely to accept neighbors that are worse than current state

## Algorithm 3.2: Simulated Annealing

```
Data: problem, max

1 cur = initial state of problem

2 for t = 1 to max do

3 | T = \text{Temperature}(t)

4 | neighbor = random neighbor of cur

5 | \Delta E = \text{how much better neighbor is than cur}

6 | if \Delta E > 0 then

7 | cur = neighbor

8 | with probability e^{\Delta E/T}, set cur = neighbor
```

# 3.3 Linear Programming

In the context where we are trying to optimize for some function or when we have real values we are often trying to minimize or maximize a cost function given a variety of constraints. This is where **linear programming** becomes useful.

Definition 3.3.1 (Linear Programming). Linear programming problems often entail the following:

- Minimize a cost function  $c_1x_1 + c_2x_2 + \ldots + c_nx_n$
- With constraints of form  $a_1x_1 + a_2x_2 + \ldots + a_nx_n \leq b$  or of form  $a_1x_1 + a_2x_2 + \ldots + a_nx_n = b$
- With bounds for each variable  $l_i \leq x_i \leq u_i$

**Problem 3.3.1.** Two machines  $X_1$  and  $X_2$  which cost \$50/hr and \$80/hr to run, respectively. We have the following constraints:

- $X_1$  requires 5 units of labor,  $X_2$  requires 2 units of labor per hour; we have total of 20 units of labor to spend; goal is to minimize cost
- $X_1$  produces 10 units of output per hour,  $X_2$  produces 12 units of output per hour; company needs 90 units of output

**Answer.** We can create a cost function:

$$50x_1 + 80x_2$$

and our constraints:

$$5x_1 + 2x_2 \le 20$$

$$10x_1 + 12x_2 \ge 90$$

or equivalently:

$$-10x_1 + -12x_2 \le -90$$

and then solve using a standard linear programming technique such as Simplex, Interior-Point, etc.

# 3.4 Constraint Satisfaction

Constraint satisfaction problems often have some number of variables that must be optimized, but they are subject to some constraints.

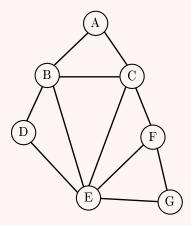
**Definition 3.4.1** (Constraint Satisfaction Problem). A general constraint satisfaction problem consists of the following:

- Set of variables  $\{x_1,\ldots,x_n\}$
- Set of domains for each variable  $\{D_1, \ldots, D_n\}$
- Set of constraints C

**Problem 3.4.1.** We must schedule exams for classes  $A, B, \ldots, G$  such that no conflicts arise for the students taking the courses. The four students' schedules are:

- Student 1: A, B, C
- Student 2: B, D, E
- Student 3: C, E, F
- Student 4: E, F, G

We can represent this graphically using an undirected graph, where the nodes represent exams and the edges indicate that the two exams *cannot* be scheduled for the same time.



Here we define our variables, domains, and constraints:

- Variables:  $\{A, B, C, D, E, F, G\}$
- Domains:  $\{Monday, Tuesday, Wednesday\}$
- Constraints:  $\{A \neq B, A \neq C, B \neq C, \dots, E \neq G, F \neq G\}$

Our goal is to find an assignment of a day to each of the classes such that we don't have any conflicts between the classes. In other words, we are aiming for **node consistency**.

**Definition 3.4.2** (Node Consistency). When all the values in a variable's domain satisfy the variable's unary constraints.

Furthermore, we are seeking arc consistency.

**Definition 3.4.3.** When all the values in a variable's domain satisfy the variable's binary constraints. More formally: to make X arc-consistent with respect to Y, remove elements from X's domain until every choice for X has a possible choice for Y.

**Definition 3.4.4** (Unary Contraint). A constraint involving a *single* variable. (e.g.  $\{A \neq Wednesday\}$ )

**Definition 3.4.5** (Binary Constraint). A constraint involving two variables. (e.g. $\{A \neq B\}$ )

Note that constraints can come of different forms. **Hard** constraints are absolute requirements, whereas **soft** constraints are preferences.

# 3.5 Arc Consistency

We first aim to define a function that, given some consistency problem csp, can make some variable X arc consistent with respect to another variable, Y.

### Algorithm 3.3: Revise

We can enforce arc consistency across an entire consistency problem:

### Algorithm 3.4: AC-3

### 3.5.1 CSPs as Search Problems

We can reframe CSPs as search problems. We do so by defining a CSP as follows:

- initial state: empty assignment (no variables)
- actions: add a  $\{variable = value\}$  to assignment
- transition model: shows how adding an assignment changes the assignment
- goal test: check if all variables assigned and constraints all satisfied
- path cost function: all paths have same cost

# 3.6 Backtracking

We can find a solution to a CSP by simply applying arbitrary assignments to variables one by one until a constraint is broken. Then, we can just backtrack and try another assignment. We eventually find a solution or check every possible assignment.

However, we can be more clever in our approach and apply the idea of **inference** to our approach. We can operate just as before, but when we find ourselves about to backtrack we can instead observe the graph and look for *arc inconsistencies*. Using the information from neighboring nodes, we can deduce what a particular node's possible assignments may be. Thus, we can backtrack less than we originally did.

### Algorithm 3.5: Backtrack

```
Data: assignment, csp
 1 if assignmennt complete then
   return assignment
3 \text{ var} = \text{SelectUnassignedVar}(assignment, csp)
 4 for value in DomainValues(var, assignment, csp) do
      if value consistent with assignment then
          add \{var = value\} to assignment
          inferences = Inference(assignment, csp) // get inferences
          if inferences \neq failure then
 8
          add inferences to assignment // add inferenced nodes
 9
          result = Backtrack(assignment, csp)
10
         if result \neq failure then
11
             return result
     remove \{var = value\} and inferences from assignment
14 return failure
```

We note that we can make the search process more efficient by being smarter about which variable we select in our SelectUnassignedVar function. We use various heuristics:

- Minimum Remaining Values (MRV) heuristic: select the variable that has the smallest domain
- Degree heuristic: select the variable that has the highest degree; works because the highest degree variable has the most constraints, removing it is most helpful

Additionally, we can refine our DomainValues function further. We utilize another set of heuristics:

• Least-constraining values heuristic: return variables in order by number of choices that are ruled out for neighboring variables; try least-constraining values first

# Appendix