Local Search (Ch. 4-4.1)



Announcements

HW1 late penalty

Test next week: will cover HW 1 & 2 -Ch 1-3 (up to A* and admissibility) -Open book/notes

Local search

Before we tried to find a path from the start state to a goal state using a "fringe" set

Now we will look at algorithms that do not care about a "fringe", but just neighbors

Some problems, may not have a clear "best" goal, yet we have some way of evaluating the state (how "good" is a state)

Local search

Today we will talk about 4 (more) algorithms:

- 1. Hill climbing
- 2. Simulated annealing
- 3. Beam search
- 4. Genetic algorithms

All of these will only consider neighbors while looking for a goal

Local search

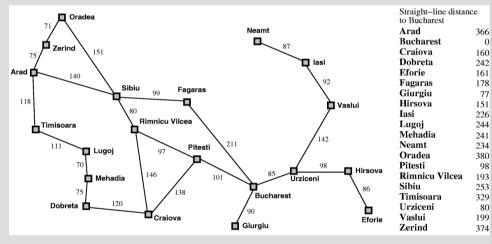
General properties of local searches:

- Fast and low memory
- Can find "good" solutions if can estimate state value
- Hard to find "optimal" path

In general these types of searches are used if the tree is too big to find a real "optimal" solution

Remember greedy best-first search?

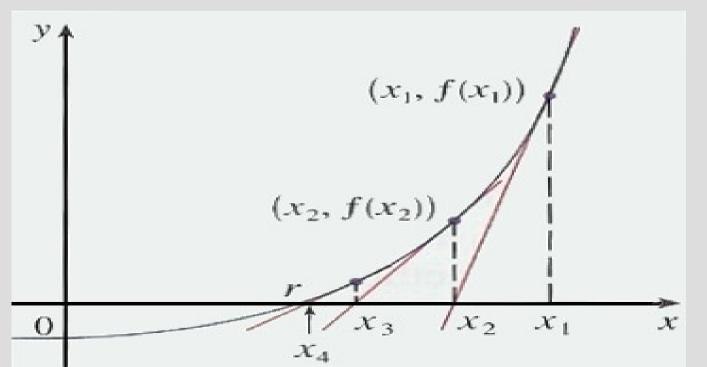
 Pick from neighbor with best heuristic
 Repeat 1...

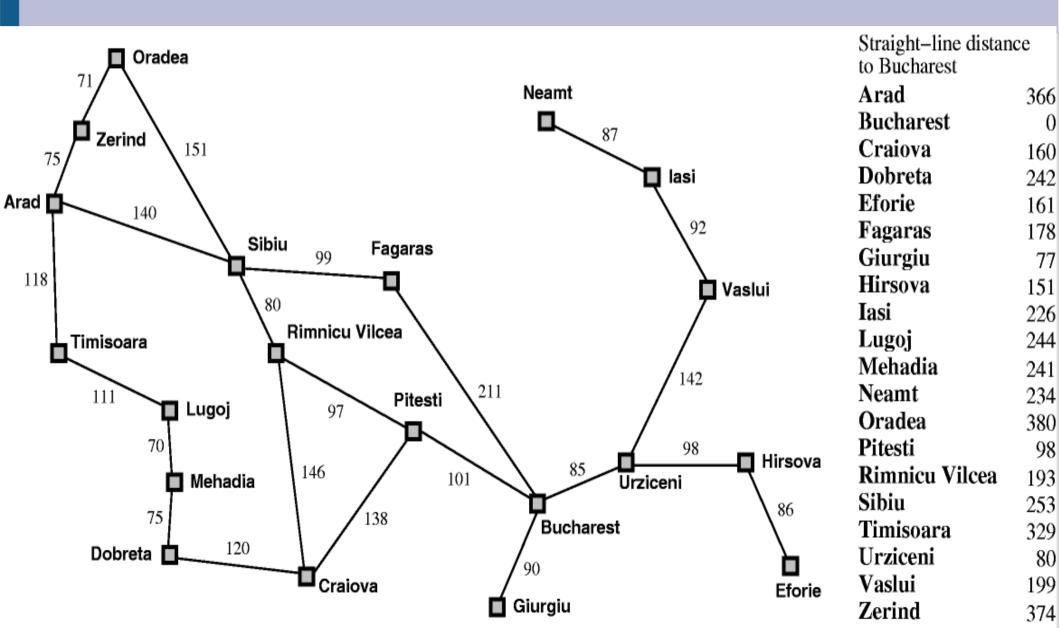


Hill climbing is only a slight variation:1. Pick best between: yourself and child2. Repeat 1...

What are the pros and cons of this?

This actually works surprisingly well, if getting "close" to the goal is sufficient (and actions are not too restrictive) Newton's method:

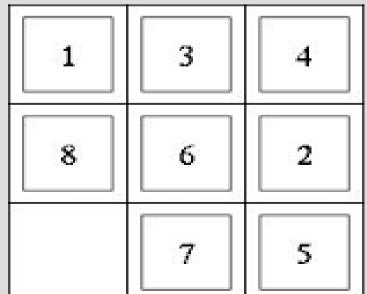




For the 8-puzzles we had 2 (consistent) heuristics:

h1 - number of mismatched pieces h2 - \sum Manhattan distance from number's current to goal position

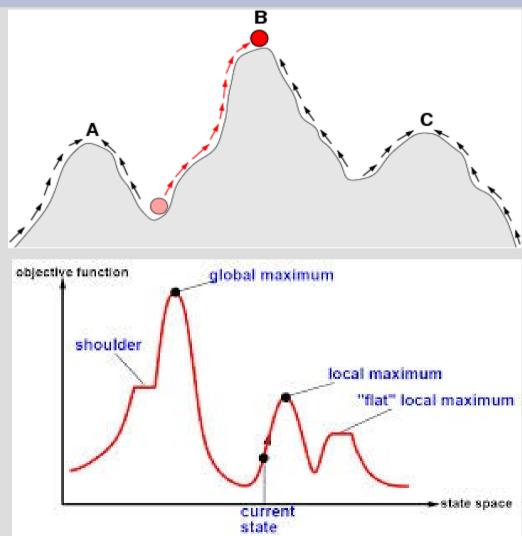
Let's try hill climbing this problem!



Can get stuck in:

- Local maximum
- Plateau/shoulder

Local maximum will have a range of attraction around it



Can get an infinite loop in a plateau if not careful (step count)

To avoid these pitfalls, most local searches incorporate some form of randomness

Hill search variants: Stochastic hill climbing - choose random move from better solutions

Random-restart hill search - run hill search until maximum found (or looping), then start at another random spot and repeat

The idea behind simulated annealing is we act more random at the start (to "explore"), then take greedy choices later

https://www.youtube.com/watch?v=qfD3cmQbn28 An analogy might be a hard boiled egg: 1. To crack the shell you hit rather hard (not too hard!)

- 2. You then hit lightly to create a cracked area around first
- 3. Carefully peal the rest

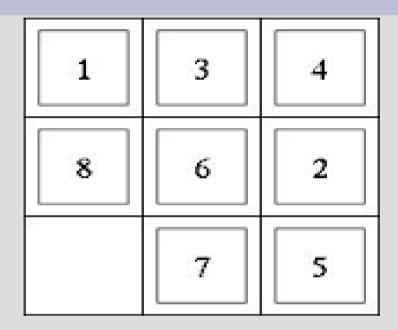


The process is:

- 1. Pick random action and evaluation result
- 2. If result better than current, take it
- 3. If result worse accept probabilistically
- 4. Decrease acceptance chance in step 3
- 5. Repeat...

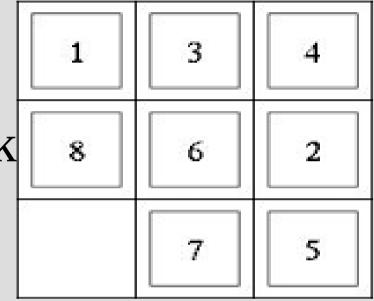
(see: SAacceptance.cpp) Specifically, we track some "temperature" T: 3. Accept with probability: $e^{\frac{result-current}{T}}$ 4. Decrease T (linear? hard to find best...)

Let's try SA on 8-puzzle:



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This example did not work well, but probably due to the temperature handling

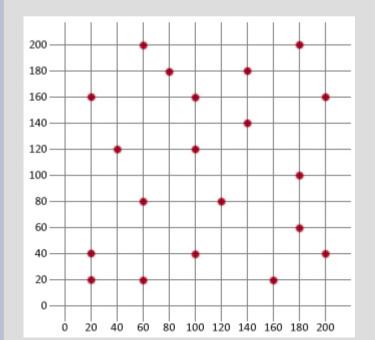


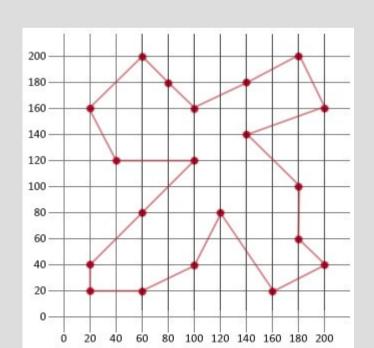
We want the temperature to be fairly high at the start (to move around the graph)

The hard part is slowly decreasing it over time

SA does work well on the traveling salesperson problem

(see: tsp.zip)





Beam search is similar to hill climbing, except we track multiple states simultaneously

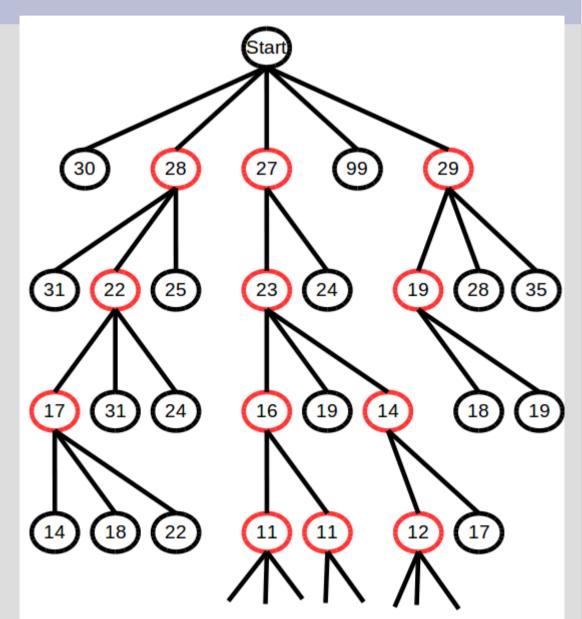
Initialize: start with K random nodes1. Find all children of the K nodes2. Add children and K nodes to pool, pick best3. Repeat...

Unlike previous approaches, this uses more memory to better search "hopeful" options

Beam search with 3 beams

Pick best 3 options at each depth to expand

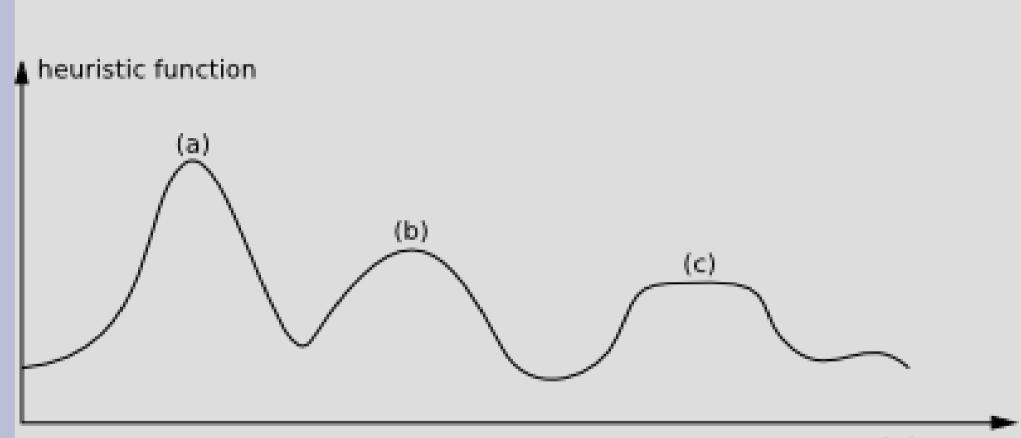
Stop like hill-climb (next pick is same as last pick)



However, the basic version of beam search can get stuck in local maximum as well

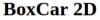
To help avoid this, stochastic beam search picks children with probability relative to their values

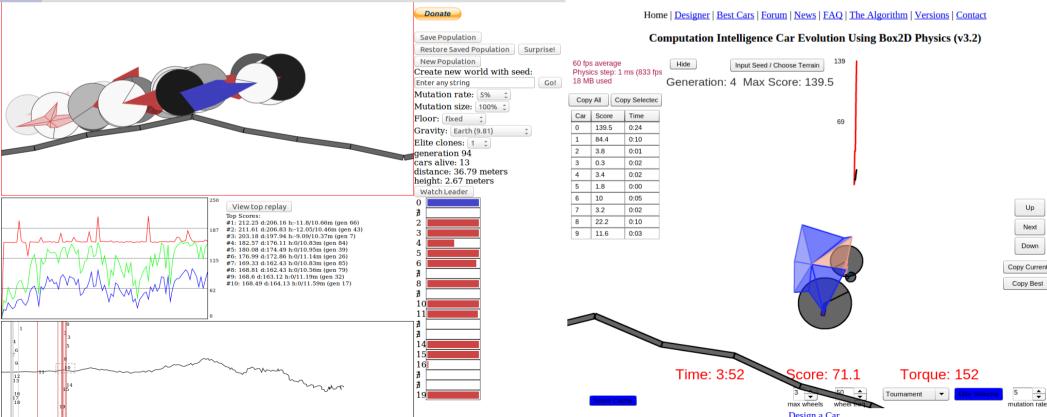
This is different that hill climbing with K restarts as better options get more consideration than worse ones



state space

Nice examples of GAs: http://rednuht.org/genetic_cars_2/ http://boxcar2d.com/





Genetic algorithms are based on how life has evolved over time

They (in general) have 3 (or 5) parts:
1. Select/generate children

1a. Select 2 random parents
1b. Mutate/crossover

2. Test fitness of children to see if they survive

3. Repeat until convergence

Selection/survival: Typically children have a probabilistic survival rate (randomness ensures genetic diversity) Mutation Crossover **Crossover:** 8 Split the parent's information into two parts, then take part 1 from parent A and 2 from B Mutation:

Change a random part to a random value

Genetic algorithms are very good at optimizing the fitness evaluation function (assuming fitness fairly continuous)

While you have to choose parameters (i.e. mutation frequency, how often to take a gene, etc.), typically GAs converge for most

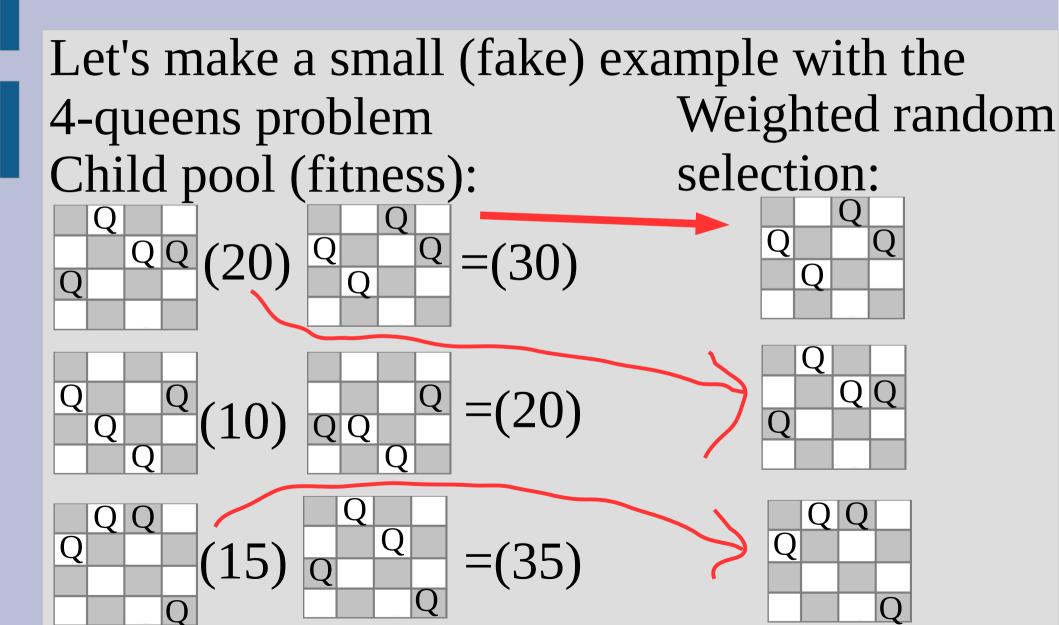
The downside is that often it takes many generations to converge to the optimal

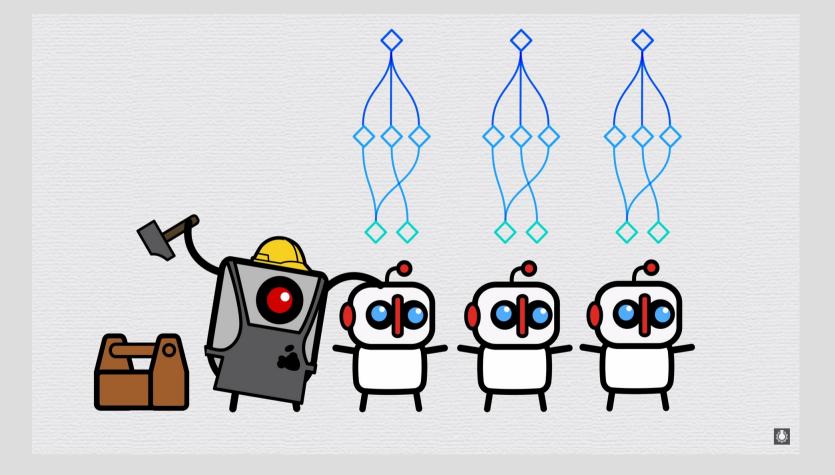
There are a wide range of options for selecting who to bring to the next generation:

- always the top (similar to hill-climbing... gets stuck a lot)
- choose purely by weighted random (i.e.
 4 fitness chosen twice as much as 2 fitness)
- choose the best and others weighted random

Can get stuck if pool's diversity becomes too little (hope for many random mutations)

Let's make a small (fake) example with the 4-queens problem Child pool (fitness): Adults: right QQ (20)Q =(30) \cap left Q Q =(20) Q (10)QQ Q 3/4 mutation $|\mathbf{Q}|$ Q Q Q (15)=(30) \cap (col 2)Ο 0





https://www.youtube.com/watch?v=R9OHn5ZF4Uo