

Chapter 5

Basic Algorithms for Allocation Problems

5.1 Heuristic Search

- Most of the allocation problems mentioned in chapter 4 are combinatorial problems.
- Therefore, they belong to the class of **discrete optimization problems**:

$$\varphi(x_1, x_2, \dots, x_n) \rightarrow \min$$

$$\text{with } \sum_{j=1}^n x_j = C; \quad x_j \in N, j = 1, \dots, n$$

- In this general form they are NP-hard.
- Thus, for many practical problems optimal solutions cannot be found in acceptable time.

- A heuristic search is a smart exploration of parts of the solution space.
- Concentration on promising areas
- No guarantee of optimality
- Compromise between **effectiveness** (how good is the solution found) and **efficiency** (how fast do we find the solution)

1	<code>solution ← initial_solution</code>	initialization
2	<code>finished ← false</code>	
3	while not finished do	main loop
4	<code>new_solution ← modify(solution)</code>	search step
5	<code>delta_phi ← phi(new_solution) - phi(solution)</code>	evaluation
6	if <code>delta_phi < 0</code>	improvements only
7	<code>then solution ← new_solution</code>	acceptance
9	<code>finished ← f(?)</code>	termination criterion
10	end while	
11	end	

Problems:

- What are search steps like?
- What are elementary search steps?
- How can we find at least a local minimum?

Digression: Combinatorics

- Combinatorics is the study of collections of objects. Specifically, **counting** objects, **arrangement**, **derangement**, etc. along with their mathematical properties.
- Counting objects is important in order to analyze algorithms and to compute discrete probabilities.
- Originally, combinatorics was motivated by gambling: counting configurations is essential to elementary probability.
 - Example: How many arrangements of a deck of 52 cards are possible?

Permutations

- A **permutation** of a set of distinct objects is an ordered arrangement of these objects.
- An ordered arrangement of r elements of a set of n elements is called an **r -permutation**
- The number of r permutations of a set of n distinct elements is

$$P(n, r) = \prod_{i=0}^{r-1} (n - i) = n(n - 1)(n - 2) \cdots (n - r + 1)$$

- It follows that $P(n, r) = \frac{n!}{(n - r)!}$
- In particular $P(n, n) = n!$
- Note here that **the order is important**. It is necessary to distinguish when the order matters and when it does not.

- How many pairs of dance partners can be selected from a group of 12 women and 20 men?
 - The first woman can partner with any of the 20 men, the second with any of the remaining 19, etc.
 - To partner all 12 women, we have

$$P(20,12) = 20!/8! = 20 \times 19 \times 18 \times \dots \times 10 \times 9$$

- Whereas permutations consider order, **combinations** are used when order does not matter.
- **Definition:** A **k -combination** of elements of a set is an **unordered** selection of k elements from the set.
 - (A combination is simply a subset of cardinality k .)
- The number of k -combinations of a set of cardinality n with $0 \leq k \leq n$ is

$$C(n, k) = \binom{n}{k} = \frac{n!}{(n-k)!k!}$$

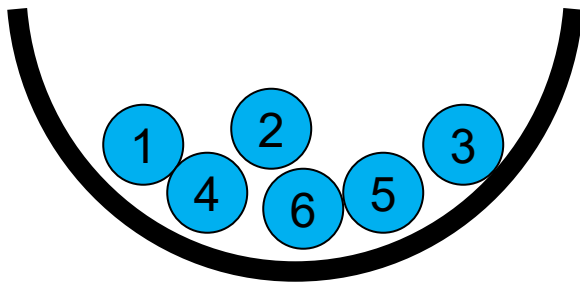
It is read ' n choose k '.

- A useful fact about combinations is that they are symmetric:

$$\binom{n}{k} = \binom{n}{n-k}$$

Ball-in-urn experiment

- Given n balls in an urn or bowl. m times a ball is taken out.
- Question: How many different possibilities exist to take out a ball m times from n balls?



Differentiation:

1. Is the ball put back each time or not?
2. Does it make a difference in which order the balls are removed from the bowl or not?

Combinatorics: Example 2 out of 4

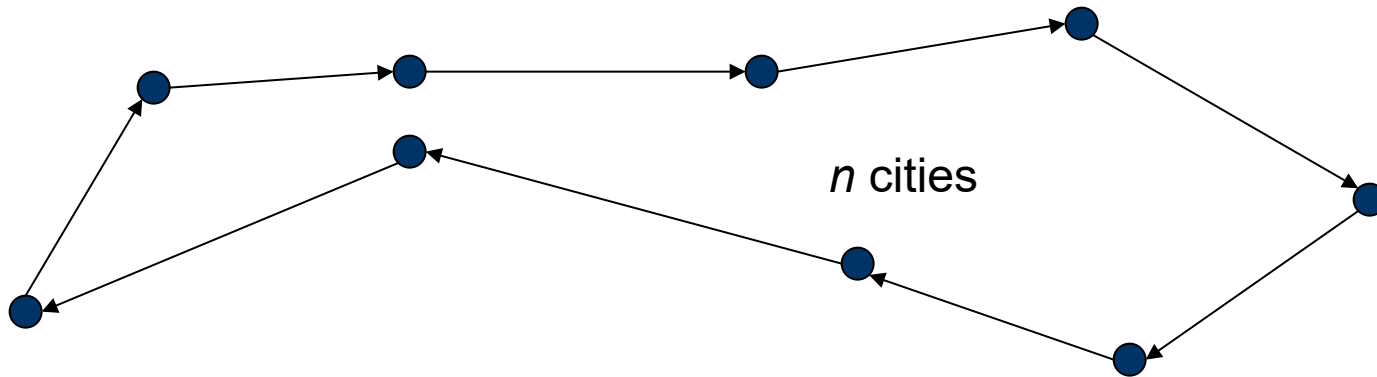
	With repetition	Without repetition
Order matters	$(1,1), (1,2), (1,3), (1,4)$ $(2,1), (2,2), (2,3), (2,4)$ $(3,1), (3,2), (3,3), (3,4)$ $(4,1), (4,2), (4,3), (4,4)$	$(1,2), (1,3), (1,4)$ $(2,1), (2,3), (2,4)$ $(3,1), (3,2), (3,4)$ $(4,1), (4,2), (4,3)$
Order does not matter	$(1,1), (1,2), (1,3), (1,4)$ $(2,2), (2,3), (2,4)$ $(3,3), (3,4)$ $(4,4)$	$(1,2), (1,3), (1,4)$ $(2,3), (2,4)$ $(3,4)$

	With repetition	Without repetition
With considering order	"k-Sample" n^k	"k-Permutation" $P(n, k) = \binom{n}{k} \cdot k! = \frac{n!}{(n-k)!}$
Without considering order	"k-Selection" $C(n+k-1, n-1) = \binom{n+k-1}{n-1}$ $= \frac{(n+k-1)!}{k!(n-1)!}$	"k-Combination" $C(n, k) = \binom{n}{k} = \frac{n!}{(n-k)! k!}$

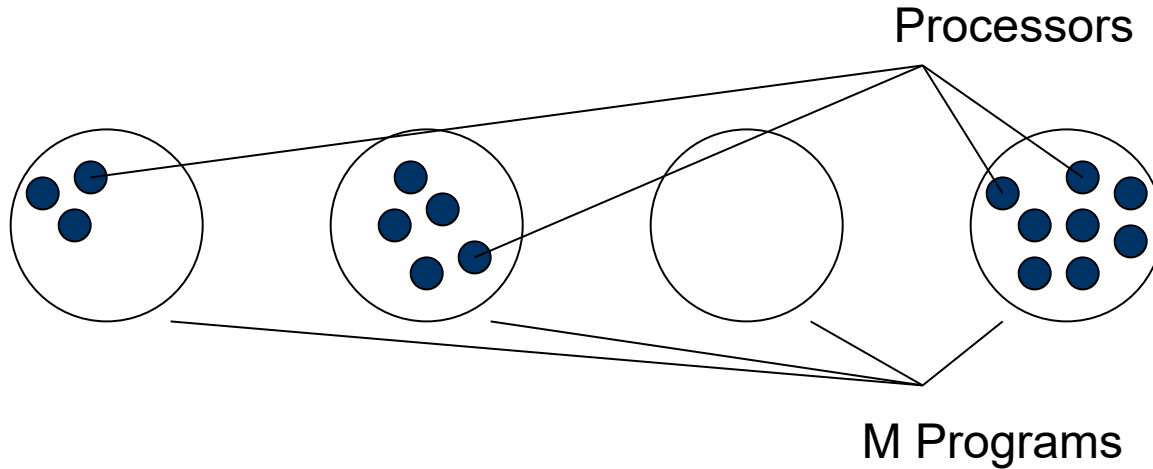
5.2 Representation

- For the general discussion of search problems, we represent a solution as a finite string from a finite alphabet.

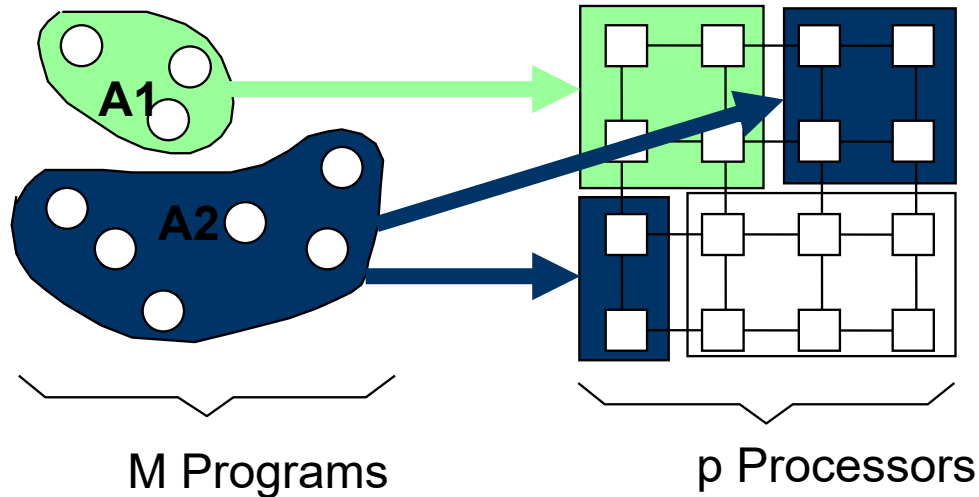
Example: Traveling Salesman Problem



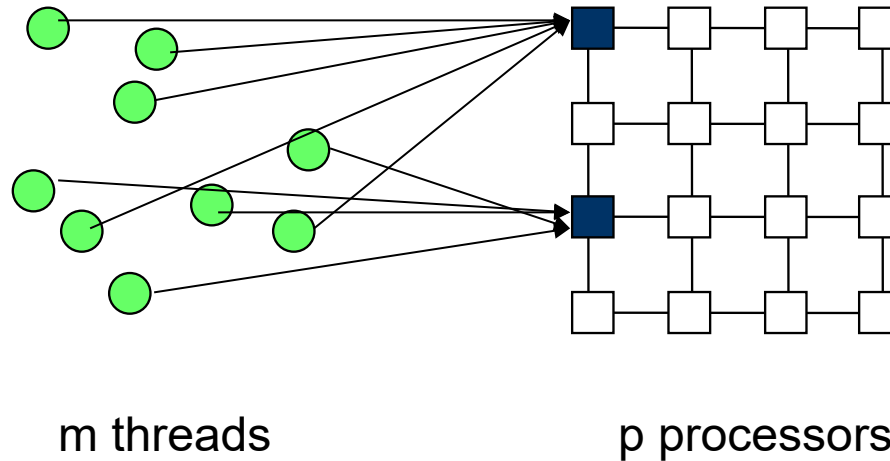
- Goal: Minimal round trip
- Coding: string of length n from the alphabet $\{1, \dots, n\}$ without repetition (permutation)
 $x_i = k$: city k is visited at i -th place.
- Size of solution space (permutation): $n!$



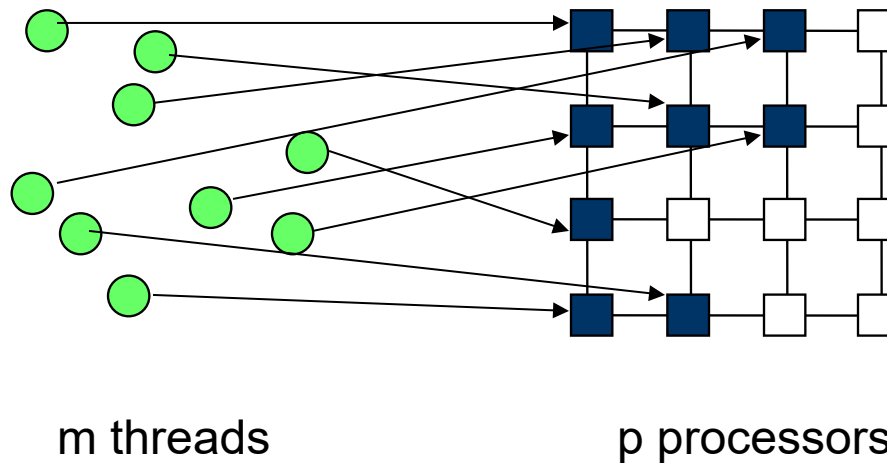
- Goal: Partitioning of p processors into M programs so that the accumulated running time is minimized
- Coding: String of size M of alphabet $\{0, 1, \dots, p\}$ with sum of digits = p
 $x_i = k$: program i is given k processors
- Size of solution space:
$$\binom{p + M - 1}{p} = \binom{p + M - 1}{M - 1}$$



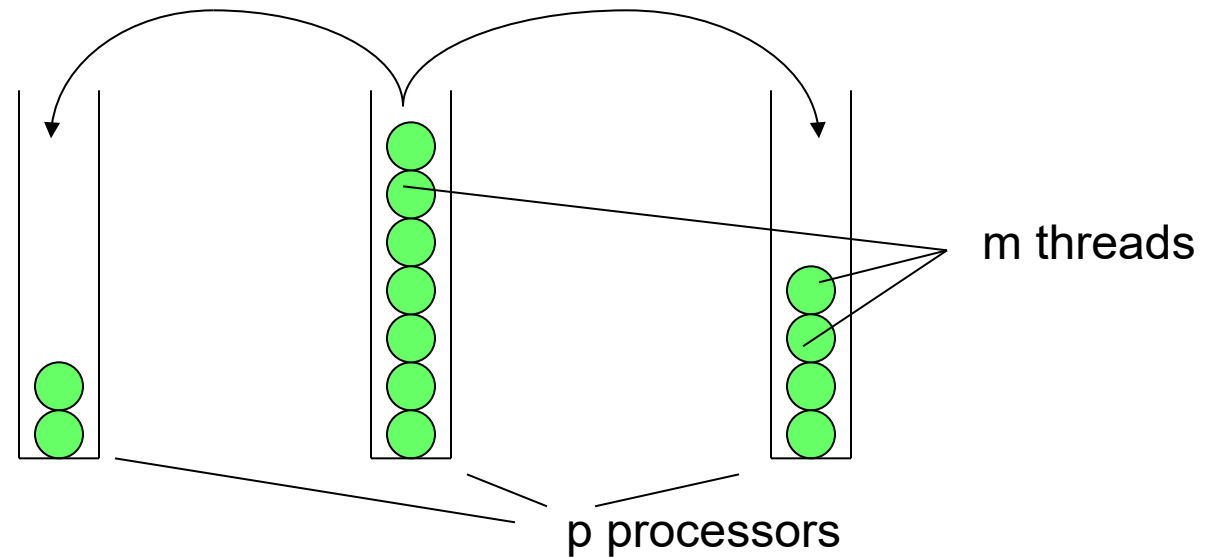
- Goal: Assignment of M programs to processor sets with minimal fragmentation and minimal interprocessor communication.
- Coding: string of size p of alphabet $\{0, 1, \dots, M\}$ with repetition
 $x_i = k$: processor i is occupied by program k
- Size of solution space: $(M + 1)^p$



- Goal: Mapping of threads to processors with minimal interprocessor communication and balanced load.
- Coding: String of size m of alphabet $\{1, \dots, p\}$ with repetition
 $x_i = k$: thread i is assigned to processor k
- Size of solution space: p^m



- Goal: Injective mapping of m threads to p processors with *minimal* communication cost.
- Coding: String of size p of alphabet $\{0, 1, \dots, m\}$ without repetition
 $x_i = k$: processor i executes thread k
- Size of solution space: $\binom{p}{m} m!$



- Goal: Balanced distribution of m threads across p processors
- Coding: String of size p of alphabet $\{0, 1, \dots, m\}$ with sum of digits of m
 $x_i = k$: processor i receives a load of k threads
- Size of solution space: (same problem as slide 13) $\binom{m+p-1}{m-1} = \binom{m+p-1}{p}$

Intermediate result

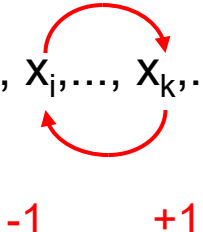
- A solution is represented by a string of size n .
- The set of feasible solutions spans the solution space.
- Solutions are points in an n -dimensional solution space.
- Associated with each solution x is the value $\Phi(x)$ of the objective function to be optimized.
- Search algorithms scan the solutions space and try to find a solution with a very high (or small) value Φ .

Elementary search steps

- A search step is a (feasible) modification of the string representing a solution.
- Elementary search steps (examples):

- Local value change $x_1, x_2, x_3, \dots, x_i \pm \Delta, \dots, x_{n-1}, x_n$

- Exchange $x_1, x_2, x_3, \dots, x_i, \dots, x_k, \dots, x_{n-1}, x_n$

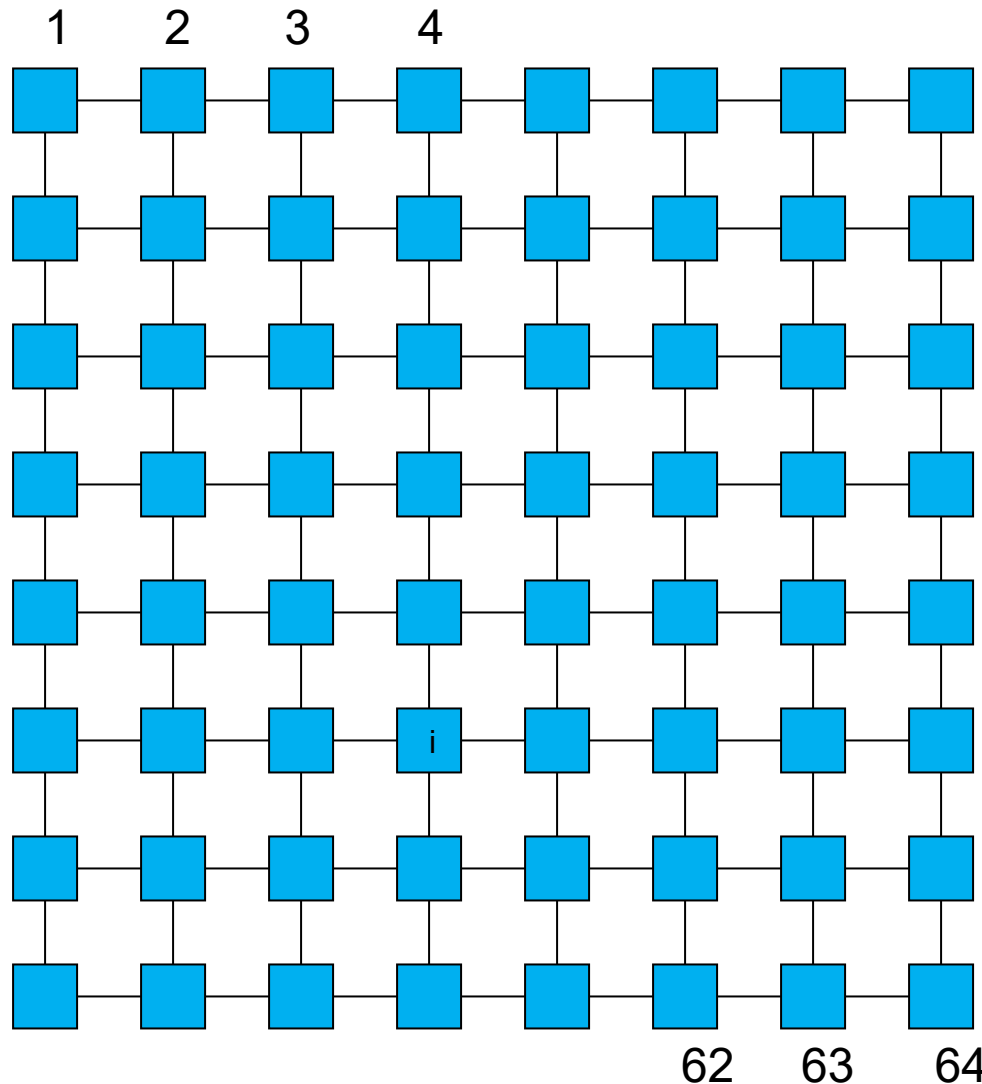


- Increment shift $x_1, x_2, x_3, \dots, x_i, \dots, x_k, \dots, x_{n-1}, x_n$

- Problem: Between the positions and between the characters distances have to be defined.

- Depending on the problem and the algorithm used, elementary steps may be defined differently.
- The neighborhood of a point (a solution) denotes all points that can be reached with one elementary step.
- $\text{neighbors}(x) := \{x' \mid \exists \text{ elementary step } x \rightarrow x'\}$
- Usually, we have symmetry:
$$x \in \text{neighbors}(y) \Leftrightarrow y \in \text{neighbors}(x)$$
- Symmetry implies that search steps are reversible.

Example: Contractive allocation of m threads on a 8x8-mesh



e.g. $x_i = 44$:

thread i is allocated to processor 44

Elementary step:

local:

move to adjacent processor

$$|\text{neighbors}(x)| = 4 \cdot m$$

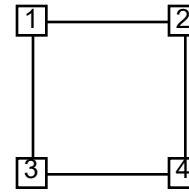
global:

move to any place

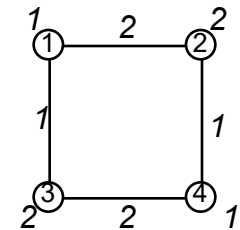
$$|\text{neighbors}(x)| = 63 \cdot m$$

Example

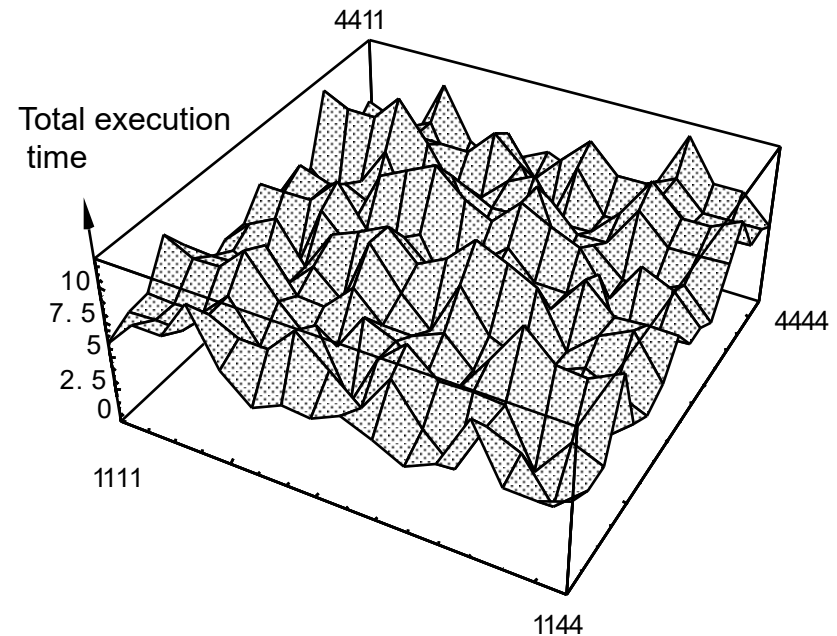
- A program consisting of 4 threads has to be assigned to a ring of 4 processors contractively.
- Objective function is the total execution time.
- Each solution is a 4-tuple (x_1, x_2, x_3, x_4) with $x_i = k$: thread i is assigned to processor k .



machine

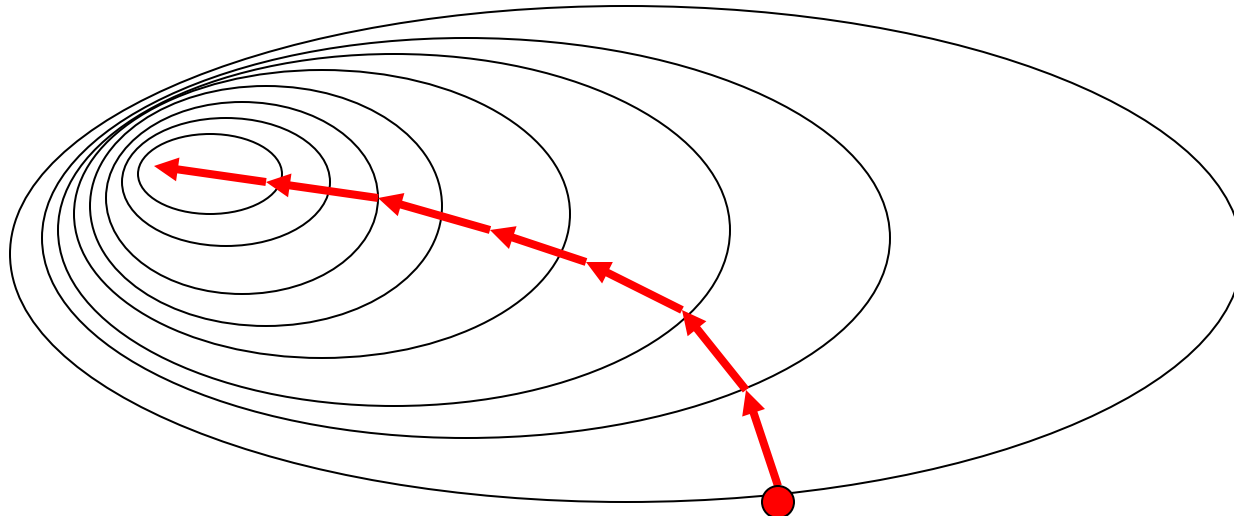


program



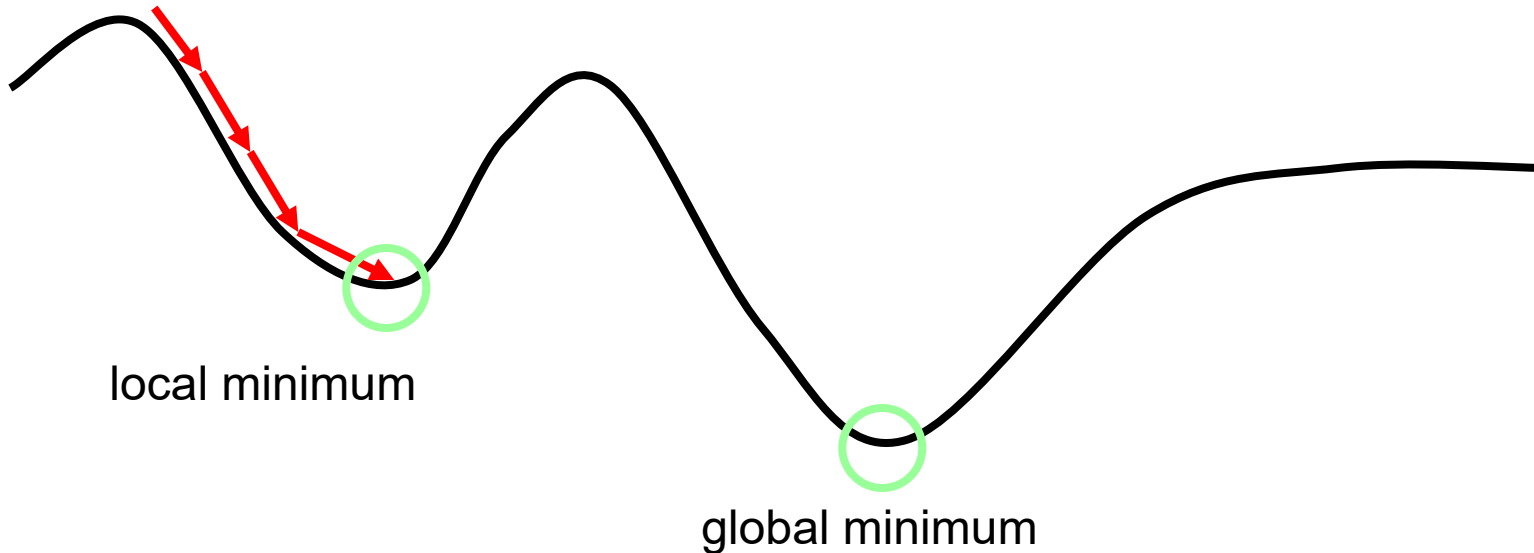
5.3 Gradient descent

- The gradient descent is an improved variant of the local search.
- Instead of performing an arbitrary elementary step, all possible elementary steps are evaluated (objective function) and that step with the largest gain is selected.
- We therefore proceed in the direction of the steepest slope (direction of gradient).



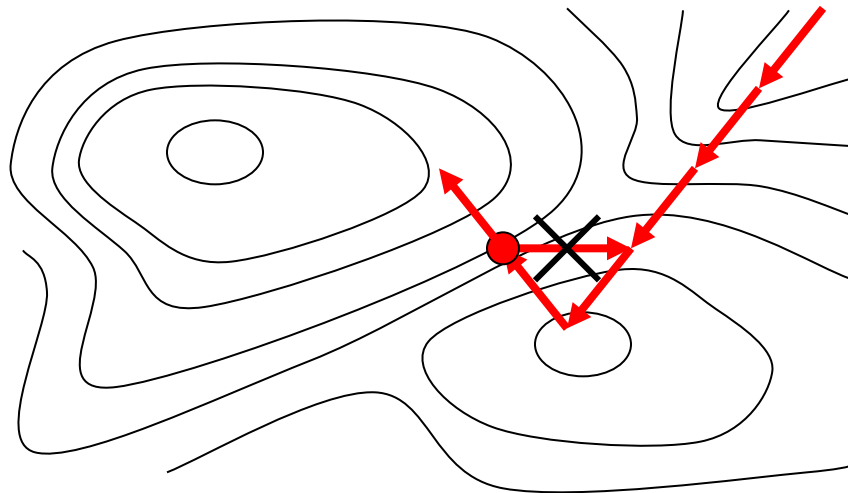
1	<code>solution ← initial solution</code>	initialization
2	<code>finished ← false</code>	
3	<code>while not finished do</code>	main loop
4	<code>for all neighbors(solution) do</code>	
5	<code>calculate phi(neighbor)</code>	test steps
6	<code>end for</code>	
7	<code>new_solution ← neighbor with min. phi(neighbor)</code>	gradient descent
8	<code>delta_phi ← phi(new_solution) - phi(solution)</code>	evaluation
9	<code>if delta_phi < 0</code>	
10	<code>then solution ← new_solution</code>	descent (improvement)
11	<code>else finished ← true</code>	local minimum
12	<code>end while</code>	
13	<code>end</code>	

- Local search algorithms exhibit the fundamental problem that they do find a local minimum but cannot escape it.
- Many heuristic approaches can be distinguished by the way they are able to leave local minima.



5.4 Taboo search

- Taboo search is another improved variant of local search.
- Occasionally also deteriorations are accepted.
- In contrast to the gradient descent it has a “memory”:
 - Best solution found so far
 - List of forbidden steps (taboo), e.g. to avoid cycles

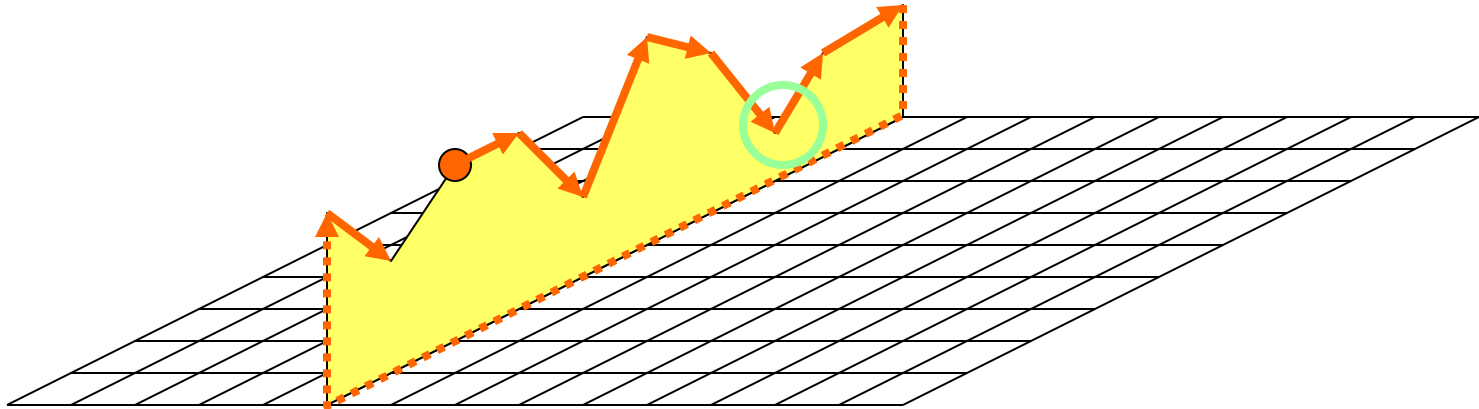


Taboo search (simplified)

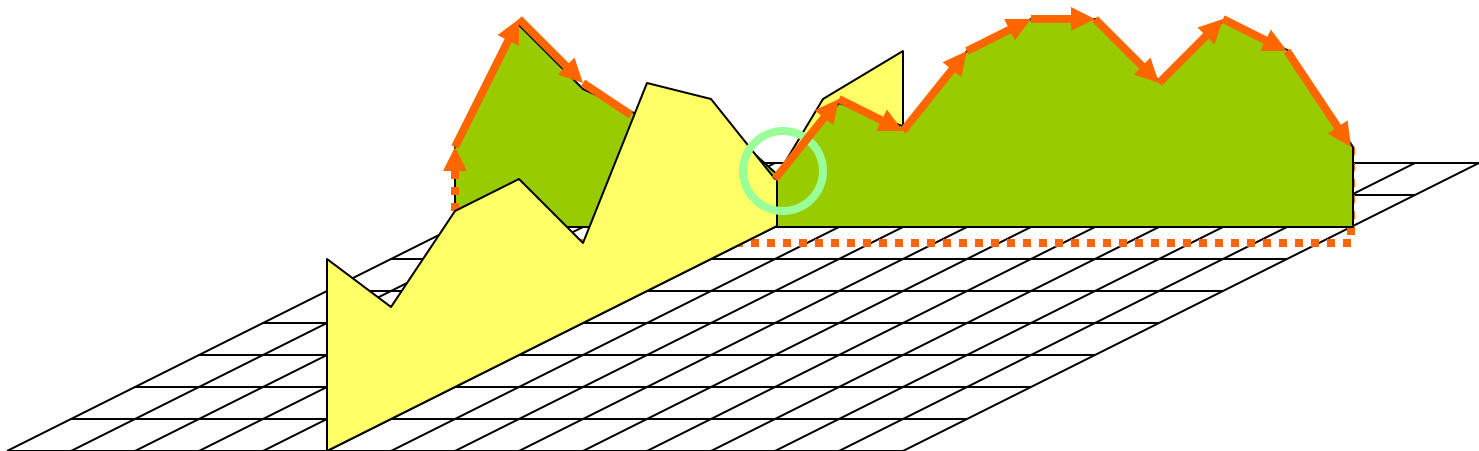
1	<code>solution ← initial solution</code>	initialization
2	<code>best_solution ← solution</code>	
3	<code>taboo_list ← {solution}</code>	
4	<code>finished ← false</code>	
5	while not finished do	main loop
6	for all neighbors(solution) do	
7	<code>calculate phi(neighbor)</code>	test steps
8	end for	
9	<code>solution ← neighbor with min. phi(neighbor) and neighbor ∉ tabulist</code>	calculate new solution
10	<code>tabulist ← tabulist ∪ {solution}</code>	update
11	if <code>phi(solution) < phi(best_solution)</code>	
12	then <code>best_solution ← solution</code>	store best solution
13	<code>finished ← f(?)</code>	termination criterion
14	end while	
15	end	

5.5 Probing paths

- Another idea to escape from local minima is to run a complete path (e.g. along some dimension) and only then accept the minimum of the path as the next solution.



- Starting at the minimum found we can then proceed with a new probing path along a new dimension.



Probing paths

1	<code>solution ← initial solution</code>	initialization
2	<code>finished ← false</code>	
3	while not finished do	main loop
4	<code>select probing_path</code>	
5	<code>i ← 0</code>	
6	<code>step[0] ← solution</code>	
7	while not (end_of_path) do	test steps
8	<code>step[i] ← next(step[i-1])</code>	
9	<code>phi_s[i] ← phi(step(i))</code>	store values of objective function
10	end while	
11	<code>i ← minimum(phi_s)</code>	index of minimum along path
12	<code>new_solution ← step[i]</code>	gradient descent
11	<code>delta_phi ← phi(new_solution) - phi(solution)</code>	evaluation
12	if <code>delta_phi < 0</code>	
13	then <code>solution ← new_solution</code>	descent (improvement)
14	else <code>finished ← true</code>	local minimum
15	end while	

5.6 Simulated Annealing (SA)

- Idea 1: Accept steps „uphill“ with some probability.
- Idea 2: At the beginning search in large areas and then gradually restrict the scope of the search.
- Approach: Prob(accept deterioration by $\Delta\varphi$) = $\exp(-\Delta\varphi/T)$
- T controls the “acceptance of deterioration” and is reduced stepwise e.g.:
 - logarithmic decrement: $T(i) := T_0 / \log(i), i = 1, 2, \dots$
 - geometric decrement: $T(i) := T_0 / q^i (q > 1), i = 1, 2, \dots$
 - linear decrement: $T(i) := T_0 (1 - i/m), i = 0, 1, 2, \dots, m$
- **Remark:** Idea goes back to computer simulation of cooling down procedures (annealing) of specific material (spin glasses), that are fluid at high temperature (high mobility of molecules) and by careful annealing achieve a homogeneous grid structure at minimum energy.

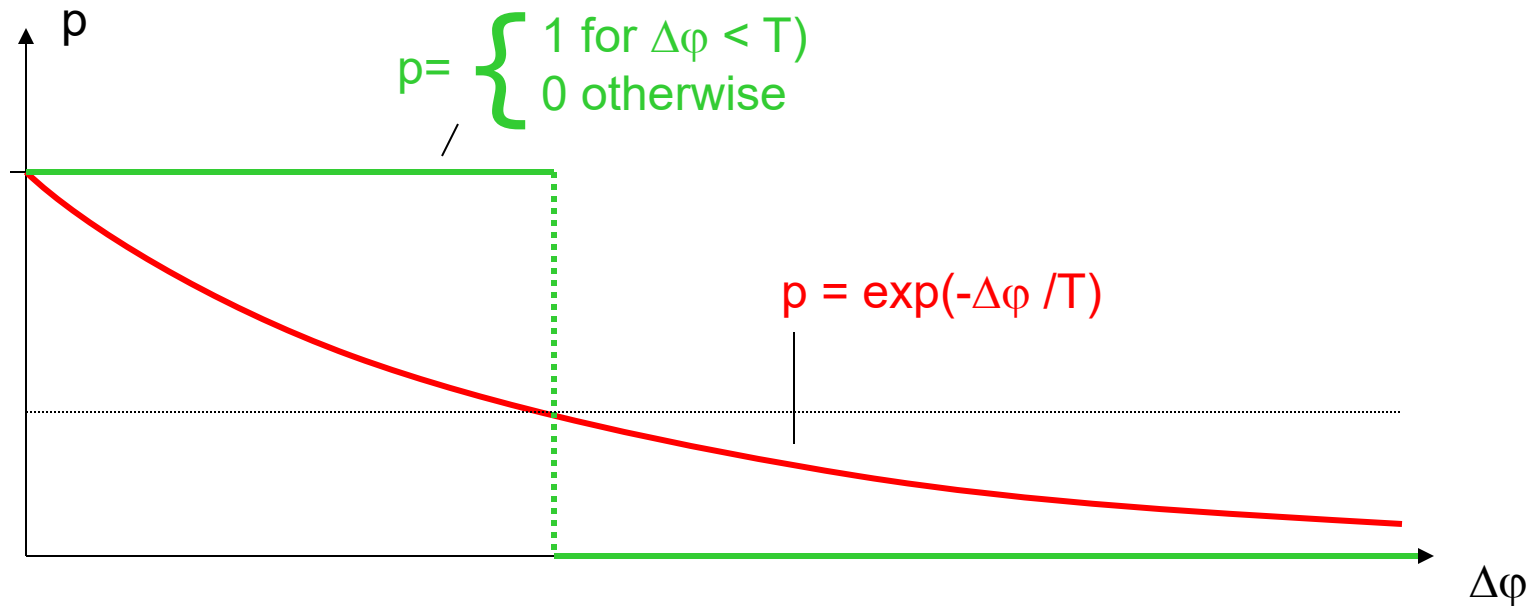
Simulated Annealing (Variant)

Also here it is rewarding to save the best solution found so far

1	<code>solution ← random initial solution</code>	
2	<code>T ← T0</code>	initialize threshold
3	<code>best_solution ← solution</code>	
4	<code>for i ← 0 to m do</code>	m=number of stages
5	<code>for j ← 1 to n do</code>	n =number of steps per stage
6	<code>new_solution ← modify(solution)</code>	
7	<code>delta_phi ← phi(new_solution) - phi(solution)</code>	
8	<code>if phi(new_solution) < phi(best_solution)</code>	
9	<code>then best_solution ← new_solution</code>	save new best solution
10	<code>if delta_phi < 0</code>	
11	<code>then solution ← new_solution</code>	improvement
12	<code>else with prob exp(-delta_phi/T)</code>	deterioration
13	<code>solution ← new_solution</code>	
14	<code>end for</code>	
15	<code>T ← decrement(T)</code>	decrement threshold
16	<code>end for</code>	

Threshold Accepting

- Experimental insight: acceptance can be controlled deterministically without loss of solution quality:
Accept, if $\Delta\varphi < T$
- Results in a variant of Simulated Annealing.



Deluge Algorithm

Further variant: Instead of relative comparing with previous solution, compare with best solution

1	<code>solution ← random initial solution</code>	
2	<code>T ← T0</code>	initialize threshold
3	<code>best_solution ← solution</code>	
4	<code>for i ← 0 to m do</code>	m=number of stages
5	<code>for j ← 1 to n do</code>	n =number of steps per stage
6	<code>new_solution ← modify(solution)</code>	
7	<code>delta_phi ← phi(new_solution) - phi(solution)</code>	
8	<code>if phi(new_solution) < phi(best_solution)</code>	
9	<code>then best_solution ← new_solution</code>	save best solution
10	<code>if phi(new_solution) < phi(best_solution) + T</code>	maximally by T worse than best solution
11	<code>then solution ← new_solution</code>	acceptance
12	<code>else solution ← best_solution</code>	variant: backtracking
13	<code>end for</code>	
14	<code>T ← decrement(T)</code>	reduce threshold
15	<code>end for</code>	

5.7 Iteration

- The solution found by a heuristic search algorithm depends on the initial solution.
- Starting with another initial value we may find another, better final solution.
- Thus: Iteration across different initial solutions.

1	<code>best_solution ← initial solution</code>	initialization
2	<code>for i ← 1 to n do</code>	iteration
3	<code>solution ← new_initial solution</code>	new initialization
4	{heuristic search}	delivers loc. minimum (solution)
5	<code>if phi(best_solution) > phi(solution)</code>	
6	<code>then best_solution ← solution</code>	improvement
7	<code>end for</code>	
8	<code>end</code>	

5.8 Parallelism

- Since the particular iterations are independent of each other, they can be executed in parallel.
- Two variants:
 - Variant A
 - Disjoint partitioning of solution space
 - Heuristic search in each partition (subspace) in parallel
 - Calculating minimum after termination of all search activities
 - Variant B
 - No partitioning
 - All „minimum searcher“ search in the whole area
 - Calculating minimum after termination of all search activities

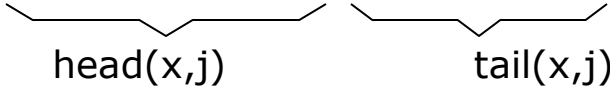
5.9 Combination of solutions

- Evolutionary algorithms:
- Idea: If we proceed with many solutions simultaneously, why not combine their „beneficial“ features.
- Reverting to Darwin's evolution strategy: crossover, mutation, „survival of the fittest“

1	<code>solution_set ← initial solution_set</code>	initialization
2	<code>finished ← false</code>	
3	while not finished do	main loop
4	<code>combination_set ← select(solution_set)</code>	survival of the fittest
5	<code>new_solution_set ← recombine(combination_set)</code>	cross over, mutation
15	<code>calculate phi(new_solution_set)</code>	evaluation
7	<code>finished ← f(?)</code>	termination criterion ?
8	end	

Genetic Algorithms (GA)

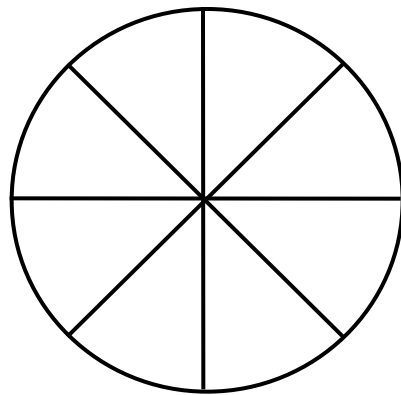
- Problem: How to find suitable operations?
 - How many solutions should be involved?
 - one: mutation
 - two: crossover
 - n: n-fold crossover
 - How do we combine or modify?
 - Permutation of strings
 - Exchange of substrings
 - Blend of substrings
- **Genetic Algorithms** are a subclass of Evolutionary Algorithms.
- They require a representation of solutions as binary strings of fixed length.
- They use two operators only :
 - binary mutation: flipping individual bits in solution string
 - binary crossover: combination of solution strings

- To explain the way GA work we need some auxiliary functions:
 - **random_select(set):**
selects an element of a n -element set with equal probability ($p=1/n$).
 - **p_select(p, set):**
selects an element of a n -element set according to a given discrete probability density function p_i ($i = 1, \dots, n$).
 - **head(x,j)** and **tail(x,j)**
$$x_1, x_2, x_3, \dots, x_j, x_{j+1}, \dots, x_n$$


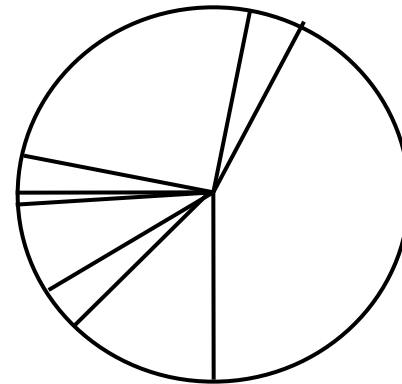
head(x,j) tail(x,j)
 - **mutation(x,j)**
Bit j in bit string x is inverted.

Comparison of the two selection procedures

- The two selection procedures can be regarded as turning the roulette wheel with the elements as sectors.
- With `random_select` all sectors are of the same size, with `p_select` the probabilities are proportional to the sector angles.



`random_select(M)`



`p_select(p, M)`

Genetic Algorithms: Example

1	<code>solution_set ← choose initial solution_set of size n</code>	initialization (maximization problem)
2	while not finished do	main loop (generations)
3	for <code>i ← 1 to n</code> do	all solution
4	<code>sp[i] ← $\phi[i] / \sum \phi[i]$</code>	selection probability <code>sp</code>
5	<code>mating ← ∅</code>	
6	for <code>i ← 1 to n</code> do	
7	<code>s ← p_select(sp, solution_set)</code>	random selection according to <code>sp</code>
8	<code>mating_pool ← mating_pool ∪ {s}</code>	
9	end for	
10	<code>mating_pairs ← ∅</code>	
11	while <code>mating_pool ≠ ∅</code> do	random mating
12	<code>s ← random_select[mating_pool]</code>	
13	<code>mating_pool ← mating_pool - {s}</code>	
14	<code>t ← random_select [mating_pool]</code>	
15	<code>mating_pool ← mating_pool - {t}</code>	
16	<code>mating_pairs ← mating_pairs ∪ {(s,t)}</code>	
17	end while	

Genetic Algorithms: Example (continued)

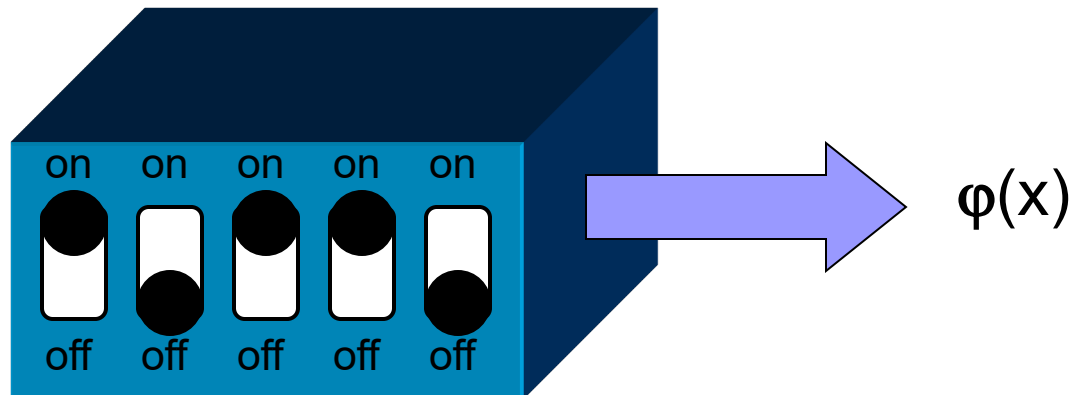
18	<code>new_solution_set $\leftarrow \emptyset$</code>	
19	<code>for all (x,y) \in mating_pairs do</code>	crossover
20	<code> j \leftarrow random_select({1,2,...,k})</code>	substring selection
21	<code> new_solution1 \leftarrow head(x,j)+tail(y,j)</code>	build generation of offsprings by
22	<code> new_solution2 \leftarrow head(y,j)+tail(x,j)</code>	crosswise exchange of substrings
23	<code> new_solution_set \leftarrow new_solution_set \cup {new_solution1,new_solution2}</code>	parents are replaced by offsprings
24	<code>end for</code>	
25	<code>for all s \in new_solution_set do</code>	
26	<code> with probability mp do</code>	modification probability mp
27	<code> j \leftarrow random_select({1,2,...,k})</code>	random feature selection
28	<code> modify(s,j)</code>	feature j is modified
29	<code> end with</code>	
30	<code>end for</code>	
31	<code>solution_set \leftarrow new_solution_set</code>	
32	<code>end while</code>	
33	<code>end GA</code>	

Example

$$\varphi(n) = n^2 \rightarrow \max ; \quad n \in \{0,1,2,\dots,31\}$$

Representation as a GA problem:

$$n = \sum_{i=0}^4 x_i 2^i \quad x_i \in \{0,1\}$$



Example (continued)

No.	Solution set	x	$\varphi(x)$	Selection probability	Selection	Mating pool after selection	No of mate (rand.)	Substring selection (random)	New solution set	x	$\varphi(x)$
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1. Iteration

1.	0 1 1 0 1	13	169	0.15	1	0 1 1 0 1	2	2	0 1 0 0 0	8	64
2.	1 1 0 0 0	24	576	0.51	2	1 1 0 0 0	1	2	1 1 1 0 1	29	841
3.	0 0 1 0 0	4	16	0.02	0	1 1 0 0 0	4	4	1 1 0 0 1	25	625
4.	1 0 0 1 1	19	361	0.32	1	1 0 0 1 1	3	4	1 0 0 1 0	18	324
S			1122								1854
Max			576								841

Example (continued)

No.	Solution set	x	$\varphi(x)$	Selection probability	Selection	Mating pool after selection	No of mate (rand.)	Substring selection (random)	New solution set	x	$\varphi(x)$
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2. Iteration

1.	0 1 0 0 0	8	64	0.03	0	1 1 1 0 1	3	3	1 1 1 0 1	29	841
2.	1 1 1 0 1	29	841	0.45	2	1 1 1 0 1	4	2	1 1 0 1 0	26	676
3.	1 1 0 0 1	25	625	0.34	1	1 1 0 0 1	1	3	1 1 0 0 1	25	625
4.	1 0 0 1 0	18	324	0.18	1	1 0 0 1 0	2	2	1 0 1 0 1	21	441
S			1854								2583
Max			841								841

Example (continued)

No.	Solution set	x	$\varphi(x)$	Selection probability	Selection	Mating pool after selection	No of mate (rand.)	Substring selection (random)	New solution set	x	$\varphi(x)$
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3. Iteration

1.	1 1 1 0 1	29	841	0.33	2	1 1 1 0 1	4	1	1 1 1 0 1	29	841
2.	1 1 0 1 0	26	676	0.26	1	1 1 0 1 0	3	3	1 1 0 0 1	25	625
3.	1 1 0 0 1	25	625	0.24	1	1 1 0 0 1	2	3	1 1 0 1 0	26	676
4.	1 0 1 0 1	21	441	0.17	0	1 1 1 0 1	1	1	1 1 1 0 1	29	841
S			2583								2983
Max			841								841

Example (continued)

No.	Solution set	x	$\varphi(x)$	Selection probability	Selection	Mating pool after selection	No of mate (rand.)	Substring selection (random)	New solution set	x	$\varphi(x)$
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4. Iteration

1.	1 1 1 0 1	29	841	0.28	2	1 1 1 0 1	2	1	1 1 1 0 1	29	841
2.	1 1 0 0 1	25	625	0.21	0	1 1 1 0 1	1	1	1 1 1 0 1	29	841
3.	1 1 0 1 0	26	676	0.23	1	1 1 0 1 0	3	3	1 1 0 0 1	25	625
4.	1 1 1 0 1	29	841	0.28	1	1 1 1 0 1	4	3	1 1 1 1 0	30	900
S			2983								3207
Max			841								900

Example (continued)

No.	Solution set	x	$\varphi(x)$	Selection probability	Selection	Mating pool after selection	No of mate (rand.)	Substring selection (random)	New solution set	x	$\varphi(x)$
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5. Iteration

1.	1 1 1 0 1	29	841	0.26	1	1 1 1 0 1	3	4	1 1 1 0 0	28	784
2.	1 1 1 0 1	29	841	0.26	1	1 1 1 0 1	4	3	1 1 1 1 0	30	900
3.	1 1 0 0 1	25	625	0.20	0	1 1 1 1 0	1	4	1 1 1 1 1	31	961
4.	1 1 1 1 0	30	900	0.28	2	1 1 1 1 0	2	3	1 1 1 0 1	29	841
S			3207								3486
Max			900								961

Example (continued)

No.	Solution set	x	$\varphi(x)$	Selection probability	Selection	Mating pool after selection	No of mate (rand.)	Substring selection (random)	New solution set	x	$\varphi(x)$
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6. Iteration

1.	1 1 1 0 0	28	784	0.22	0	1 1 1 1 1	3	2	1 1 1 1 1	31	961
2.	1 1 1 1 0	30	900	0.26	1	1 1 1 1 0	4	4	1 1 1 1 1	31	961
3.	1 1 1 1 1	31	961	0.28	2	1 1 1 1 1	1	2	1 1 1 1 1	31	961
4.	1 1 1 0 1	29	841	0.24	1	1 1 1 0 1	2	4	1 1 1 0 0	28	784
S			3486								3667
Max			961								961

Summary of Techniques

- Proceeding along the most promising direction (gradient descent)
- Memory, e.g., prevention of cycles (Taboo search)
- Evaluation of sequence of steps instead of single steps (probing paths)
- Acceptance of deterioration (Simulated Annealing)
- Backtracking
- Iteration with new initial solution
- Search in parallel
- Combination of solutions (Genetic Algorithms)

Further references

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