



PROBLEM SOLVING AND SEARCH IN ARTIFICIAL INTELLIGENCE

Lecture 3 Local Search

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Agenda

- 1 Introduction
- 2 Traditional Search Methods (e.g. Best First Search, A* Search, Heuristics)
- 3 Local Search, Stochastic Hill Climbing, Simulated Annealing
- 4 Tabu Search
- 5 Answer-set Programming (ASP)
- 6 Constraint Satisfaction (CSP)
- 7 Evolutionary Algorithms/ Genetic Algorithms
- 8 Structural Decomposition Techniques (Tree/Hypertree Decompositions)

Hill-climbing Methods

- Hill climbing methods use an **iterative improvement technique**.
- Technique is applied to a single point - the **current point** - in the search space.
- During each iteration, a new point is selected from the neighbourhood of the current point.
- If **new point provides better value** (in light of evaluation function) the new point **becomes the current point**.
- Otherwise, some other **neighbour is selected and tested** against the current point.
- The method **terminates if no further improvement is possible** (or, in the iterated version, if we run out of time).

Basic Hill-Climber

Algorithm basic hill-climber

```
initialize best
local  $\leftarrow$  FALSE
select a current point  $v_c$  at random
evaluate  $v_c$ 
repeat
  select all new points in the neighbourhood of  $v_c$ 
  select the point  $v_n$  from the set of new points with the best value of evaluation function eval
  if eval( $v_n$ ) is better than eval( $v_c$ ) then
     $v_c \leftarrow v_n$ 
  else
    local  $\leftarrow$  TRUE
  end if
until local
best  $\leftarrow v_c$ 
```

Weaknesses of Hill-climbing Algorithms

- 1 They usually **terminate at solutions** that are only **locally optimal**.
- 2 No information about how much the local optimum **deviates from the global optimum**, or from other local optima.
- 3 The obtained optimum depends on the **initial configuration**.
- 4 In general, it is **not** possible to provide an **upper bound** for the computation time.

But, they are **easy to apply**. All that is needed is:

- the representation,
- the evaluation function, and
- a measure that defines the neighbourhood around a given solution.

Iterated Hill-Climber

Algorithm iterated hill-climber

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t ← 0
initialize best
repeat
  local ← FALSE
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  evaluate  $v_c$ 
  repeat
    select all new points in the neighbourhood of  $v_c$ 
    select the point  $v_n$  from the set of new points with the best value of evaluation function eval
    if  $eval(v_n)$  is better than  $eval(v_c)$  then
       $v_c \leftarrow v_n$ 
    else
      local ← TRUE
    end if
  until local
  t ← t + 1
  if  $v_c$  is better than best then
    best ←  $v_c$ 
  end if
until t = MAX
```

Balance Between Exploration and Exploitation

Effective search techniques provide a mechanism for balancing two conflicting objectives:

- **exploiting** the best solutions found so far, and
- at the same time **exploring** the search space.

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Hill-climbing Techniques

Exploit the best available solution for possible improvement but **neglect exploring** a large portion of the search space.

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Random Search

Explores the search space thoroughly (points are sampled from the search space with equal probabilities) but **foregoes exploiting promising regions**.

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- **exploiting** the best solutions found so far, and
- at the same time **exploring** the search space.

There is no way to choose a single search method that can serve well in every case!

Local Search

- 1 Pick a solution from the search space and evaluate its merit. Define this as the **current** solution.
- 2 Apply **transformations** to the current solution to generate a neighbourhood of solutions and evaluate their merit.
- 3 Pick the **"best"** solution of the neighbourhood.
- 4 If the **new solution is "better"** than the current solution then **exchange** it with the current solution.
- 5 **Repeat** sets 2 and 3 **until no transformation** in the given set **improves** the current solution.

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 - Then, current solution has **no effect on the probabilities of selecting any new solution**.
 - The search becomes essentially **enumerative**.
 - Could be even **worse**: one might resample points that have already been tried.

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 - Then, current solution has **no effect on the probabilities of selecting any new solution**.
 - The search becomes essentially **enumerative**.
 - Could be even **worse**: one might resample points that have already been tried.
- Another extreme would be to always return the current solution - **this gets you nowhere!**

Local Search ctd.

- Searching within some **local neighbourhood** of current solution is a useful compromise.
- Then, current solution imposes a **bias on where we can search next**.
- If we find something better, we can update the current point to the new solution.
- If the **size of the neighbourhood is small**, the search might be **very quick**, but we might get trapped at **local optimum**.
- If the size of neighbourhood is **very large**, there is less chance to get stuck, but the **efficiency may suffer**.
- **The type of transformation we apply determines the size of neighbourhood.**

Local Search and the SAT

Local search algorithms are **surprisingly good** at finding satisfying assignments for certain classes of SAT formulas. **GSAT** is one of the **best-known** (randomized) **local search algorithms for SAT**.

Algorithm GSAT

```
for  $i \leftarrow 1$  step 1 to MAX-TRIES do
   $T \leftarrow$  a randomly generated truth assignment
  for  $j \leftarrow 1$  step 1 to MAX-FLIPS do
    if  $T$  satisfies the formula then
      return( $T$ )
    else
      make a flip
  end if
end for
return("no satisfying assignment found")
end for
```

Local Search and the SAT

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    else
      make a flip
  end if
end for
return("no satisfying assignment found")
end for
```

- "make a flip" flips the variable in T that results in the largest decrease in the number of unsatisfied clauses.
- MAX-TRIES, determines the number of new search sequences.
- MAX-FLIPS, determines the maximum number of moves per try.

Local Search and the SAT ctd.

- GSAT begins with randomly generated truth assignment.
- If assignment satisfies the problem, the algorithm terminates.
- Else, it flips each of the variables from TRUE to FALSE or FALSE to TRUE and records the decrease in the number of unsatisfied clauses.
- After trying all possible flips, it updates current solution to solution with largest decrease in unsatisfied clauses.
- If this new solution satisfies the problem, we are done.
- Otherwise, the algorithm starts flipping again.

Local Search and the SAT ctd.

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Interesting feature of the algorithm:

- Best available flip might **increase** the number of unsatisfied clauses.
- Selection is only made from neighbourhood of current solution. If every neighbour (defined as being one flip away) is worse than current solution, then GSAT takes the one that is the least bad.
- **Has the chance to escape local optimum!**
 - But, it might **oscillate** between points and never escape from some plateaus.

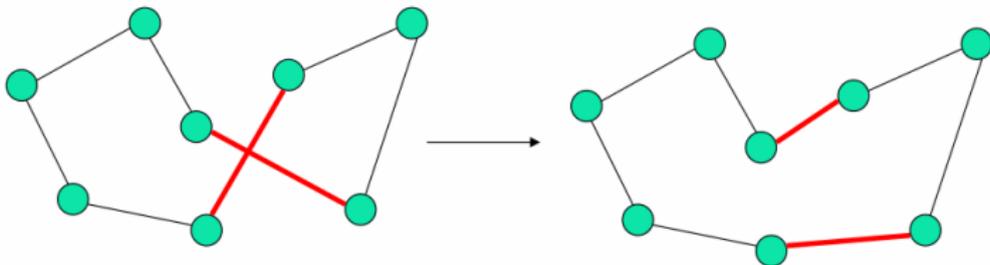
Local Search and the TSP

- There are many local search algorithms for TSP.
- The simplest is called *2-opt*.
- Starts with **random permutation** of cities (call this tour T) and tries to improve it.
- **neighbourhood** of T is defined as the **set of all tours** that can be **reached by changing two nonadjacent edges** in T .
- This move is called a *2-interchange*.

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2-interchange move



2-opt Algorithm

- A new tour T' replaces T if it is better.
 - Note: we replace the tour **every** time we find an improvement.
 - Thus, we terminate the search in the neighbourhood of T when the **first improvement** is found.
- If none of the tours in neighbourhood of T is better, then T is called **2-optimal** and algorithm terminates.
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- If none of the tours in neighbourhood of T is better, then T is called **2-optimal** and algorithm terminates.
- As GSAT, algorithm should be **restarted** from several random permutations.
- Can be **generalized to k -opt**, where either k or upto k edges are selected.
- Trade-off between size of neighbourhood and efficiency of the search:
 - If k is **small** the entire neighbourhood can be searched quickly, but increases likelihood of suboptimal answer.
 - For **larger values of k** , the number of solutions in neighbourhood become enormous (grows **exponentially** with k). Seldomly used for $k > 3$.

Escaping Local Optima

- Traditional problem-solving strategies either
 - **guarantee** discovering global solution, but are **too expensive**, or
 - have a tendency of "**getting stuck**" in **local optima**.
- There is almost no chance to speed up algorithms that guarantee finding global solution.
 - Problem of finding polynomial-time algorithms for real problems (as they are NP-hard).
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Simulated Annealing

Additional parameter (called **temperature**) that change the probability of moving from one point of the search space to another.

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Simulated Annealing

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Tabu Search

Memory, which forces the algorithm to explore new areas of the search space.

Local Search Revisited

Algorithm local search

$x =$ some initial starting point in \mathcal{S}

while $\text{improve}(x) \neq \text{"no"}$ **do**

$x = \text{improve}(x)$

end while

return(x)

Local Search Revisited

Algorithm local search

```
 $x$  = some initial starting point in  $\mathcal{S}$   
while improve( $x$ )  $\neq$  "no" do  
     $x$  = improve( $x$ )  
end while  
return( $x$ )
```

- **improve(x)** returns new point y from neighbourhood of x , i.e., $y \in N(x)$, if y is better than x ,
- otherwise, returns a string "no". In that case, x is a local optimum in \mathcal{S} .

Simulated Annealing

Algorithm simulated annealing

```
 $x =$  some initial starting point in  $\mathcal{S}$   
while not termination-condition do  
   $x =$  improve?( $x, T$ )  
  update( $T$ )  
end while  
return( $x$ )
```

Simulated Annealing vs. Local Search

There are three important differences:

- 1 How the procedure halts.
 - Simulated annealing is executed until some **external termination condition** is satisfied.
 - Local search is performed until no improvement is found.
- 2 $\text{improve?}(x, T)$ doesn't have to return a better point from the neighbourhood of x . It returns an **accepted** solution $y \in N(x)$, where acceptance is based on the current temperature T .
- 3 Parameter T is updated periodically, and the value of T influences the outcome of the procedure "improve?".

Iterated Hill-Climber Revisited

Algorithm iterated hill-climber

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t ← 0
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repeat
  local ← FALSE
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    else
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    end if
  until local
  t ← t + 1
  if  $v_c$  is better than best then
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  end if
until t = MAX
```

Modification of Iterated Hill-Climber

- **Instead** of checking **all** strings in the neighbourhood of v_c and selecting the best one, **select only one point**, v_n from this neighbourhood.
- Accept this new point, i.e., $v_c \leftarrow v_n$ with some probability that depends on the relative merit of these two points, i.e., the difference between the values returned by the evaluation function for these two points.

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⇒ **Stochastic hill-climber**

Stochastic Hill-Climber

Algorithm stochastic hill-climber

$t \leftarrow 0$

select a current point v_c at random

evaluate v_c

repeat

select the string v_n from the neighbourhood of v_c

select v_n with probability $\frac{1}{1+e^{\frac{eval(v_c)-eval(v_n)}{T}}}$

$t \leftarrow t + 1$

until $t = MAX$

Analyzing Stochastic Hill-Climber

- Probabilistic formula for accepting a new solution is based on **maximizing the evaluation function**.
- It has **only one loop**. No repeated calls from different random points.
- Newly selected point is **accepted** with probability p . Thus, the rule of moving from current point v_c to new neighbour, v_n , is probabilistic.
- New accepted point can be **worse** than current point.
- $$p = \frac{1}{1 + e^{\frac{eval(v_c) - eval(v_n)}{T}}}$$
- **Probability of acceptance** depends on the **difference in merit** between these two competitors, i.e., $eval(v_c) - eval(v_n)$, and on the value of an additional parameter T .
- T remains **constant** during the execution of the algorithm.

Role of Parameter T

Example:

- $eval(v_c) = 107$ and $eval(v_n) = 120$
- $eval(v_c) - eval(v_n) = -13$, new point v_n is better than v_c

Role of Parameter T

Example:

- $eval(v_c) = 107$ and $eval(v_n) = 120$
- $eval(v_c) - eval(v_n) = -13$, new point v_n is better than v_c
- What is probability of accepting new point based on **different values of T** ?

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- What is probability of accepting new point based on **different values of T** ?

T	$e^{\frac{-13}{T}}$	p
1	0.000002	1.00
5	0.0743	0.93
10	0.2725	0.78
20	0.52	0.66
50	0.77	0.56
10^{10}	0.9999...	0.5...

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- The greater T , the smaller the importance of the relative merit of the competing points!
- If T is **huge** (e.g., $T = 10^{10}$), the probability of acceptance approaches 0.5.
The search becomes random!
- If T is **very small** (e.g., $T = 1$), we have an **ordinary hill-climber!**

Role of new String

Suppose $T = 10$ and $eval(v_c) = 107$. Then, probability of acceptance depends only on the value of the new string.

$eval(v_n)$	$eval(v_c) - eval(v_n)$	$e^{\frac{eval(v_c) - eval(v_n)}{T}}$	p
80	27	14.88	0.06
100	7	2.01	0.33
107	0	1.00	0.50
120	-13	0.27	0.78
150	-43	0.01	0.99

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- If new point has **same merit** as current point, i.e., $eval(v_c) = eval(v_n)$, the probability of acceptance is 0.5.
- If new point is **better**, the probability of acceptance is greater than 0.5.
- The probability of acceptance **grows together with the (negative) difference** between these evaluations.

Simulated Annealing

- **Main difference** to stochastic hill-climber is that simulated annealing **changes the parameter T** during the run.
- **Starts with high values of T** making procedure more similar to random search, and then **gradually decreases** value of T .
- Towards end of the run, values of T are quite small, like an ordinary hill-climber.
- In addition, **new points are always accepted** if they are **better** than current point.

Simulated Annealing ctd.

Algorithm simulated annealing

```
t ← 0
initialize T
select a current point vc at random
evaluate vc
repeat
  repeat
    select new point vn in the neighbourhood of vc
    if eval(vc) < eval(vn) then
      vc ← vn
    else if random[0, 1) <  $e^{\frac{eval(v_n) - eval(v_c)}{T}}$  then
      vc ← vn
    end if
  until (termination-condition)
  T ← g(T, t)
  t ← t + 1
until (halting-criterion)
```

Simulated Annealing ctd.

- Is also known as Monte Carlo annealing, statistical cooling, probabilistic hill-climbing, stochastic relaxation, and probabilistic exchange algorithm.
- Based on an analogy taken from **thermodynamics**.
 - To **grow a crystal**, the raw material is heated to a molten state.
 - The temperature of the **crystal melt** is reduced until the crystal structure is **frozen in**.
 - Cooling should not be done too quickly, otherwise some irregularities are locked in the crystal structure.

Analogies Between Physical System and Optimization Problem

Physical System	Optimization Problem
state	feasible solution
energy	evaluation function
ground state	optimal solution
rapid quenching	local search
temperature	control parameter T
careful annealing	simulated annealing

Problem-Specific Questions

As with any search algorithm, simulated annealing requires answers for the following problem-specific questions.

- What is a solution?
- What are the neighbours of a solution?
- What is the cost of a solution?
- How do we determine the initial solution?

Answers yield the structure of the search space together with the definition of a neighbourhood, the evaluation function, and the initial starting point.

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Further Questions

- How do we determine the initial "temperature" T ?
- How do we determine the cooling ration $g(T, t)$?
- How do we determine the termination condition?
- How do we determine the halting criterion?

STEP 1:

$T \leftarrow T_{max}$
select v_c at random

STEP 2:

pick a point v_n from the neighbourhood of v_c

if $eval(v_n)$ is better than $eval(v_c)$ **then**

select if ($v_c \leftarrow v_n$)

else

select it with probability $e^{-\frac{\Delta eval}{T}}$

end if

repeat

this step

until k_T times

STEP 3:

set $T \leftarrow rT$

if $T \geq T_{min}$ **then**

goto STEP 2

else

goto STEP 1

end if

Where, T_{max} initial temperature, k_T number of iterations, r cooling ratio, and T_{min} frozen temperature.

Algorithm SA-SAT

```
tries  $\leftarrow$  0
repeat
   $v \leftarrow$  random truth assignment
   $j \leftarrow 0$ 
  repeat
    if  $v$  satisfies the clauses then
      return  $v$ 
       $T = T_{max} \cdot e^{-j \cdot r}$ 
      for  $k = 1$  to the number of variables do
        compute the increase (decrease)  $\delta$  in the number of clauses made true if  $v_k$  was flipped
        flip variable  $v_k$  with probability  $(1 + e^{-\frac{\delta}{T}})^{-1}$ 
         $v \leftarrow$  new assignment if the flip is made
      end for
       $j \leftarrow j + 1$ 
    end if
  until  $T < T_{min}$ 
  tries  $\leftarrow$  tries+1
until tries = MAX-TRIES
```

SA for SAT ctd.

- Outermost loop variable called "tries" keeps track of the number of **independent attempts** to solve the problem.
- T is set to T_{max} at the beginning of each attempt ($j \leftarrow 0$) and a new random truth assignment is made.
- Inner repeat loop **tries different assignments by probabilistically flipping** each of the Boolean variables.
- Probability of a flip depends on the **improvement δ of the flip** and the current temperature.
- If the improvement is negative, the flip is **unlikely to be accepted** and vice versa.
- r represents a decay rate for the temperature, the rate where it drops from T_{max} to T_{min} .
- The drop is caused by incrementing j , as $T = T_{max} \cdot e^{-j \cdot r}$.

SA-SAT vs. GSAT

- **Major difference:** GSAT can make a **backward move** (decrease in number of unsatisfied clauses) if other moves are not available.
- GSAT cannot make two backward moves in a row, as one backward move implies existence of next improvement move!
- SA-SAT can make an **arbitrary sequence of backward moves**, thus **escape local optima!**
- SA-SAT appeared to satisfy at least as many formulas as GSAT, with less work.
- **Applications** of SA: travelling salesman problem, production scheduling, Timetabling problems and image processing

Summary

- Hill-climbing methods face a danger of getting trapped in local optima and need to be started from different points.
- Simulated annealing is designed to escape local optima and can make uphill moves at any time.
- Hill-climbing and SA work on complete solutions.
- SA has many parameters to worry about (temperature, rate of reductions, ...).
- The more sophisticated the method, the more you have to use your judgement as to how it should be utilised.

References



Zbigniew Michalewicz and David B. Fogel.

How to Solve It: Modern Heuristics, volume 2. Springer, 2004.