

# Metaheuristics in Scheduling

## Local Search and Genetic Algorithms

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# Metaheuristics in Scheduling

- $NP$ -hard scheduling problem  $\rightarrow$  Branch-and-Bound?
  - ... usually only valid for small instances!
- $NP$ -hard scheduling problem  $\rightarrow$  Approximation?
  - ... does not work in general!
- $NP$ -hard scheduling problem  $\rightarrow$  Heuristic?
  - ... is there a *general* approach to design a non-trivial heuristic?

We discuss two general techniques for solving optimization problems heuristically.

# Local Search Algorithms

Recap: *Discrete Optimization Problem* (Minimization)

A (*Discrete*) *Optimization Problem* is given by its problem description  $\Pi = (\mathcal{I}, \mathcal{S})$ , where

- $\mathcal{I}$  is the set of instances, and
- $\mathcal{S}(x)$  is the (discrete) set of feasible solutions for an instance  $x \in \mathcal{I}$ ,

together with an objective function  $f : \mathcal{S}(x) \rightarrow \mathbb{R}$  that evaluates each feasible solution.

We then seek—given an instance  $x$ —a feasible solution  $y \in \mathcal{S}(x)$  with minimum objective function value.

# Basic Structure of Local Search

Suppose we are given an instance  $x \in \mathcal{I}$ .

- $\mathcal{S} = \mathcal{S}(x)$  is a discrete set

*Local Search* is an iterative procedure that *moves* from one solution in  $\mathcal{S}$  to the next (until some stopping criterion is satisfied).

- ... think of it as the discrete analogy of *hill-climbing*

# Neighborhoods on the Solution Space

In order to move systematically through the solution set, the possible moves from one solution to another is restricted by

- *neighborhood structures*  $\mathcal{N} : \mathcal{S} \rightarrow 2^{\mathcal{S}}$ .

For each solution  $s \in \mathcal{S}$ , the set  $\mathcal{N}(s)$  describes the subset of solutions which can be reached from  $s$  in the next step.

- $\mathcal{N}(s)$  is called *neighborhood of  $s$*

# Neighborhood Graph

A neighborhood structure  $\mathcal{N}$  may be represented by a directed graph  $G = (V, A)$  where

- $V = \mathcal{S}$
- $(u, v) \in A \iff v \in \mathcal{N}(u)$ .

This graph is called the *neighborhood graph*.

Note: in general, it is not possible to store this graph completely!

- $\mathcal{S}$  usually has exponential size w.r.t. the instance.

# Allowed Modifications

In order to avoid problems with size of the neighborhood graph, a neighborhood is usually described by *operators*:

- Let  $F : \mathcal{S} \rightarrow \mathcal{S}$  be a function,
- for each feasible  $s \in \mathcal{S}$ ,  $F(s)$  is a subset consisting only of feasible solutions, we call  $F$  thus an *allowed modification*.

For every  $s \in \mathcal{S}$ , we can define a neighborhood structure for a set  $AM$  of allowed modifications as follows

$$\mathcal{N}(s) := \{F(s) \mid F \in AM\}.$$

# Connectivity of the Neighborhood Graph

Suppose the neighborhood graph  $G = (V, A)$  is connected

- $\Rightarrow$  for every (starting) solution  $s \in \mathcal{S}$ , there exists a directed path to *every other solution* in  $\mathcal{S}$ .
- In particular, we can provide a sequence of operations to  $s$  that result in an optimal solution  $s^* \in \mathcal{S}$ .



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Well, this is usually overkill!

We only need the latter condition:

- A neighborhood  $\mathcal{N}$  is called *OPT-connected* if, from each solution  $s \in \mathcal{S}$ , an optimal solution can be reached by a finite sequence  $s, s_1, \dots, s_k, s_{k+1}$  of solutions  $s_i \in \mathcal{S}$  s.t.  $s_{i+1} \in \mathcal{N}(s_i)$  for  $i = 1, \dots, k$ , and  $s_{k+1}$  optimal.

# Example: Exchange Neighborhood

Recall: many (active) schedule are completely described by permutations.

*Swap Neighborhood* for a permutation  $\pi$ :

- Swap two adjacent elements in  $\pi$ .

*Exchange Neighborhood* for a permutation  $\pi$ :

- New permutation  $\pi'$  with

$$\pi'(a) := \pi(b) \wedge \pi'(b) := \pi(a)$$

for two indexes  $a$  and  $b$ .

(In order to obtain an allowed modification, we need to restrict all permutations to feasibility, this is problem specific!)

# Local Search Method

Given a solution  $s \in \mathcal{S}$ ,

- in each iteration, we choose a solution  $s' \in \mathcal{N}(s)$  (or the allowed modification that yields  $s'$ ), and
- based on the objective function values  $f(s)$  and  $f(s')$ , we choose a starting solution for the next iteration.

According to different criteria for the choice of the next solution, different types of local search methods emerge.

$\mathcal{N}$  OPT-connected  $\Rightarrow$  independent of starting solution, we are *able* to reach optimality.

$\mathcal{N}$  *not* OPT-connected  $\Rightarrow$  may happen that we are unable to even reach an optimal solution at all.

# Iterative Improvement

For a local search approach, the simplest choice is to always take a neighboring solution with smallest objective value.

Algo.: *Iterative Improvement*

- ① Generate initial solution  $s \in \mathcal{S}$
- ② WHILE  $\exists s' \in \mathcal{N}(s) \mid f(s') < f(s)$  DO
  - ① Choose best solution  $s' \in \mathcal{N}(s)$ ;
  - ②  $s := s'$ ;

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... terminates with *local minimum*  $s^*$  w.r.t. neighborhood  $\mathcal{N}$

- ... we could start the algorithm several times with different solutions
- ... we could also accept solutions with increasing objective value
  - need strategies to avoid cycling!

# Simulated Annealing

Idea: avoid cycling by randomization, i.e. simulate the annealing process from physics

- choose a solution  $s' \in \mathcal{S}$  randomly
- accept solution only with a certain probability

In the  $i$ -th iteration,  $s'$  is accepted with probability

$$\min\left\{1, e^{\frac{f(s')-f(s)}{t_i}}\right\}$$

where  $(t_i)$  is a sequence of positive control values with

$$\lim_{i \rightarrow \infty} t_i = 0.$$

# Simulated Annealing

## Algo.: *Simulated Annealing*

- ①  $i := 0$
- ② Generate initial solution  $s \in \mathcal{S}$
- ③  $best := f(s)$
- ④ REPEAT
  - ① Generate randomly a solution  $s' \in \mathcal{N}(s)$
  - ② IF  $\text{Rand}(0,1) < \min\{1, e^{\frac{f(s')-f(s)}{t_i}}\}$  THEN
    - ①  $s := s'$
    - ② IF  $f(s') < best$  THEN
    - ③  $s^* := s$
    - ④  $best := f(s')$
- ⑤  $i := i + 1$
- ⑥ UNTIL some stopping condition is satisfied



# Threshold Acceptance

- often,  $(t_i)$  is defined (in analogy to physics) as

$$t_{i+1} := \alpha t_i, 0 < \alpha < 1$$

- search may be stopped after number of iterations, certain number of non-improving solutions, time limit, ...

Other variant of S.A. is given by *Threshold Acceptance*, where

- acceptance rule for  $s' \in \mathcal{N}$  is accepted if difference  $f(s) - f(s')$  is within some limit  $l_i$
- $l_i$  also decreases with number of iterations

Another *deterministic* strategy to avoid cycling is to store all visited solutions in a so-called *tabu-list*  $T$ .

- $\Rightarrow$  a neighbor is only accepted if it is not contained in  $T$

## Algo.: *Tabu Search*

- 1 Generate initial solution  $s \in \mathcal{S}$
- 2  $best := f(s); s^* := s$
- 3  $T := \emptyset$
- 4 REPEAT
  - 1  $Cand(s) := \{s' \in \mathcal{N}(s) \mid \text{move from } s \text{ to } s' \text{ is not tabu} \}$
  - 2 Choose a solution  $s' \in Cand(s)$
  - 3 Update  $T$
  - 4  $s := s'$
  - 5 IF  $f(s') < best$  THEN
    - 6  $s^* := s$
    - 7  $best := f(s')$
- 5 UNTIL some stopping condition is satisfied

# Tabu Search

Another *deterministic* strategy to avoid cycling is to store all visited solutions in a so-called *tabu-list*  $T$ .

- $\Rightarrow$  a neighbor is only accepted if it is not contained in  $T$

... due to memory constraints, this may not be possible!

- $T$  may contain only the  $|T| \leq B$  visited solutions
  - only cycles of length greater than  $B$  may occur
  - if  $B$  sufficiently large, the probability of cycling becomes small
- $T$  may not contain complete solution descriptions, but only *attributes* of already visited solutions
  - all solutions having one of the stored attributes are tabu
  - solution will not be re-visited as long as its attributes are stored in  $T$
- Disadvantage: also new solutions may be declared tabu!
  - *aspiration criteria*: accept solution even if they are tabu
  - e.g. based on objective function value

## Algo.: *Tabu Search*

- ① Generate initial solution  $s \in \mathcal{S}$
- ②  $best := f(s); s^* := s$
- ③  $T := \emptyset$
- ④ REPEAT
  - ①  $Cand(s) := \{s' \in \mathcal{N}(s) \mid \text{move from } s \text{ to } s' \text{ is not tabu OR } s' \text{ satisfies the aspiration criterion} \}$
  - ② Choose a solution  $s' \in Cand(s)$
  - ③ Update  $T$
  - ④  $s := s'$
  - ⑤ IF  $f(s') < best$  THEN
  - ⑥      $s^* := s$
  - ⑦      $best := f(s')$
- ⑤ UNTIL some stopping condition is satisfied

# Neighbor Selection

Depending on the size of the neighborhood, several selection strategies for  $Cand(s)$  emerge:

- *best-fit*: explore entire neighborhood and take best neighbor
- *first-fit*: explore neighborhood and take first neighbor that improves current solution
  - if no such neighbor exists, take the best one from  $Cand(s)$
- ...

# Tabu List Management

For tabu-list management, two types are distinguished.

*static* tabu-lists

- constant size

*dynamic* tabu-lists

- variable length
- if a solution is found that improves the current leader, the list is emptied as we have never visited this part of the solution space before
- improving phase of T.S.: *decrease* length of list
- non-improving phase: *increase* length of list

Generally speaking, a tabu-list serves a *short-term* memory of the local search procedure.

Besides short-term memory (T.L.), also *long-term* memory may be kept that is used for *diversification*.

Here, properties of promising solutions not explored further are stored which are then used in a restarting phase:

- If within a certain number of iterations, the current leader is not improved (*intensification*), then
- the search process is stopped and restarted with a new solution (*diversification*).

Note: a restart from a randomly generated solution would neglect all information of the previous search process.



# Application of Local Search

Arriving at a local search algorithm for a specific problem:

- Define problem specific ingredients of local search:
  - most importantly: the neighborhood
- *Tune* the chosen local search approach.

Claim:

The *problem specific ingredients* are far more important than the *tuning*.

# Efficiency of Local Search

*Local efficiency* (one iteration):

- quality of  $s'$  or  $\mathcal{N}(s)$
- computational time to calculate and evaluate  $s'$
- size of  $\mathcal{N}(s)$

Note: large size of neighborhood needs not result in large computational time (see c.f. research on VLSN: efficient search for optimal solution w.r.t. neighborhood)

*Global efficiency*:

- number of iterations, computational time
- quality of final solution
  - related to *price of anarchy* (game theory)

# Applying Tabu Search to the Job Shop Problem

## Recap: Job-Shop Problem $J||C_{\max}$

- $n$  jobs  $j = 1, \dots, n$  consisting of  $n_j$  operations
- $m$  machines
- each operation  $O_{ij}$  has machine  $\mu_{ij}$  and processing time  $p_{ij}$

## Recap: Disjunctive Graph Model and Complete Selections

- complete selection  $\rightarrow$  all arcs in model fixed
  - cycle-free  $\iff$  feasible solution
- ... can use permutation of operations per machine to describe complete selection uniquely
- ... can use longest path calculation to determine starting time of each operation (critical path)

Use  $\pi = (\pi_1, \dots, \pi_m)$  to describe the set  $\mathcal{S}$  of solutions.

# TS-JS: Neighborhood Structures

Apply the Swap-Neighborhood approach based on the following lemma.

Lemma: Let  $s$  be a complete selection, and let  $P$  be a longest path in  $G(s)$ .

Let  $(v, w)$  be an arc of  $P$  such that  $v$  and  $w$  are processed on the same machine. Then,  $s'$  obtained by reversing  $v$  and  $w$  is again a complete selection.

(Proof on the board)

We call the resulting neighborhood  $\mathcal{N}_1$ .

# TS-JS: Neighborhood Structures

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Theorem:  $\mathcal{N}_1$  is OPT-connected.

(Proof on the board)

Consider a feasible solution, i.e. complete selection,  $s = \pi$ .

Definition (*Block*): Let  $G(s) = (V, C \cup D)$  be the graph induced by the complete selection  $s$ , and let  $P$  be a critical path in  $G(s)$ . A sequence  $u_1, \dots, u_k$  of successive nodes in  $P$  is called *block* if the following two properties hold:

- (i) The sequence contains at least two nodes.
- (ii) The sequence represents a maximal number of operations to be processed on the same machine.

We denote the  $j$ -th block on a given critical path  $P$  by  $B_j$ .

Lemma: Let  $s$  be a complete selection corresponding to a feasible solution for the job-shop problem. If there exists another selection  $s'$  such that  $L(s') < L(s)$  holds, then in  $s'$  at least one operation from some block  $B$  of  $G(s)$  has to be processed before the first or after the last operation of  $B$ .  
(Proof on the board.)

Lemma: Let  $s$  be a complete selection corresponding to a feasible solution for the job-shop problem. If there exists another selection  $s'$  such that  $L(s') < L(s)$  holds, then in  $s'$  at least one operation from some block  $B$  of  $G(s)$  has to be processed before the first or after the last operation of  $B$ .

Consequence:  $s, s'$  with  $L(s') < L(s)$ , then one of the following holds

- at least one operation of one block  $B$  in  $G(s)$ , different from the first operation in  $B$ , has to be processed *before all other* operations of  $B$  in the schedule given by  $G(s')$
- at least one operation of one block  $B$  in  $G(s)$ , different from the last operation in  $B$ , has to be processed *after all other* operations of  $B$  in the schedule given by  $G(s')$



Consider  $(u, v)$  on critical path w.r.t.  $s$ .

Disadvantage of  $\mathcal{N}_1$ :

- If  $(u, v)$  belong to a block, but do not contain first or last operation of this block, no improvement occurs.
- $\Rightarrow$  generally, several moves are needed to improve solution

Let  $(v, w)$  be processed on the same machine, and denote by  $PM(v)(SM(w))$  the immediate predecessor (successor) of  $v(w)$  (if exists).

Consider as moves all permutations of  $\{PM(v), v, w\}$  and  $\{v, w, SM(w)\}$  where  $(v, w)$  is reversed and that are feasible ( $\mathcal{N}_2$ ).

Clearly,  $\mathcal{N}_1 \subseteq \mathcal{N}_2 \Rightarrow \mathcal{N}_2$  is OPT-connected

Directly using the block-lemma, we obtain:

$\mathcal{N}_3$  is defined as the neighborhood, where operations of a block are shifted at the beginning or the end of the respective block.

open question: Is  $\mathcal{N}_3$  OPT-connected?

$\mathcal{N}_3$  can be extended to a neighborhood  $\mathcal{N}_4$  which is OPT-connected in the following way:

- Let  $P$  be a critical path in  $G(s)$ .
- $s'$  is derived from  $s$  by moving one operation  $j$  of a block  $B$  of  $P$  different from the first (last) operation in  $B$  before (after) all operations of  $B$  (if feasible).
- Otherwise (i.e. above not feasible),  $j$  is moved to the position inside  $B$  closest to first (last) operation, that is still feasible.

Note:  $\mathcal{N}_3 \subseteq \mathcal{N}_4$

Lemma:  $\mathcal{N}_4$  is OPT-connected  
(proof on the board)

# Organization of the Tabu-List

$\mathcal{N}_1$  up to  $\mathcal{N}_4$  work by reversing an arc  $(v, w)$  in  $G(s)$ :

- attribute = arc reversed by recent moves
- a solution is defined to be tabu if an arc belonging to the attribute set is contained in it

As supporting data-structure, we use a matrix  $A = (a_{ij})$ :

- $a_{ij}$  = count of the iteration in which arc  $(i, j)$  was last reversed
- we forbid a swap of  $(i, j)$  if the count + length of the tabu-list is greater than the current iteration

$\Rightarrow$  the tabu-list length can be arbitrarily chosen  
(memory does not increase with length)

On  $A$ , we use a dynamic tabu-list management:

- improving phase: *decrease* length of list
- non-improving phase: *increase* length of list

also include an aspiration criterion, e.g. based on a lower bound

## *Genetic Algorithm*

- general search technique inspired by *biological evolution*
  - ' survival of the fittest '
- work on a set *POP* of solutions (*population*)
  - instead of single solution as in local search
- single solution  $s \in POP$  is called *chromosome*
  - usually encoded by a sequence of symbols (DNA)
- for each feasible solution  $s$ ,  $fit(s)$  is a measure of adaption (*fitness value*)
  - $fit(s)$  is often related to the objective function  $f(s)$

Starting from an initial population, 'parent' solutions are selected and new 'child' solutions are created by genetic operators:

- *crossover*
  - mix subsequences of parent chromosomes
- *mutation*
  - perturbate a chromosome

Size of the population is controlled by fitness value.

## Algo.: *Genetic Algorithm*

- ① Generate initial population  $POP$
- ② Compute fitness of each individual  $s \in POP$
- ③ REPEAT
  - ① Choose two parent solutions  $s^M, s^F \in POP$
  - ② Create a child solution  $s^C$  from  $s^M, s^F$  by crossover
  - ③ Mutate  $s^C$  with certain probability
  - ④ Compute fitness of  $s^C$
  - ⑤ Add  $s^C$  to  $POP$  and reduce  $POP$  by selection
- ④ UNTIL some stopping criterion is satisfied

Different variations are also possible

- several children may be generated simultaneously
- population may be divided into two sets
  - matching and creation of two new children → new population
- mutations may be realized by local search

In this approach, also infeasible solutions (chromosomes) may be present, this can be encoded in the fitness value.