Knowledge representation and reasoning

- First-order logic: syntax and semantics
- Soundness and completeness of proof procedures
- Converting first-order formulas into clausal form
- Unification and MGU
- Resolution proof: forward chaining and refutation
- Answer extraction
- Knowledge graph

Alphabet

Logical symbols (fixed meaning and use):

- Punctuation: (,),,,.
- Connectives and quantifiers: $=, \neg, \land, \lor, \forall, \exists$
- Variables: $x, x_1, x_2, ..., x', x'', ..., y, ..., z, ...$

Non-logical symbols (domain-dependent meaning and use):

- Predicate (谓词) symbols
 - arity: number of arguments
 - arity 0 predicates: propositional symbols
- Function symbols
 - arity 0 functions: constant symbols



Terms (项)

- Every variable is a term
- If t_1, \ldots, t_n are terms and f is a function symbol of arity n, then $f(t_1, \ldots, t_n)$ is a term

Formulas

- If t_1, \ldots, t_n are terms and P is a predicate symbol of arity n, then $P(t_1, \ldots, t_n)$ is an atomic formula
- ullet If t_1 and t_2 are terms, then $(t_1=t_2)$ is an atomic formula
- If α and β are formulas, and v is a variable, then $\neg \alpha, (\alpha \land \beta), (\alpha \lor \beta), \exists v.\alpha, \forall v.\alpha$ are formulas

Interpretations

An interpretation (\mathbb{R}) is a pair $M = \langle D, I \rangle$

- D is the domain, can be any non-empty set
- ullet I is a mapping from the set of predicate and function symbols
- If P is a predicate symbol of arity n, I(P) is an n-ary relation over D, i.e., $I(P) \subseteq D^n$
 - If p is a 0-ary predicate symbol, i.e., a propositional symbol, $I(p) \in \{true, false\}$
- If f is a function symbol of arity n, I(f) is an n-ary function over D, i.e., $I(f):D^n\to D$
 - If c is a 0-ary function symbol, i.e., a constant symbol, $I(c) \in D$



Denotation (指称) of terms

- Terms denote elements of the domain
- \bullet A variable assignment μ is a mapping from the set of variables to the domain D
- $\bullet \|v\|_{M,\mu} = \mu(v)$
- $||f(t_1,\ldots,t_n)||_{M,\mu} = I(f)(||t_1||_{M,\mu},\ldots,||t_n||_{M,\mu})$

Satisfaction: atomic formulas

 $M, \mu \models \alpha$ is read " M, μ satisfies (满足) α "

- $M, \mu \models P(t_1, \dots, t_n) \text{ iff } \langle ||t_1||_{M,\mu}, \dots, ||t_n||_{M,\mu} \rangle \in I(P)$
- $M, \mu \models (t_1 = t_2) \text{ iff } ||t_1||_{M,\mu} = ||t_2||_{M,\mu}$

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Intro to Al

7/6

Satisfaction: propositional connectives

- $M, \mu \models \neg \alpha \text{ iff } M, \mu \not\models \alpha$
- $\bullet \ M, \mu \models (\alpha \land \beta) \ \text{iff} \ M, \mu \models \alpha \ \text{and} \ M, \mu \models \beta$
- $M, \mu \models (\alpha \lor \beta)$ iff $M, \mu \models \alpha$ or $M, \mu \models \beta$

Satisfaction: quantifiers

 $\mu[v\mapsto d]$ denotes a variable assignment just like $\mu,$ except that it maps v to d

- $M, \mu \models \exists v. \alpha$ iff for some $d \in D$, $M, \mu[v \mapsto d] \models \alpha$
- $M, \mu \models \forall v. \alpha$ iff for all $d \in D$, $M, \mu[v \mapsto d] \models \alpha$

Let α be a sentence. Then whether $M, \mu \models \alpha$ is independent of μ .

Thus we simply write $M \models \alpha$



Logical entailment (蕴涵)

- ullet $S \models \alpha$ iff for every M, if $M \models S$ then $M \models \alpha$
- $S \models \alpha$ is read: S entails α or α is a logical consequence (推论) of S
- A special case: $\emptyset \models \alpha$, simply written $\models \alpha$, read " α is valid" (有效的)

Inference (推理) procedure

- We want a mechanical procedure to check if $KB \models \alpha$
- Called an inference procedure
- Sound (合理的) if whenever it says yes, then $KB \models \alpha$
- Complete (完备的) if whenever $KB \models \alpha$, then it says yes

Conversion to Clausal Form: 8 steps

- Eliminate Implications.
- Move Negations inwards (and simplify ¬¬).
- Standardize Variables.
- Skolemize.
- Convert to Prenex (前缀) Form.
- O Distribute disjunctions over conjunctions.
- Flatten nested conjunctions and disjunctions.
- Convert to Clauses.

Unifiers

- A unifier (合一项) of two formulas f and g is a substitution σ that makes f and g syntactically identical.
- Note that not all formulas can be unified substitutions only affect variables.
- e.g., P(f(x),a) and P(y,f(w)) cannot be unified, as there is no way of making a=f(w) with a substitution.

Computing MGUs

Given two atomic formulas f and g

- **1** $\sigma = \{\}; S = \{f, g\}$
- 2 If S contains an identical pair of formulas, stop and return σ as the MGU of f and g.
- **3** Else find the disagreement set $D = \{e_1, e_2\}$ of S
- If $e_1=V$ a variable, and $e_2=t$ a term not containing V (or vice-versa) then let $\sigma=\sigma\{V=t\}$; $S=S\{V=t\}$; Goto 2
- **5** Else stop, f and g cannot be unified.

Note: to update σ , we must compose σ with $\{V=t\}$. A common error is to just add V=t to σ .



First-order Resolution

From the two clauses $\{\rho_1\} \cup c_1$ and $\{\neg \rho_2\} \cup c_2$, where there exists a MGU σ for ρ_1 and ρ_2 , infer the clause $(c_1 \cup c_2)\sigma$

Theorem. $S \vdash ()$ iff S is unsatisfiable

Answer extraction

- We can also answer wh- questions
- Replace query $\exists x P(x)$ by $\exists x [P(x) \land \neg answer(x)]$
- Instead of deriving (), derive any clause containing just the answer predicate

什么是知识图谱

- 知识图谱是一张有向图,图中的节点表示实体或概念, 而图中的边则由属性或关系构成
- 知识图谱可以看作是三元组(triple)的集合
 - (实体1-关系-实体2): 中国-首都-北京
 - (实体-属性-属性值): 北京-人口-2069万

Search

- Problem solving by search: formalization
- Uninformed search: Breadth-First, Uniform-Cost, Depth-First, Depth-Limited, and Iterative- Deepening
- Heuristic search: Greedy best-first, A*
- Properties of search: completeness, optimality, time and space complexity
- Path/cycle checking
- Game tree search: MiniMax, alpha-beta pruning
- Simulated annealing and genetic algorithms



The formalism

To formulate a problem as a search problem we need the following components:

- Formulate a state space over which to search. The state space necessarily involves abstracting the real problem.
- Formulate actions that allow one to move between different states. The actions are abstractions of actions you could actually perform.
- Identify the initial state that best represents your current state
- Identify the goal or desired condition one wants to achieve.

A solution to the problem is a sequence of actions that can transform the initial state into a state where the goal condition holds.

Tree search

- Frontier is the set of states we haven't yet explored/expanded, and want to explore
- Initial call has Frontier = the set of initial state

```
TreeSearch(Frontier, Sucessors, Goal?)

If Frontier is empty return failure

Curr = select state from Frontier

If (Goal?(Curr)) return Curr.

Frontier' = (Frontier - {Curr}) U Successors(Curr)

return TreeSearch(Frontier', Successors, Goal?)
```

Summary of uninformed search

Criterion	Breadth- First	Uniform- Cost	Depth- First	Depth- Limited	Iterative Deepening	Bidirectional (if applicable)
Complete?	Yes^a	$Yes^{a,b}$	No	No	Yes^a	$\mathrm{Yes}^{a,d}$
Time	$O(b^d)$	$O(b^{1+\lfloor C^*/\epsilon\rfloor})$	$O(b^m)$	$O(b^{\ell})$	$O(b^d)$	$O(b^{d/2})$
Space	$O(b^d)$	$O(b^{1+\lfloor C^*/\epsilon\rfloor})$	O(bm)	$O(b\ell)$	O(bd)	$O(b^{d/2})$
Optimal?	Yes^c	Yes	No	No	Yes^c	$\mathrm{Yes}^{c,d}$

Figure 3.21 Evaluation of tree-search strategies. b is the branching factor; d is the depth of the shallowest solution; m is the maximum depth of the search tree; l is the depth limit. Superscript caveats are as follows: a complete if b is finite; b complete if step costs b for positive b optimal if step costs are all identical; b if both directions use breadth-first search.

Note. The table is from the textbook where for BFS, goal test is done when a node is generated, so both time and space complexity is $O(b^d)$ instead of $O(b^{d+1})$.

A* search: Summary

- Define an evaluation function f(n) = g(n) + h(n)
- We use f(n) to order the nodes on the frontier.
- h(n) is admissible if for all nodes n, $h(n) \leq h^*(n)$
- h(n) is consistent/monotone if for any nodes n_1 and n_2 , $h(n_1) \leq c(n1 \rightarrow n2) + h(n_2)$
- Consistency implies admissibility
- Admissibility implies optimality
- Exponential time and space complexity



Path / cycle checking

- ullet Path checking: when we expand n to obtain child c, ensures that the state c is not equal to the state reached by any ancestor of c along this path
- ullet Cycle checking: keep track of all states previously expanded during the search; when we expand n to obtain child c, ensure that c is not equal to any previously expanded state
- For uniform-cost search, cycle checking preserves optimality
- For A* with monotone heuristics, cycle checking preserves optimality

Summary: Two-Player Zero-Sum Game

- Two players A (Max) and B (Min)
- Set of states S
- An initial state $I \in S$
- ullet Terminal positions $T\subseteq S$
- Successor function
- Utility (效益) function $V: T \to \mathbf{R}$.

The MiniMax Strategy

- Assume that the other player will always play their best move
- you always play a move that will minimize the payoff that could be gained by the other player.

Alpha beta pruning

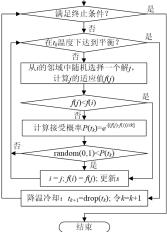
- Mark Max nodes with the change of alpha values, and Min nodes with the change of beta values
- Do alpha cut on a Max node whenever the current value

 the value of an ancestor Min node
- ullet Do beta cut on a Min node whenever the current value \leq the value of an ancestor Max node

模拟退火基本流程



初始化: 随机选择一个解i, 计算适应值f(i)。 设置代数k=0, 初始温度 $t_0=T$, 令最优解s=i。

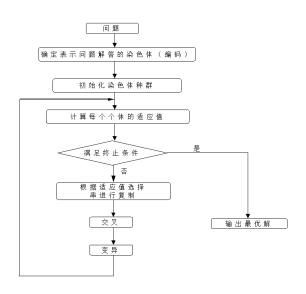


```
//功能,模拟很火管法伪代码
//说明: 本例以求问题最小值为目标
//参数: T为初始温度; L为内层循环次数
procedure SA
  //Initialization
     Randomly generate a solution X_0, and calculate its
     fitness value f(X_0);
  X_{best} = X_0; k = 0; t_k = T;
  while not stop
     //The search loop under the temperature t_k
     for i=1 to L //The loop times
       Generate a new solution X_{now} based on the current
        solution X_k, and calculate its fitness value f(X_{new}).
       if f(X_{now}) < f(X_k)
           X_k = X_{new};
           if f(X_k) < f(X_{best}) X_{best} = X_k;
           continues:
       end if
       Calculate P(t_k)=e^{-[(f(X)new)-f(Xk))/tk]};
        if random(0,1) < P
           X_k = X_{nove};
       end if
     end for
     //Drop down the temperature
     t_{k+1} = drop(t_k); k = k+1;
  end while
```

print X_{best}

end procedure

遗传算法的一般步骤



常见问题

- 旅行商问题 (Traveling Salesman Problem, TSP): 给定一 系列城市和每对城市之间的距离,求解访问每一座城市一次 并回到起始城市的最短回路
- 背包问题(Knapsack problem): 给定一组物品, 每种物品都有自己的重量和价值, 在限定的总重量内, 选择物品使得物品的总价值最高。

Planning

- Closed world assumption
- STRIPS representation of actions
- STRIPS planning
- Relaxed plan heuristics

Closed World Assumption (CWA)

- The knowledge base used to represent a state of the world is a list of positive ground atomic facts. (Like a database.)
- Closed World Assumption (CWA) is the assumption that
 - the constants mentioned in KB are all the domain objects.
 - if a ground atomic fact is not in our list of "known" facts, its negation must be true.
- This gives complete information about the state of the system.

STRIPS Actions

- STRIPS represents an action using 3 lists.
 - A list of action preconditions.
 - A list of action add effects.
 - A list of action delete effects.
- These lists contain variables, so that we can represent a whole class of actions with one specification.
- Each ground instantiation of the variables yields a specific action.

Classical planning

- Given
 - a CW-KB representing the initial state,
 - A set of STRIPS operators mapping a state to a new state
 - a goal condition
- Determine a sequence of actions that transforms the initial CW-KB to a CW-KB satisfying the goal

Relaxed problem

- We make the assumption that
 - the precondition of each action is a set of positive facts, and
 - the goal is a set of positive facts.
- Recall that we can obtain heuristics by solving a relaxed version of 8-puzzle in which we relax one of the restrictions:
 - move to adjacent field only
 - move to blank field only
- The idea here is similar: consider what happens if we ignore the delete lists of actions.
- This yields a "relaxed problem" that can produce a useful heuristic estimate.



Reasoning under uncertainty

- Bayesian networks: graphs + tables, inference
- Variable elimination algorithm
- Use D-separation to determine independence

Bayesian Networks

A BN over variables $\{X_1, X_2, \dots, X_n\}$ consists of:

- a DAG (directed acyclic graph) whose nodes are the variables
- a set of CPTs (conditional probability tables) $Pr(X_i|Par(X_i))$ for each X_i

Inference in Bayes Nets

Given

1) a Bayes net

$$Pr(X_1, X_2,..., X_n)$$

= $Pr(X_n \mid Par(X_n)) * Pr(X_{n-1} \mid Par(X_{n-1}) * \cdots * Pr(X_1 \mid Par(X_1))$

2) some Evidence, E

E = {a set of values for some of the variables}

We want to

compute the new probability <u>distribution</u>

$$Pr(X_k | E)$$

That is, we want to figure out

$$Pr(X_k = d \mid E)$$
 for all $d \in Dom[X_k]$



The VE Algorithm

Given a Bayes Net with CPTs F, query variable Q, evidence variables $\bf E$ (observed to have values e), and remaining variables $\bf Z$. Compute $Pr(Q|\bf E)$

- Replace each factor $f \in F$ that mentions a variable(s) in **E** with its restriction $f_{\mathbf{E}=e}$ (this might yield a "constant" factor)
- ② For each Z_j in the order given eliminate $Z_j \in \mathbf{Z}$ as follows:
 - ① Let f_1, f_2, \ldots, f_k be the factors in F that include Z_j
 - 2 Compute new factor $g_j = \sum_{Z_j} f_1 \times f_2 \times \ldots \times f_k$
 - f 3 Remove the factors f_i from ${\sf F}$ and add new factor g_j to ${\sf F}$
- **3** The remaining factors refer only to the query variable Q. Take their product and normalize to produce $Pr(Q|\mathbf{E})$.



D-separation

- A set of variables E d-separates X and Y if it blocks every undirected path in the BN between X and Y.
- If evidence E d-separates X and Y, then X and Y are conditionally independent given evidence E
- So what is blocking?

Let P be an **undirected path** from X to Y in a BN.

Let **E** be a set of variables.

We say *E blocks path P* iff there is some node Z on the path such that:

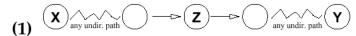
Case 1: one arc on P goes into Z and one goes out of Z, and Z∈E; or

Case 2: both arcs on P leave Z, and Z∈E; or

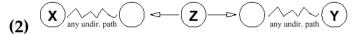
Case 3: both arcs on P enter Z and *neither Z, nor any of its descendants*, are in **E**.



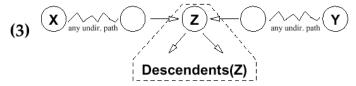
Blocking: Graphical View



If Z in evidence, the path between X and Y blocked



If Z in evidence, the path between X and Y blocked



If Z is **not** in evidence and **no** descendent of Z is in evidence, then the path between X and Y is blocked

Machine learning

- Decision-tree learning
- Naive Bayes and Bayes network parameter learning
- K-means and EM
- Chain rule for computing partial derivatives
- Linear and logistic regression
- Backpropagation
- Reinforcement learning: Value iteration, Q-learning and SARSA
- CNN



Decision tree learning

- Aim: find a small tree consistent with the training examples
- Idea: choose "most significant" attribute as root of (sub)tree

 $\begin{array}{ll} \textbf{function} & \textbf{DECISION-TREE-LEARNING} (examples, attributes, parent_examples) & \textbf{returns} \\ \textbf{a} \text{ tree} & \end{array}$

```
else if all examples have the same classification then return the classification else if attributes is empty then return PLURALITY-VALUE(examples) else A \leftarrow \operatorname{argmax}_{a \in attributes} \text{ IMPORTANCE}(a, examples) \\ tree \leftarrow a \text{ new decision tree with root test } A \\ \text{for each value } v_k \text{ of } A \text{ do} \\ exs \leftarrow \{e : e \in examples \text{ and } e.A = v_k\} \\ subtree \leftarrow \text{DECISION-TREE-LEARNING}(exs, attributes - A, examples) \\ \text{add a branch to } tree \text{ with label } (A = v_k) \text{ and subtree } subtree \\ \text{return } tree
```

if examples is empty then return PLURALITY-VALUE(parent_examples)

Plurality-value(examples) returns the majority classification of the examples

Entropy

• The entropy of a random variable V with values v_k , each with probability $P(v_k)$:

$$H(V) = -\sum_{k} P(v_k) \log_2 P(V_k)$$

 The entropy of a Boolean random variable that is true with probability q:

$$B(q) = -(q \log_2 q + (1 - q) \log_2 (1 - q))$$

 If a training set contains p positive examples and n negative examples, then the entropy of the goal attribute on the whole set is

$$H(Goal) = B(\frac{p}{p+n})$$



Information gain

- An attribute A with d distinct values divides the training set E into subsets E_1, \ldots, E_d .
- Each subset E_k has p_k positive examples and n_k negative examples,
- ullet So the expected entropy remaining after testing attribute A is

$$Remainder(A) = \sum_{k=1}^{d} \frac{p_k + n_k}{p + n} B(\frac{p_k}{p_k + n_k}).$$

 The information gain (IG) from the attribute test on A is the expected reduction in entropy:

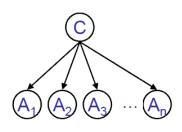
$$Gain(A) = B(\frac{p}{p+n}) - Remainder(A)$$

Choose the attribute with the largest IG



Naive Bayes models

- ullet Want to predict a class C based on attributes A_1,\ldots,A_n
- Parameters:
 - $\theta = P(C = true)$
 - $\theta_{i1} = P(A_i = true | C = true)$
 - $\theta_{i2} = P(A_i = true | C = false)$
- Assumption: A_i 's are independent given C



Naive Bayes learning

- Notation: $p = \#(c), n = \#(-c), p_i^+ = \#(c, a_i),$ $n_i^+ = \#(c, -a_i), p_i^- = \#(-c, a_i), n_i^- = \#(-c, -a_i)$
- $P(d|h) = \theta^p (1-\theta)^n \Pi_i \theta_{i1}^{p_i^+} \theta_{i2}^{p_i^-} (1-\theta_{i1})^{n_i^+} (1-\theta_{i2})^{n_i^-}$
- $\theta = p/(p+n)$, $\theta_{i1} = p_i^+/(p_i^+ + n_i^+)$, $\theta_{i2} = p_i^-/(p_i^- + n_i^-)$,
- $P(C|a_1,\ldots,a_n) = \alpha P(C) \Pi_i P(a_i|C)$
- Choose the most likely class



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Bayesian network parameter learning (ML)

- Parameters $\theta_{V,pa(V)=v}$:
 - CPTs: $\theta_{V,pa(V)=v} = P(V|pa(V)=v)$
- Data d:
 - $d_1 : \langle V_1 = v_1, V_2 = v_2, ..., V_n = v_{n,1} \rangle$
 - $d_2 : \langle V_1 = v_{1,2}, V_2 = v_{2,2}, ..., V_n = v_{n,2} \rangle$
- Maximum likelihood:
 - Set $\theta_{V,pa(V)=v}$ to the relative frequencies of the values of V given the values \mathbf{v} of the parents of V $\theta_{V,pa(V)=v} = \#(V,pa(V)=v) / \#(pa(V)=v)$



k均值算法

用于硬聚类

- 选择k个重心(centroid)。
- ② 寻找最近的重心并且更新聚类分配。将每个数据点都分配给 离它最近的重心的聚类。距离的度量通常是欧式距离。
- 将重心移动到它们的聚类的中心。每个聚类的重心的新位置 是通过计算该聚类中所有数据点的平均位置得到的。
- 重复第2和3步,直到每次迭代时重心的位置不再显著变化 (即直到该算法收敛)。

高斯混合模型

- 高斯混合模型的参数是:
 - $w_i = P(C = i)$ (每个成分的权重),
 - μ_i (每个成分的均值),
 - Σ_i (每个成分的协方差).

EM算法

随机初始化模型参数,重复以下两步直到收敛;

- ① E步
 - 计算数据x;是由成分i生成的概率

$$p_{ij} = P(C = i | \mathbf{x}_j) = \alpha P(\mathbf{x}_j | C = i) P(C = i),$$

其中
$$P(\mathbf{x}_j|C=i)$$
是第 i 个高斯分布, $P(C=i)=w_i$

- $\Diamond n_i = \sum_i p_{ij}$, 即当前分配到成分i的数据点的期望数量
- ② M步: 计算新的均值, 协方差, 和权重 ${m \mu}_i \leftarrow \sum_i p_{ij} {m x}_j / n_i$ $\Sigma_i \leftarrow \sum_i p_{ij} (\mathbf{x}_j - \boldsymbol{\mu}_i) (\mathbf{x}_j - \boldsymbol{\mu}_i)^\top / n_i$ $w_i \leftarrow n_i/N$

其中 N 是数据点的总数量.



Linear regression

- $h_w(x) = w \cdot x = \sum_i w_i x_i$
- Squared error loss: $Loss(w) = (y h_w(x))^2$
- Chain rule: $\partial g(f(x))/\partial x = g'(f(x))\partial f(x)/\partial x$
- $\partial Loss(w)/\partial w_i = -2(y h_w(x))x_i$
- $w_i \leftarrow w_i + \alpha(y h_w(x))x_i$



Logistic regression

- A logistic function is the sigmoid of a linear function
- Logistic regression: regression with logistic functions
- $g(x) = 1/(1 + e^{-x})$
- $\bullet \ h_w(x) = g(w \cdot x)$
- g' = g(1 g)
- $Loss(w) = (y h_w(x))^2$
- $\partial Loss(w)/\partial w_i = -2(y h_w(x))g'(w \cdot x)x_i$ = $-2(y - h_w(x))h_w(x)(1 - h_w(x))x_i$
- $w_i \leftarrow w_i + \alpha(y h_w(x))h_w(x)(1 h_w(x))x_i$



The algorithm

```
initialize w arbitrarily  \begin{aligned} \textbf{repeat} \\ & \text{for each } e \text{ in examples do} \\ & p \leftarrow g(w \cdot x(e)) \\ & \delta \leftarrow y(e) - p \\ & \text{for each } i \text{ do} \\ & w_i \leftarrow w_i + \alpha \delta p(1-p) x_i \end{aligned}  until some stopping criterion is satisfied return w
```

Forward and backward phases for backpropagation

Forward phase:

- Propagate inputs forward to compute the output of each unit
- Output a_j at unit j: $a_j = g(in_j)$ where $in_j = \sum_i w_{ij} a_i$

Backward phase:

- Propagate errors backward
- For an output unit j: $\Delta_j = g'(in_j)(y_j a_j) = a_j(1 a_j)(y_j a_j)$
- For an hidden unit i: $\Delta_i = g'(in_i) \sum_j w_{ij} \Delta_j = a_i (1 a_i) \sum_j w_{ij} \Delta_j$

Weight updating: $w_{ij} \leftarrow w_{ij} + \alpha a_i \Delta_j$



Markov Decision Processes

An MDP consists of:

- set S of states.
- set A of actions.
- P(s'|s,a) specifies the probability of transitioning to state s' given that the agent is in state s and does action a.
- R(s, a, s') is the expected reward received when the agent is in state s, does action a and ends up in state s'.
- $0 \le \gamma \le 1$ is discount factor.

Formulation of reinforcement planning

- A policy $\pi: S \times A \to [0,1]$, where $\pi(s,a)$ is the probability of taking action a in state s
- Value function $V:S \to \mathbb{R}$, where $V_{\pi}(s) = \mathbb{E}_{\pi}[G|S=s]$, meaning the expected discount award of following π in state s
- Given an environment MDP (S,A,P,R,γ) , where P and R are not known to the agent, find a policy π that maximizes $V_\pi(s_0)$

Q-learning

initialize Q[S,A] arbitrarily observe current state s repeat forever: select and carry out an action a observe reward r and state s' $Q[s,a] \leftarrow Q[s,a] + \alpha \left(r + \gamma \max_{a'} Q[s',a'] - Q[s,a] \right)$ $s \leftarrow s'$

SARSA (state-action-reward-state-action)

```
initialize Q[S,A] arbitrarily observe current state s select action a using a policy based on Q repeat forever:
    carry out action a observe reward r and state s' select action a' using a policy based on Q Q[s,a] \leftarrow Q[s,a] + \alpha \left(r + \gamma Q[s',a'] - Q[s,a] \right) s \leftarrow s' a \leftarrow a'
```

Q值更新公式的汇总

• 值迭代: 已知 P(s'|s,a) 和 R(s,a,s')

$$Q[s, a] \leftarrow \sum_{s'} P(s'|s, a) (R(s, a, s') + \gamma \max_{a'} Q[s', a'])$$

• Q-learning: 使用经验 $\langle s, a, r, s' \rangle$

$$Q[s, a] \leftarrow Q[s, a] + \alpha \left(r + \gamma \max_{a'} Q[s', a'] - Q[s, a]\right)$$

● Sarsa: 使用经验 ⟨s,a,r,s',a'⟩

$$Q[s, a] \leftarrow Q[s, a] + \alpha \left(r + \gamma Q[s', a'] - Q[s, a] \right)$$



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CNN: basic ideas

- Local connectivity: connect each neuron to only a local region of the previous layer
- Parameter sharing: weights shared across multiple local regions
- Downsampling to make images smaller

Convolutional Layer

- Accepts a volume of size $W_1 \times H_1 \times D_1$
- Requires four hyperparameters:
 - ullet Number of filters K,
 - their spatial extent F,
 - the stride S,
 - ullet the amount of zero padding P.
- Produces a volume of size $W_2 \times H_2 \times D_2$ where:
 - $W_2 = (W_1 F + 2P)/S + 1$
 - $H_2 = (H_1 F + 2P)/S + 1$
 - $D_2 = K$
- In the output volume, the d-th depth slice is the result of performing a convolution of the d-th filter over the input volume, and then offset by d-th bias.



Pooling layer

- Accepts a volume of size $W_1 \times H_1 \times D_1$
- Requires two hyperparameters:
 - ullet their spatial extent F,
 - \bullet the stride S,
- Produces a volume of size $W_2 \times H_2 \times D_2$ where:
 - $W_2 = (W_1 F)/S + 1$
 - $H_2 = (H_1 F)/S + 1$
 - D2 = D1
- For Pooling layers, it is not common to pad the input using zero-padding.

There are only two commonly seen variations of the max pooling layer found in practice: A pooling layer with F=3,S=2 (also called overlapping pooling), and more commonly F=2,S=2.