

DM811
Heuristics for Combinatorial Optimization

Lecture 8
Local Search (cntd.)

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Summary: Local Search Algorithms

(as in [Hoos, Stützle, 2005])

For given problem instance π :

1. search space S_π
2. neighborhood relation $\mathcal{N}_\pi \subseteq S_\pi \times S_\pi$
3. evaluation function $f_\pi : S \rightarrow \mathbf{R}$
4. set of memory states M_π
5. initialization function $\text{init} : \emptyset \rightarrow S_\pi \times M_\pi$
6. step function $\text{step} : S_\pi \times M_\pi \rightarrow S_\pi \times M_\pi$
7. termination predicate $\text{terminate} : S_\pi \times M_\pi \rightarrow \{\top, \perp\}$

1. Local Search Revisited
Components

2. Examples

Search Space

Defined by the solution representation:

- permutations
 - linear (scheduling)
 - circular (TSP)
- arrays (assignment problems: GCP)
- sets or lists (partition problems: graph partitioning, max indep. set)

Neighborhood function

Also defined as: $\mathcal{N} : S \times S \rightarrow \{T, F\}$ or $\mathcal{N} \subseteq S \times S$

- neighborhood (set) of candidate solution s : $N(s) := \{s' \in S \mid \mathcal{N}(s, s')\}$
- neighborhood size is $|N(s)|$
- neighborhood is symmetric if: $s' \in N(s) \rightarrow s \in N(s')$
- neighborhood graph of (S, N, π) is a directed graph: $G_{\mathcal{N}, \pi} := (V, A)$
with $V = S_\pi$ and $(uv) \in A \Leftrightarrow v \in N(u)$
(if symmetric neighborhood \rightsquigarrow undirected graph)

Notation: N when set, \mathcal{N} when collection of sets or function

A neighborhood function is also defined by means of an operator.

An operator Δ is a collection of operator functions $\delta : S \rightarrow S$ such that

$$s' \in N(s) \implies \exists \delta \in \Delta, \delta(s) = s'$$

Definition

k -exchange neighborhood: candidate solutions s, s' are neighbors iff s differs from s' in at most k solution components

Examples:

- 1-exchange (flip) neighborhood for SAT
(solution components = single variable assignments)
- 2-exchange neighborhood for TSP
(solution components = edges in given graph)

Definition:

- **Local minimum:** search position without improving neighbors wrt given evaluation function f and neighborhood \mathcal{N} ,
i.e., position $s \in S$ such that $f(s) \leq f(s')$ for all $s' \in N(s)$.
- **Strict local minimum:** search position $s \in S$ such that
 $f(s) < f(s')$ for all $s' \in N(s)$.
- *Local maxima* and *strict local maxima*: defined analogously.

Evaluation (or cost) function:

- function $f_\pi : S_\pi \rightarrow \mathbb{Q}$ that maps candidate solutions of a given problem instance π onto rational numbers (most often integer), such that global optima correspond to solutions of π ;
- used for assessing or ranking neighbors of current search position to provide guidance to search process.

Evaluation vs objective functions:

- *Evaluation function*: part of LS algorithm.
- *Objective function*: integral part of optimization problem.
- Some LS methods use evaluation functions different from given objective function (e.g., guided local search).

Constraint-based local search

From [B4]

What is a violation?

Constraint specific:

- decomposition-based violations
number of violated constraints, eg: alldiff
- variable-based violations
min number of variables that must be changed to satisfy c .
- value-based violations
for constraints on number of occurrences of values
- arithmetic violations
- combinations of these

Constraint-based local search

From [B4]

Local Search Revisited
Examples

Arithmetic constraints

- $l \leq r \rightsquigarrow \text{viol} = \max(l - r, 0)$
- $l = r \rightsquigarrow \text{viol} = |l - r|$
- $l \neq r \rightsquigarrow \text{viol} = 1$ if $l = r$, 0 otherwise

Combinatorial constraints

- **alldiff**(x_1, \dots, x_n):
Let a be an assignment with values $V = \{a(x_1), \dots, a(x_n)\}$ and $c_v = \#_a(v, x)$ be the number of variables with the same value. Possible definitions for violations are:
 - $\text{viol} = \sum_{v \in V} I(\max(c_v - 1, 0) > 0)$ value-based
 - $\text{viol} = \max_{v \in V} \max(c_v - 1, 0)$ value-based
 - $\text{viol} = \sum_{v \in V} \max(c_v - 1, 0)$ value-based
 - # variables with same value, variable-based, here leads to same definitions as previous three

Note:

- Local search implements a **walk** through the neighborhood graph
- Procedural versions of **init**, **step** and **terminate** implement sampling from respective probability distributions.
- Local search algorithms can be described as **Markov processes**:
behavior in any **search state** $\{s, m\}$ depends only
on current position s
higher order MP if (limited) memory m .

LS Algorithm Components

Search step (or **move**):

pair of search positions s, s' for which

s' can be reached from s in one step, i.e., $\mathcal{N}(s, s')$ and

$\text{step}(\{s, m\}, \{s', m'\}) > 0$ for some memory states $m, m' \in M$.

- **Search trajectory**: finite sequence of search positions $\langle s_0, s_1, \dots, s_k \rangle$ such that (s_{i-1}, s_i) is a *search step* for any $i \in \{1, \dots, k\}$ and the probability of initializing the search at s_0 is greater than zero, i.e., $\text{init}(\{s_0, m\}) > 0$ for some memory state $m \in M$.
- **Search strategy**: specified by `init` and `step` function; to some extent independent of problem instance and other components of LS algorithm.
 - random
 - based on evaluation function
 - based on memory

1. Local Search Revisited
Components
2. Examples

Iterative Improvement

Resume

- does not use memory
- **init**: uniform random choice from S or construction heuristic
- **step**: uniform random choice from improving neighbors

$$\Pr(s, s') = \begin{cases} 1/|I(s)| & \text{if } s' \in I(s) \\ 0 & \text{otherwise} \end{cases}$$

where $I(s) := \{s' \in S \mid \mathcal{N}(s, s') \text{ and } f(s') < f(s)\}$

- terminates when no improving neighbor available

Note: Iterative improvement is also known as iterative descent or hill-climbing.

Pivoting rule decides which neighbors go in $I(s)$

- **Best Improvement** (aka *gradient descent*, *steepest descent*, *greedy hill-climbing*): Choose maximally improving neighbors, i.e., $I(s) := \{s' \in N(s) \mid f(s') = g^*\}$, where $g^* := \min\{f(s') \mid s' \in N(s)\}$.

Note: Requires evaluation of all neighbors in each step!

- **First Improvement:** Evaluate neighbors in fixed order, choose first improving one encountered.

Note: Can be more efficient than Best Improvement but not in the worst case; order of evaluation can impact performance.

Examples

Iterative Improvement for SAT

- **search space S** : set of all truth assignments to variables in given formula F (solution set S' : set of all models of F)
- **neighborhood relation \mathcal{N}** : 1-flip neighborhood
- **memory**: not used, i.e., $M := \{0\}$
- **initialization**: uniform random choice from S , i.e., $\text{init}(\emptyset, \{a\}) := 1/|S|$ for all assignments a
- **evaluation function**: $f(a) :=$ number of clauses in F that are *unsatisfied* under assignment a (Note: $f(a) = 0$ iff a is a model of F .)
- **step function**: uniform random choice from improving neighbors, i.e., $\text{step}(a, a') := 1/|I(a)|$ if $a' \in I(a)$, and 0 otherwise, where $I(a) := \{a' \mid \mathcal{N}(a, a') \wedge f(a') < f(a)\}$
- **termination**: when no improving neighbor is available i.e., $\text{terminate}(a, \top) := 1$ if $I(a) = \emptyset$, and 0 otherwise.

Examples

Random order first improvement for SAT

URW-for-SAT($F, \text{maxSteps}$)

input: propositional formula F , integer maxSteps

output: a model for F or \emptyset

choose assignment φ of truth values to all variables in F
uniformly at random;

$\text{steps} := 0$;

while $\neg(\varphi$ satisfies $F)$ and $(\text{steps} < \text{maxSteps})$ **do**

 | select x uniformly at random from $\{x' \mid x' \text{ is a variable in } F \text{ and}$
 | changing value of x' in φ decreases the number of unsatisfied clauses}
 | $\text{steps} := \text{steps} + 1$;

if φ satisfies F **then**

 | **return** φ

else

 | **return** \emptyset

queensLS00.co

```
import cotls;
int n = 16;
range Size = 1..n;
UniformDistribution distr(Size);

Solver<LS> m();
var{int} queen[Size](m,Size) := distr.get();
ConstraintSystem<LS> S(m);

S.post(alldifferent(queen));
S.post(alldifferent(all(i in Size) queen[i] + i));
S.post(alldifferent(all(i in Size) queen[i] - i));
m.close();

int it = 0;
while (S.violations() > 0 && it < 50 * n) {
  select(q in Size, v in Size : S.getAssignDelta(queen[q],v) < 0) {
    queen[q] := v;
    cout << "chng @ " << it << ": queen[" << q << "] = " << v << " viol: " << S.violations() << endl;
  }
  it = it + 1;
}
cout << queen << endl;
```

```
import cotls;
int n = 16;
range Size = 1..n;
UniformDistribution distr(Size);

Solver<LS> m();
var{int} queen[Size](m,Size) := distr.get();
ConstraintSystem<LS> S(m);

S.post(alldifferent(queen));
S.post(alldifferent(all(i in Size) queen[i] + i));
S.post(alldifferent(all(i in Size) queen[i] - i));
m.close();

int it = 0;
while (S.violations() > 0 && it < 50 * n) {
  selectMin(q in Size,v in Size)(S.getAssignDelta(queen[q],v)) {
    queen[q] := v;
    cout << "chng @ " << it << ": queen[" << q << "] := " << v << " viol: " << S.violations() <<
      endl;
  }
  it = it + 1;
}
cout << queen << endl;
```

In Comet

First Improvement

queensLS2.co

```
import cotls;
int n = 16;
range Size = 1..n;
UniformDistribution distr(Size);

Solver<LS> m();
var{int} queen[Size](m,Size) := distr.get();
ConstraintSystem<LS> S(m);

S.post(alldifferent(queen));
S.post(alldifferent(all(i in Size) queen[i] + i));
S.post(alldifferent(all(i in Size) queen[i] - i));
m.close();

int it = 0;
while (S.violations() > 0 && it < 50 * n) {
  selectFirst(q in Size, v in Size: S.getAssignDelta(queen[q],v) < 0) {
    queen[q] := v;
    cout << "chng @ " << it << ": queen[" << q << "] := " << v << " viol: " << S.violations() <<
      endl;
  }
  it = it + 1;
}
cout << queen << endl;
```

In Comet

Min Conflict Heuristic

```
import cotls;
int n = 16;
range Size = 1..n;
UniformDistribution distr(Size);

Solver<LS> m();
var{int} queen[Size](m,Size) := distr.get();
ConstraintSystem<LS> S(m);

S.post(alldifferent(queen));
S.post(alldifferent(all(i in Size) queen[i] + i));
S.post(alldifferent(all(i in Size) queen[i] - i));
m.close();

int it = 0;
while (S.violations() > 0 && it < 50 * n) {
  select(q in Size : S.violations(queen[q])>0) {
    selectMin(v in Size)(S.getAssignDelta(queen[q],v)) {
      queen[q] := v;
      cout << "chng @ " << it << ": queen[" << q << "] := " << v << " viol: " << S.violations() <<
        endl;
    }
  }
  it = it + 1;
}
cout << queen << endl;
```

```
function void conflictSearch (Constraint<LS> c, int itLimit) {  
    int it = 0;  
    var{int}[] x = c.getVariables();  
    range Size = x.getRange();  
    while (!c.isTrue() && it < itLimit) {  
        selectMax(i in Size)(c.violations(x[i]))  
            selectMin(v in x[i].getDomain())(c.getAssignDelta(x[i],v))  
                x[i] := v;  
        it = it + 1;  
    }  
}
```

```
import cotls;  
int n = 16;  
range Size = 1..n;  
UniformDistribution distr(Size);  
  
Solver<LS> m();  
var{int} queen[Size](m,Size) := distr.get();  
ConstraintSystem<LS> S(m);  
  
S.post(alldifferent(queen));  
S.post(alldifferent(all(i in Size) queen[i] + i));  
S.post(alldifferent(all(i in Size) queen[i] - i));  
m.close();  
  
conflictSearch(S,50*n);  
cout << queen << endl;
```

Examples

Random-order first improvement for the TSP

- **Given:** TSP instance G with vertices v_1, v_2, \dots, v_n .
- **search space:** Hamiltonian cycles in G ;
- **neighborhood relation N :** standard 2-exchange neighborhood
- **Initialization:**
 - search position := fixed canonical tour $\langle v_1, v_2, \dots, v_n, v_1 \rangle$
 - P := random permutation of $\{1, 2, \dots, n\}$
- **Search steps:** determined using first improvement w.r.t. $f(s)$ = cost of tour s , evaluating neighbors in order of P (does not change throughout search)
- **Termination:** when no improving search step possible (local minimum)

Iterative Improvement for TSP

TSP-2opt-first(s)

input: an initial candidate tour $s \in S(\epsilon)$

output: a local optimum $s \in S_\pi$

$\Delta = 0$;

for $i = 1$ to $n - 2$ **do**

if $i = 1$ **then** $n' = n - 1$ **else** $n' = n$

for $j = i + 2$ to n' **do**

$\Delta_{ij} = d(c_i, c_j) + d(c_{i+1}, c_{j+1}) - d(c_i, c_{i+1}) - d(c_j, c_{j+1})$

if $\Delta_{ij} < 0$ **then**

 UpdateTour(s, i, j)

is it really?