

DPV Chapter 9

# Dealing with NP-Completeness

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Uncredited diagrams are from DPV or homemade.

# So the problem you want to solve is NP-Complete...

## Now what?

- ✗ Give up.
- ✗ Burn cycles and try to solve it exactly.
- ✗ Try the first thing that comes into your head and hope it produces correct answers and is fast enough to get by.
- ✓ Open a different tool box. (Chapter 9 of DPV.)

# Backtracking = exhaustive search + pruning

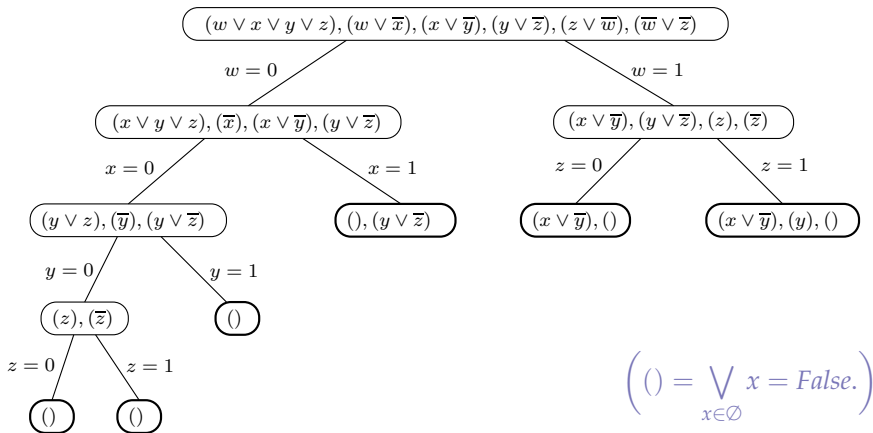
## Example: SAT via Backtracking

Let  $\varphi = (w \vee x \vee y \vee z) \wedge (w \vee \bar{x}) \wedge (x \vee \bar{y}) \wedge (y \vee \bar{z}) \wedge (z \vee \bar{w}) \wedge (\bar{w} \vee \bar{z})$ .

# Backtracking = exhaustive search + pruning

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# Backtracking: The general scheme

First we need a fast test for subproblems such that

$$\text{test}(P) = \begin{cases} \text{failure,} & \text{if subproblem } P \text{ has no solution;} \\ \text{success,} & \text{if a solution to } P \text{ is found;} \\ \text{uncertainty,} & \text{otherwise.} \end{cases}$$

Then:

```
Start with some problem  $P_0$ 
 $S \leftarrow \{P_0\}$  // the set of active subproblems
while ( $S \neq \emptyset$ ) do
  Choose a  $P \in S$ ;  $S \leftarrow S - \{P\}$ 
  Expand  $P$  into subproblems  $P_1, \dots, P_k$ 
  for  $i \leftarrow 1$  to  $k$  do
    case  $\text{test}(P_i)$  of
      success: announce solution and halt
      failure: discard  $P_i$ 
      uncertainty: add  $P_i$  to  $S$ 
Announce that there is no solution.
```

For SAT:

- ▶ *Choose*  $\equiv$  pick a clause
- ▶ *Expand*  $\equiv$  pick a variable in the clause

# Branch-and-Bound

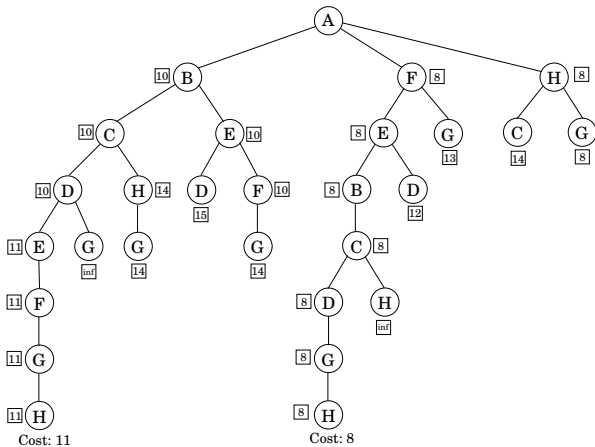
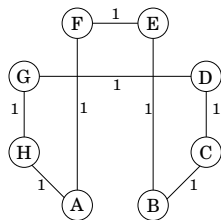
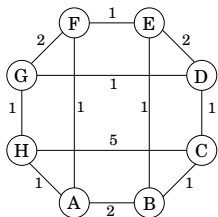
- ▶ B&B = the backtracking idea for optimization problems
- ▶ We consider minimization problems.
- ▶ First we need a fast way to compute *lower bounds* for the cost.
- ▶ Then:

```
Start with some problem  $P_0$   
 $S \leftarrow \{P_0\}$  // the set of active subproblems  
 $bestSoFar \leftarrow \infty$   
while ( $S \neq \emptyset$ ) do  
  Choose a  $P \in S$ ;  $S \leftarrow S - \{P\}$   
  Expand  $P$  into subproblems  $P_1, \dots, P_k$   
  for  $i \leftarrow 1$  to  $k$  do  
    if ( $P_i$  is a complete solution) then update  $bestSoFar$   
    else if ( $lowerbound(P_i) < bestSoFar$ ) then add  $P_i$  to  $S$   
return  $bestSoFar$ 
```

# Branch-and-Bound Applied to TSP, 1

- ▶  $G = (V, E)$  each  $e \in E$  with length  $d_e > 0$ .
- ▶ Fix an  $a \in V$ .
- ▶ Partial solution:  $[a, S, b]$  = a path from  $a$  to  $b$ ,  $S$  = the verts in this path
- ▶ Initial subproblem:  $[a, \{a\}, a]$ .
- ▶ Extension:  $[a, S \cup \{x\}, x]$  where  $x \in (V - S)$  and  $(b, x) \in E$ .
- ▶ *lowerbound* ( $[a, S, b]$ )
  - = a lower bound on the cost of completing the partial tour  $[a, S, b]$
  - = the sum of:
    - + the cheapest edge from  $a$  to  $V - S$ .
    - + the cheapest edge from  $b$  to  $V - S$ .
    - + the cost of a minimum spanning tree of  $V - S$ .
- ?? Why is this a lower bound on the cost of completing the partial tour  $[a, S, b]$ ?

# Branch-and-Bound Applied to TSP, 2



- ▶ 28 partial solutions examined.
- ▶  $7! = 5,040$  partial solutions in a brute-force search.



# Approximation Algorithms

- ▶ Instead of seeking an optimum solution, try “close to optimum”
- ▶ The question is how close is good enough.
- ▶  $opt(I)$  = the value of an optimum solution for instance  $I$ .
- ▶ **Convention:** Assume  $opt(I)$  is always a positive integer.
- ▶ **Convention:** Focus on **min**imization problems.
- ▶ Suppose  $\mathcal{A}(I)$  is the solution for  $I$  an algorithm  $\mathcal{A}$  returns.
- ▶ The *approximation ratio* for  $\mathcal{A}$  is

$$\alpha_{\mathcal{A}} = \max_I \frac{\mathcal{A}(I)}{Opt(I)} \geq 1.$$

- ▶ For **max**imization problems, take:

$$\alpha_{\mathcal{A}} = \max_I \frac{Opt(I)}{\mathcal{A}(I)} \geq 1.$$

- ▶ (The closer  $\alpha_{\mathcal{A}}$  is to 1 the better.)

## Recall from Chapter 5: Set Cover, 1

Suppose  $B$  is a set and  $S_1, \dots, S_m \subseteq B$ .

### Definition

- (a) A **set cover** of  $B$  is a  $\{S'_1, \dots, S'_k\} \subseteq \{S_1, \dots, S_m\}$  with  $B \subseteq \bigcup_{i=1}^k S'_i$
- (b) A **minimal set cover** of  $B$  is a set cover of  $B$  using as few of the  $S_i$ -sets as possible.

### The Set Cover Problem (SCP)

**Given:**  $B$  and  $S_1, \dots, S_m$  as above.

**Find:** A minimal set cover of  $B$ .

## Recall from Chapter 5: Set Cover, 1

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### The Set Cover Problem (SCP)

**Given:**  $B$  and  $S_1, \dots, S_m$  as above.

**Find:** A minimal set cover of  $B$ .

### Example

For:  $B = \{1, \dots, 14\}$  and

$$S_1 = \{1, 2\}$$

$$S_2 = \{3, 4, 5, 6\}$$

$$S_3 = \{7, 8, 9, 10, 11, 12, 13, 14\}$$

$$S_4 = \{1, 3, 5, 7, 9, 11, 13\}$$

$$S_5 = \{2, 4, 6, 8, 10, 12, 14\}$$

the solution to SCP is  $\{S_4, S_5\}$ .

## Recall from Chapter 5: Set Cover, 2

### A Greedy Approximation to the Set Cover Problem

// **Input:**  $B$  and  $S_1, \dots, S_m \subseteq B$  as above.

// **Output:** A set cover of  $B$  which is close to minimal.

$C \leftarrow \emptyset$

**while** (some element of  $B$  is not yet covered) **do**

    Pick the  $S_i$  with the largest number of uncovered  $B$ -elements

$C \leftarrow C \cup \{S_i\}$

**return**  $C$

### Example

$$B = \{1, \dots, 14\}$$

$$S_1 = \{1, 2\}$$

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$$S_3 = \{7, 8, 9, 10, 11, 12, 13, 14\}$$

$$S_4 = \{1, 3, 5, 7, 9, 11, 13\}$$

$$S_5 = \{2, 4, 6, 8, 10, 12, 14\}$$

On this, the algorithm returns  
 $\{S_1, S_2, S_3\}$ .

## Recall from Chapter 5: Set Cover, 3

### A Greedy Approx. to SCP

```
// Input:  $B$  and  $S_1, \dots, S_m \subseteq B$   
// Output: A near min. set cover  
 $\mathcal{C} \leftarrow \emptyset$   
while (all of  $B$  is not covered) do  
    Pick the  $S_i$  with the largest  
        number of uncovered  $B$ -elms  
     $\mathcal{C} \leftarrow \mathcal{C} \cup \{S_i\}$   
return  $\mathcal{C}$ 
```

### Claim

Suppose  $B$  contains  $n$  elements  
and the min. cover has  $k$  sets.

**Then** the greedy algorithm will  
use at most  $k \log_e n$  sets.

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So  $n_0 = n$ .

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- ▶ So there must be some set with at least  $n_t/k$  many elements.

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After iteration  $t$ :

- ▶ there are  $n_t$  elms left.
- ▶  $k$  many sets cover them
- ▶ So there must be some set with at least  $n_t/k$  many elements.
- ▶ So by the greedy choice,

$$\begin{aligned}n_{t+1} &\leq n_t - \frac{n_t}{k} = n_t \left(1 - \frac{1}{k}\right) \\ &= n_0 \left(1 - \frac{1}{k}\right)^t.\end{aligned}$$

## Recall from Chapter 5: Set Cover, 4

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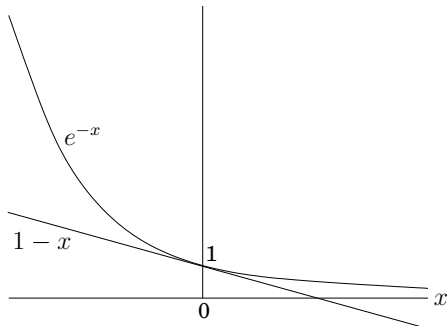
**Then** the greedy algorithm will use at most  $k \log_e n$  sets.

*Proof:* Let

$n_t$  = the number of uncovered elms after  $t$ -many while loop iterations

We know:  $n_{t+1} \leq n \left(1 - \frac{1}{k}\right)^t$ .

**Fact:**  $1 - x \leq e^{-x}$  for all  $x$ ,  
with equality iff  $x = 0$ .



## Recall from Chapter 5: Set Cover, 5

### A Greedy Approx. to SCP

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We know:  $n_{t+1} \leq n \left(1 - \frac{1}{k}\right)^t$ .

**Fact:**  $1 - x \leq e^{-x}$  for all  $x$ ,  
with equality iff  $x = 0$ .

$\therefore$  At  $t \geq k \log_e n$ ,  $n_t < ne^{-\log_e n} = 1$ ,  
i.e., we must have covered all of  $B$ .

So the greedy algorithm is optimal within a  $\log_e n$  factor.

That is,

$$\alpha_{\mathcal{A}} = \max_I \frac{\mathcal{A}(I)}{\text{Opt}(I)} \leq \log_e n.$$

# Approximating Vertex Cover, 1

## Vertex Cover (as an optimization problem)

**Given:**  $G = (V, E)$  an undirected graph

**Find:**  $S \subseteq V$  such that  $S$  touches every edge.

**Goal:** Minimize  $|S|$ .

- ▶ Vertex Cover is a special case of Set Cover.
- ▶ Therefore, it can be approximated within a  $O(\log n)$  factor.
- ▶ *However*, it turns out we can do much better.

# Approximating Vertex Cover, 2

## Definition

Suppose  $G = (V, E)$  an undirected graph.

- (a) A *matching* is an  $M \subseteq E$  such that any two edges in  $M$  have no endpoints in common.
- (b)  $M$  is a *maximum matching* when for each  $e \in (E - M)$ ,  $M \cup \{e\}$  fails to be a matching.

## Observations

- ▶ Maximal matchings are easy to construct. *(How?)*
- ▶ Fix  $G$ .
- ▶ If  $C$  is a vertex cover and  $M$  is a maximum matching, then each  $(u, v) \in M$  must have at least one of  $u$  and  $v$  in  $C$ . *(Why?)*
- ∴ (the size of a min. vertex cover for  $G$ )  $\geq$  (the size of a max. matching for  $G$ )
- ▶ If  $M$  is a maximal matching, then  $S = \{u \mid u \text{ is an endpoint of an } e \in M\}$  is a vertex cover. *(Why?)*
- ∴  $|S| = 2|M| \geq$  (the size of a min. vertex cover for  $G$ )  $\geq |M|$ .



# Approximating Vertex Cover, 3

## An approximation algorithm for Vertex Cover

**input**  $G = (V, E)$

Find a maximal matching  $M \subseteq E$ .

**return**  $S = \{u \mid u \text{ is an endpoint of an } e \in M\}$

- ▶ By the Observations, the approximation ratio of this algorithm is  $\alpha_A \leq 2$ .
- ▶ In fact, you can find examples where the ratio is exactly 2.
- ∴ The approximation ratio of *this* algorithm is  $\alpha_A = 2$ .
- ▶ *What about other algorithms?*

## Amazing Fact (Dinur and Safra, 2005)

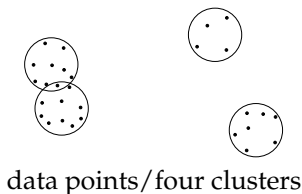
Minimum vertex cover cannot be approximated within a factor of 1.3606 for any sufficiently large vertex degree unless  $P=NP$ .

# Clustering, 1

## Definition

A *metric* on a space  $X$  is a function  $d: X \times X \rightarrow \mathbb{R}^{\geq 0}$  such that, for all  $x, y, z \in X$ :

1.  $d(x, y) \geq 0$ .
2.  $d(x, y) = 0 \implies x = y$ .
3.  $d(x, y) = d(y, x)$ .
4.  $d(x, y) \leq d(x, z) + d(z, y)$ .



## $k$ -Clustering

**Input:** Points  $X = \{x_1, \dots, x_n\}$ , metric  $d$ , integer  $k > 0$ .

**Output:** A partition of  $X$  into  $k$  clusters  $C_1, \dots, C_k$ .

**Goal:** Minimize the diameter of the clusters:  $\max_j \max_{x, x' \in C_j} d(x, x')$ .

- ▶  $k$ -Clustering is NP-complete.
- ▶  $k$ -Clustering is important in lots of areas (e.g., data mining).  
See [http://en.wikipedia.org/wiki/K-means\\_clustering](http://en.wikipedia.org/wiki/K-means_clustering)

# Clustering, 2

## Approximation Algorithm for $k$ -Clustering

Pick any point  $p_1 \in X$  to start

**for**  $i \leftarrow 2$  **to**  $k$  **do**

$p_i \leftarrow$  a point in  $X$  that is farthest away from  $p_1, \dots, p_{i-1}$

// i.e.,  $p_i$  maximizes:  $\min\{d(\cdot, p_j) : j = 1, \dots, i-1\}$

Create  $k$  clusters:  $C_i = \{x \in X : p_i \text{ is the closest center}\}$

**Claim:** For the above algorithm,  $\alpha_{\mathcal{A}} \leq 2$ .

*Proof:*

- ▶ Let  $x$  be the point farthest from  $p_1, \dots, p_k$ .
- ▶ Let  $r =$  the distance of  $x$  to the nearest  $p_i$ .
- ∴ Every point must be within  $r$  from its cluster center.
- ∴ The diameter of the clusters is  $\leq 2r$ .
- ▶ The points  $p_1, \dots, p_k$  and  $x$  are all  $\geq r$  distant from one another.
- ▶ Any partition of  $X$  into  $k$  cluster must put two of  $p_1, \dots, p_k, x$  into the same cluster. (By the PHP.)
- ∴ These clusters must have diameter  $\geq r$ . QED

# Traveling Salesman with metric distances, 1

## Traveling Salesman Problem

**Given:**  $n$  vertices and all  $n \cdot (n - 1)/2$ -many distances between them.

**Find:** An ordering of  $1, \dots, n$ :  $\pi(1), \pi(2), \dots, \pi(n)$  so that the tour's cost  $d(\pi(1), \pi(2)) + d(\pi(2), \pi(3)) + \dots + d(\pi(n), \pi(1))$  is minimal.

**Question:** Suppose we require the distances to come from a metric. Does this help make the problem easier? **Answer:** Yes!

## Definition (Repeated)

A *metric* on a space  $X$  is a function  $d: X \times X \rightarrow \mathbb{R}^{\geq 0}$  such that, for all  $x, y, z \in X$ :

1.  $d(x, y) \geq 0$ .
2.  $d(x, y) = 0 \implies x = y$ .
3.  $d(x, y) = d(y, x)$ .
4.  $d(x, y) \leq d(x, z) + d(z, y)$ .



# RECALL: Rudrata/Hamiltonian Cycle $\trianglelefteq$ TSP

## Rudrata/Hamiltonian Cycle Problem

**Given:**  $G = (V, E)$ , an undirected graph.

**Find:** A simple cycle that visits each vertex of  $G$ .

## Traveling Salesman Problem (TSP)

**Given:**  $V'$ ,  $n$  vertices; all  $\frac{n \cdot (n-1)}{2}$ -many distances between them; and  $b$ , a budget

**Find:**  $\pi$ , an ordering of  $1, \dots, n$ , such that  $\sum_{i=1}^n d_{\pi(i), \pi(1+(i \bmod n))} \leq b$ .

**Construction** of  $I(G, C)$ .

Given  $G = (V, E)$  and  $C \geq 1$ , define

$$V' = V$$

$$d_{i,j} = \begin{cases} 1, & \text{if } (i, j) \in E; \\ 1 + C, & \text{otherwise.} \end{cases}$$

$$b = |V|$$

**Claim:**  $(V, E)$  has a R/H cycle

$\iff (V', d)$  has tour of cost  $\leq b$ .

**If  $C \gg 1$ :**

- ▶ Gap: **either** a solution of cost  $n$ ,  
**or** solutions with costs  $\geq n + C$ ,  
**but** none inbetween.

$\therefore$  An approx. solution to (the full) TSP would let us solve Ham. Cycle in polytime!

**How?** (See next page.)

# Approximating General TSP

## Claim

An approximate solution to TSP would give us polytime solution of Rudrata Path.

## Proof

- ▶ Suppose that we had  $\mathcal{A}$ , a polytime approximation algorithm for TSP with approximation factor  $\alpha_{\mathcal{A}}$ .
- ▶ Suppose  $G$  is any instance of Rudrata Path.
- ▶ Construct  $I(G, C)$  where  $C = \alpha_{\mathcal{A}} \cdot n$  and run  $\mathcal{A}$  on it.
- ▶ If  $G$  has a Rudrata path, then  $OPT(I(G, C)) = n$  and  $\mathcal{A}$  finds a TSP tour of cost  $\alpha_{\mathcal{A}} \cdot OPT(I(G, C)) = \alpha_{\mathcal{A}} \cdot n$ .
- ▶ If  $G$  has no Rudrata path, then  $\mathcal{A}$  must return a tour of cost  $> \alpha_{\mathcal{A}} \cdot n$ .
- ▶ Since  $\mathcal{A}$  is supposed to run in polytime, this means we can decide Rudrata path in polytime!!!!

## Corollary

- ▶ *If TSP has a polytime approximation algorithm, then  $P=NP$ .*
- ▶ *If  $P \neq NP$ , then TSP has no polytime approximation algorithm.*

# Approximating Knapsack, 1

## Knapsack without repetition

### Given:

- A knapsack with capacity  $W$ .
- Items  $1, \dots, n$
- Item  $i$  has weight  $w_i$  & value  $v_i$ .

### Find: a set $M \subseteq \{1, \dots, n\} \ni$

- $\sum_{i \in M} w_i \leq W$  and
- $\sum_{i \in M} v_i$  is maximized.

- ▶ By Chapter 6, there is a dynamic programming solution to Knapsack that runs in  $O(n \cdot W) = O(n \cdot 2^{|W|})$  time.
- ▶ There is a similar dynamic programming solution to Knapsack that runs in  $O(n \cdot V) = O(n \cdot 2^{|V|})$  time, where  $V = \sum_{i=1}^n v_i$ .
- ▶ We use the  $O(n \cdot V)$  version as the basis for an approximation algorithm.



# Approximating Knapsack, 2

```
function ksApprox( $\vec{v}, \vec{w}, W, \epsilon$ ) //  $\epsilon$  = an approximation factor
// Assume each  $w_i \leq W$ .
 $v_{\max} \leftarrow \max\{v_i : i = 1, \dots, n\}$ .
for  $i = 1, \dots, n$  do  $\hat{v}_i \leftarrow \lfloor \frac{v_i \cdot n}{v_{\max} \cdot \epsilon} \rfloor$ . // Rescale the values
Run the dynamic programming algorithm using the  $\hat{v}_i$  values.
return the resulting choices of items
```

## Runtime Analysis

- ▶ Since each  $\hat{v}_i \leq n/\epsilon$ , we have  $\hat{v}_1 + \dots + \hat{v}_n \leq n^2/\epsilon$ .
- ▶ So the DP algorithm runs in  $O(n^3/\epsilon)$  time.

# Approximating Knapsack, 3

```
function ksApprox( $\vec{v}, \vec{w}, W, \epsilon$ ) //  $\epsilon$  = an approximation factor
// Assume each  $w_i \leq W$ .
 $v_{\max} \leftarrow \max\{v_i : i = 1, \dots, n\}$ .
for  $i = 1, \dots, n$  do  $\hat{v}_i \leftarrow \left\lfloor \frac{v_i \cdot n}{v_{\max} \cdot \epsilon} \right\rfloor$ . // Rescale the values
Run the dynamic programming algorithm using the  $\hat{v}_i$  values.
return the resulting choices of items
```

**Approximation Analysis** Suppose:

- ▶  $S$  is an optimal solution to the original problem with total value  $K^*$ .
- ▶  $\hat{S}$  is the solution produces for the scaled problem.

Then: 
$$\sum_{i \in \hat{S}} \hat{v}_i = \sum_{i \in \hat{S}} \left\lfloor \frac{v_i \cdot n}{v_{\max} \cdot \epsilon} \right\rfloor \geq \sum_{i \in \hat{S}} \left( \frac{v_i \cdot n}{v_{\max} \cdot \epsilon} - 1 \right) = K^* \cdot \frac{n}{v_{\max} \cdot \epsilon} - n.$$

So, the value of  $\hat{S}$  is at least  $K^* \cdot \frac{n}{v_{\max} \cdot \epsilon} - n$ . Hence,

**Correction:** The boxed part is what I missed in class.

$$\sum_{i \in \hat{S}} v_i \geq \frac{v_{\max} \cdot \epsilon}{n} \sum_{i \in \hat{S}} \hat{v}_i \geq \frac{v_{\max} \cdot \epsilon}{n} \left( K^* \cdot \frac{n}{v_{\max} \cdot \epsilon} - n \right) = K^* - v_{\max} \cdot \epsilon \geq K^*(1 - \epsilon).$$

# The approximability hierarchy

- ❖ No finite approximation ratio is possible.  
E.g., TSP.
- ❖ An approximation ratio of about  $\log n$  is possible.  
E.g., Set Cover.
- ❖ A constant approximation ratio is possible, *but* there are limits to how small this can be.  
E.g., Vertex Cover,  $k$ -Clustering, and metric TSP.  
*The proofs of these lower limit results are really hard!!!*
- ❖ A constant approximation ratio is possible, and in fact you can make  $\alpha_{\mathcal{A}}$  arbitrarily close to 1.  
E.g., Knapsack.

**NOTE:** All of the above assumes  $P \neq NP$ .

- ❖ If  $P=NP$ , all the problems can be solved exactly in polytime.

## Local search heuristics: The general scheme

```
 $s \leftarrow$  any initial solution  
while there is a solution  $s'$  in the neighborhood of  $s$  with  $cost(s') < cost(s)$  do  
   $s \leftarrow s'$   
return  $s$ 
```

For any application of this scheme to a particular problem,  
the key question *what is a good notion of neighborhood?*

# Local search heuristics: Traveling Salesman, 1

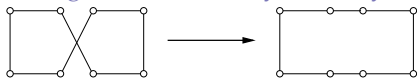
- ▶ Assume we have a complete graph on  $n$  vertices  
(with a cost assigned to each edge).
- ▶ So there are  $(n - 1)!$  many tours.
- ▶ Two tours differ by at least two edges. (Why?)
- ▶ So let us try:  
Tours  $T_1$  and  $T_2$  are neighbors when they differ by two edges.



- ▶ With this choice of “neighbor”:
  1. What is the overall running time?
  2. Does this always return an optimal answer?

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- ▶ With this choice of “neighbor”:
  1. What is the overall running time?
  2. Does this always return an optimal answer?
- ▶ Answers:
  1. Hard to say.
  2. Of course not.

# Local search heuristics: Traveling Salesman, 2

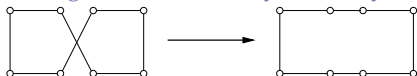
- ▶ Tours  $T_1$  and  $T_2$  are neighbors when they differ by two edges.



- ▶ With this choice of “neighbor”:  
**What is the overall running time?**
  - Each tour has  $O(n^2)$  neighbors, so making the choice is not too expensive.
  - But, the algorithm may well go through exponentially many iterations.

# Local search heuristics: Traveling Salesman, 3

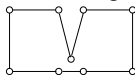
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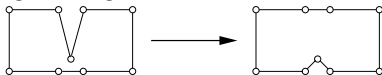
- ▶ With this choice of “neighbor”:

Does this always return an optimal answer?

- The final answer will be *locally optimal*, but not necessarily optimal.
- The problem is that this notion of neighbor is too myopic. E.g.,



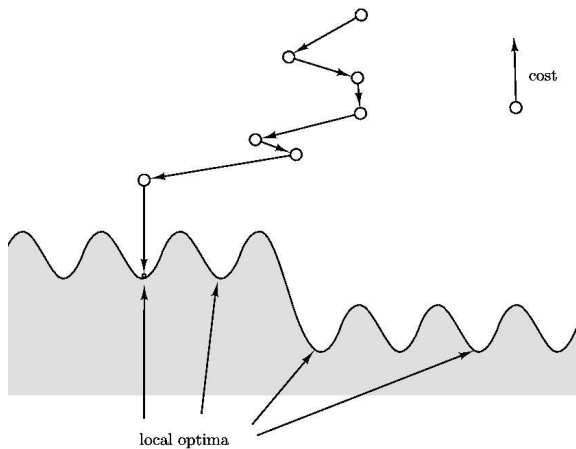
- ▶ If we allow three-edge changes, then:



but then a tour has  $O(n^3)$  neighbors and the choice part of the algorithm slows down.



# Local search heuristics: Optima, Local vs. global



# Local search: Graph partitioning, 1

## Graph partitioning

**Given:**  $G = (V, E)$ , an undirected graph with nonnegative edge wghts, and  $\alpha \in (0, 1/2]$ .

**Return:** A partition of  $V$  into  $A$  and  $B$  with

$$|A|, |B| \geq \alpha|V|.$$

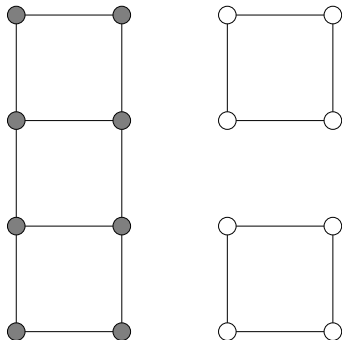
**Goal:** Minimize the capacity of the  $(A, B)$ -cut.

**Note:** The general problem is reducible to the special case of  $\alpha = 1/2$ .

**Strategy:**

- ▶ Start with a partition with  $|A| = |B|$ .
- ▶ Neighbors of  $(A, B) =$

$$\{(A - \{a\} + \{b\}, B - \{b\} + \{a\}) : a \in A, b \in B\}.$$



## Dealing with NP-Completeness

2019-04-17

## └ Local search: Graph partitioning, 1

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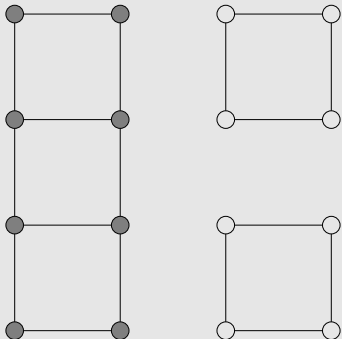
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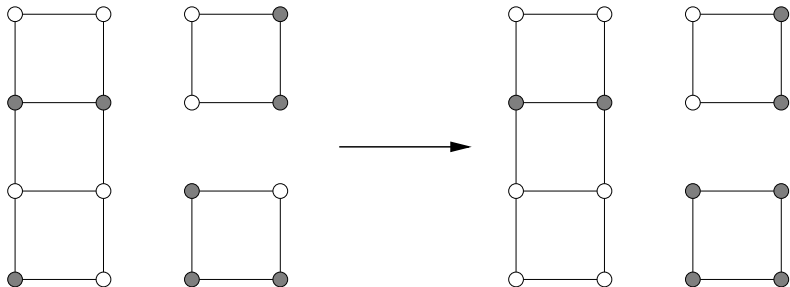


- $A =$  gray verts,  $B =$  white verts.
- Weights 0 and 1
- Optimal partition as cost 0.

## Local search: Graph partitioning, 2

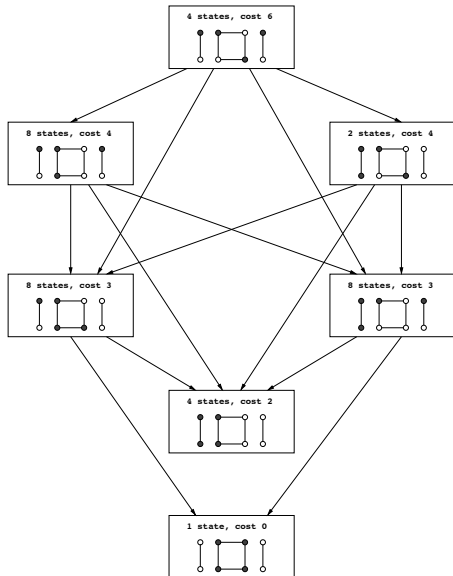
- ▶ Start with a partition with  $|A| = |B|$ .
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# Local search: Graph partitioning, 3

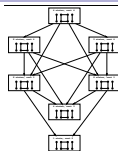
- ▶ The problem with this notion of neighbor is that there are stubborn local minima.



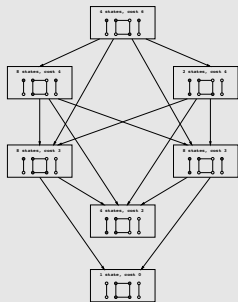
## Dealing with NP-Completeness

2019-04-17

## └ Local search: Graph partitioning, 3



► The problem with this notion of neighbor is that there are stubborn local minima.



- Search space for a graph with 8 nodes.
- Then entire space has 35 solutions, but the picture has grouped these into seven groups to cut the clutter.
- There are five local optima.

# Dealing with local optima: Randomized Restarts

```
 $L \leftarrow$  an empty list  
repeat  $k$  times  
   $s \leftarrow$  a randomly chosen initial solution  
  while (there is a solution  $s'$  in the neighborhood of  $s$  with  $cost(s') < cost(s)$ ) do  
     $s \leftarrow s'$   
    add  $s$  to  $L$   
end-repeat  
return the best solution in  $L$ 
```

This can shake free of bad local optima.

# Dealing with local optima: Simulated Annealing

```
s ← a randomly chosen initial solution
repeat
  s' ← a randomly chosen solution in the neighborhood of s
  Δ ← cost(s') - cost(s)
  if (Δ < 0) then s ← s'
  else with probability e-Δ/T do s ← s'
until we decide we are done
```

- ▶  $T \equiv$  temperature
- ▶ If  $T \approx 0$  this is roughly the previous scheme.
- ▶ If  $T$  is big, then  $s$  jumps around a lot.
- ▶ We vary  $T$ , initially large (hot), and gradually small (cooler).

