POSL: A Parallel-Oriented Solver Language

THESIS FOR THE DEGREE OF DOCTOR OF COMPUTER SCIENCE

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Submitted: dd/mm/2016

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Part I

PRESENTATION

A Parallel-Oriented Language for Modeling Meta-Heuristic-Based Solvers

In this chapter POSL is introduced as the main contribution, and a new way to solve CSPs. Its characteristics and advantages are summarized, and a general procedure to be followed is described, in order to build parallel solvers using POSL, followed by a detailed description of each of the single steps.

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In this chapter we present the different steps to build communicating parallel solvers with POSL. First of all, the algorithm we have conceived to solve the target problem is decomposed it into small modules of computation, which are implemented as separated functions. We name them computation modules (see Figure 1.1a, blue shapes). At this point it is crucial to find a good decomposition of its algorithm, because it will have a significant impact in its future re-usage and variability. The next step is to decide which information is interesting to receive from other solvers. This information is encapsulated into another kind of component called communication module, allowing data transmission between solvers (see Figure 1.1a, red shapes). A third stage is to ensemble the modules through POSL's inner language (the interested reader is referred to Appendix [...]) to create independent solvers. The parallel-oriented language based on operators provided by POSL (see Figure 1.1b, green shapes) allows not only the information exchange, but also executing components in parallel. In this stage the information that is interesting to be shared with other solvers is sent using operators. After that we can connect them using communication operators. We call this final entity a solvers set (see Figure 1.1c).

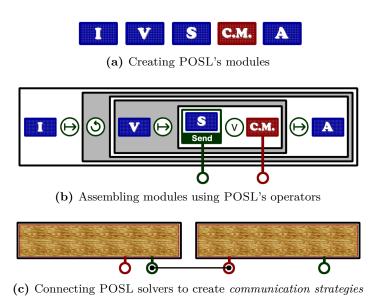


Figure 1.1: Solver construction process using POSL

In the following sections we explain in details all these steps, but first, we explain how to model the target benchmark using POSL.

Modeling the target benchmark

1.1

Target problems are modeled in POSL using the C++ programing language, respecting some rules of the object-oriented design. First of all, the benchmark must inherit from

the class **Benchmark** provided by POSL. This class does not have any method to be overridden or implemented, but receives in its constructor three objects, instances from classes that the user must create. Those classes must inherit from **SolutionCostStrategy**, **RelativeCostStrategy** and **ShowStrategy**, respectively. In these classes the most important functionalities of the benchmark model are defined.

<u>SolutionCostStrategy</u>: In this class the strategy to compute the cost of a configuration is implemented. POSL is based on improving step by step an initial configuration, taking into account a cost function provided by the user through the model (by implementing the function solutionCost(dots)). The kind of problems that POSL solves is the class of Constraint Satisfaction Problems, so this cost function must return an integer taking into account the problem constraints. Given a configuration s, the cost function, as a mandatory rule, must return 0 if and only if s is a solution of the problem, i.e., s fulfill all the problem constraints. An example of cost function is one that returns the number of violated constraints. However, the more expressive the function cost is, the better the performance of POSL leading to the solution.

The method to be implemented in this class is:

int solutionCost(std::vector<int> & c) → Computes the cost of a given configuration (c).

RelativeCostStrategy: In this class the user implements the strategy to compute the cost of a given configuration with respect to another. If the cost of some configuration has been calculated, sometimes it is possible to store some information in order to compute the cost of another configuration, if the differences between them are known. If it is possible, the algorithms is defined in this class. If it is not possible, this class must have the same functionality of SolutionCostStrategy.

The methods to implement in this class are:

- void initializeCostData(std::vector<int> & c) → Initializes the information related to the cost (auxiliary data structures, the current configuration (c), the current cost, etc.)
- void updateConfiguration(std::vector<int> & c) → Updates the information related to the cost.
- int relativeSolutionCost(std::vector<int> & c) \rightarrow Returns the relative cost of the configuration c with respect to the current configuration.
- int currentCost() \rightarrow Property that returns the cost of the current configuration.

- int costOnVariable(int variable_index) → Returns a measure of the contribution of a variable to the total cost of a configuration.
- int sickestVariable() \rightarrow Returns the variable contributing the most to the cost.

<u>SolutionCostStrategy</u>: This class represents the way a benchmark shows a configuration, in order to provide more information about the structure. For example, a configuration of the instance 3–3–2 of the *Social Golfers Problem* (see bellow for more details about this benchmark) can be written as follows:

```
[1, 2, 3, 4, 5, 6, 7, 8, 9, 3, 4, 5, 6, 7, 8, 9, 1, 2]
```

This text is, nevertheless, very difficult to be read if the instance is larger. Therefore, it is recommended that the user implements this class in order to give more details and to make it easier to interpret the configuration. For example, for the same instance of the problem, a solution could be presented as follows:

```
Golfers: players-3, groups-3, weeks-2
6 8 7
1 3 5
4 9 2
--
7 2 3
4 8 1
5 6 9
--
```

The method to be implemented in this class is:

• std::string showSolution(std::shared_ptr<Solution> s) → Returns a string to be written in the standard output.

Once we have modeled the target benchmark, it can be solved using POSL. In the following sections we describe how to use this parallel-oriented language to solve *Constraint Satisfaction Problems*.

First stage: creating POSL's modules

There exist two types of basic modules in POSL: computation modules and communication modules. A computation module is a function which received an input, then executes an internal algorithm, and returns an output. A communication module is also a function receiving and returning information, but in contrast, the communication module can receive

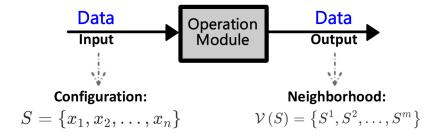


Figure 1.2: An example of a computation module computing a neighborhood

information from two different sources: through input parameters or from outside, i.e., by communicating with a module from another solver.

1.2.1 | Computation Module

A computation module is the most basic and abstract way to define a piece of computation. It is a function which receives an instance of a POSL data type as input, then executes an internal algorithm, and returns an instance of a POSL data type as output. The input and output types will characterize the computation module signature. It can be dynamically replaced by (or combined with) other computation modules, since they can be shared among solvers working in parallel. They are joined through abstract solvers.

Definition 1 (Computation Module) A computation module Cm is a mapping defined by:

$$Cm: D \to I$$
 (1.1)

where D and I can be either a set of configurations, a set of sets of configurations, a set of values of some data type, etc.

Consider a local search meta-heuristic solver. One of its *computation modules* can be the function returning the set of configurations composing the neighborhood of a given configuration:

$$Cm_{neighborhood}: D_1 \times D_2 \times \cdots \times D_n \to 2^{D_1 \times D_2 \times \cdots \times D_n}$$

where D_i represents the definition domains of each variable of the input configuration.

Figure 1.2 shows an example of *computation module*: which receives a configuration S and then computes the set \mathcal{V} of its neighbor configurations $\{S^1, S^2, \dots, S^m\}$.

1.2.1.1 Creating new computation modules

To create new *computation modules* we use C++ programing language. POSL provides a hierarchy of data types to work with (See anexes) and some abstract classes to inherit from, depending on the type of *computation module* that the user wants to create. These abstract classes represent *abstract computation module* and define a type of action to be executed. In the following we present the most important ones:

- AOM_FirstConfigurationGeneration → Represents computation modules generating a first configuration. The user must implement the method spcf_execute(ComputationData) which returns a pointer to a Solution, that is, an object containing all the information concerning a partial solution (configuration, variable domains, etc.)
- AOM_NeighborhoodFunction → Represent computation modules creating a neighborhood of a given configuration. The user must implement the method spcf_execute(Solution) which returns a pointer to an object Neighborhood, containing a set of configurations which constitute the neighborhood of a given configuration, according to certain criteria. These configuration are efficiently stored.
- AOM_SelectionFunction → Represents computation modules selecting a configuration from a neighborhood. The user must implement the method spcf_execute(Neighborhood) which returns a pointer to an object DecisionPair, containing two solutions: the current and the selected one.
- AOM_DecisionFunction → Represents computation modules deciding which of the two solutions will be the current configuration for the next iteration. The user must implement the method spcf_execute(DecisionPair) which returns a pointer to an object Solution.

1.2.2 Communication modules

A communication module is also a function receiving and returning information, but in contrast, the communication module can also receive information by communicating with a module from another solver. A communication module is the component managing the information reception in the communication between solvers (we will talk about information transmission in the next section). They can interact with computation modules through operators (see Figure 1.3).

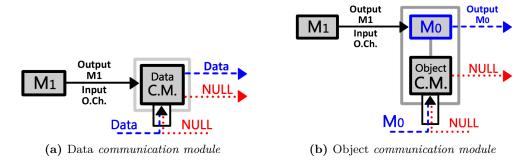


Figure 1.3: Communication module

A communication module can receive two types of information from an external solver: data or computation modules. It is important to notice that by sending/receiving computation modules, we mean sending/receiving only required information to identify and being able to instantiate the computation module.

In order to distinguish from the two types of *communication modules*, we will call Data Communication Module to the *communication module* responsible for the data reception (Figure 1.3a), and Object Communication Module to the one responsible for the reception and instantiation of *computation modules* (Figure 1.3b).

Definition 2 (Data Communication Module) A Data Communication Module Ch is a component that produces a mapping defined as follows:

$$Ch: U \to I$$
 (1.2)

It returns the information I coming from an external solver, no matter what the input U is.

Definition 3 (Object Communication Module) If we denote by \mathbb{M} the space of all the computation modules defined by Definition 1.1, then an Object Communication Module Ch is a component that produces a computation module coming from an external solver as follows:

$$Ch: \mathbb{M} \to \mathbb{M}$$
 (1.3)

Users can implement new computation and connection modules but POSL already contains many useful modules for solving a broad range of problems.

Due to the fact that communication modules receive information coming from outside without having control on them, it is necessary to define the NULL information, in order to denote the absence of information. If a Data Communication Module receives a piece of information, is returned automatically. If a Object Communication Module receives a computation module, it is instantiated and executed with the communication module's input and its result is

returned. In both cases, if no available information exists (no communications performed), the $communication\ module$ returns the NULL object.

Second stage: assembling POSL's modules

Modules mentioned above are defined respecting the signature of some predefined abstract module. For example, the module showed in Figure 1.2 is a computation module based on an abstract module that receives a configuration and returns a neighborhood. In that sense, an example of a concrete computation module (or just computation module) can be a function receiving a configuration, and returning a neighborhood constituted by N configurations which only differ from the input configuration in one entry.

In this stage an *abstract solver* is coded using POSL. It takes abstract modules as *parameters* and combines them through operators. Through the *abstract solver*, we can also decide which information to send to other solvers by using some operators to send the result of a computation module (see below). In the following we present a formal and more detailed specification of POSL's operators.

The abstract solver is the solver's backbone. It joins the computation modules and the communication modules coherently. It is independent from the computation modules and communication modules used in the solver. It means that they can be changed or modified during the execution, without altering the general algorithm, but still respecting the main structure. Each time we combine some of them using POSL's operators, we are creating a compound module. Here we formally define the concept of module and compound module.

Definition 4 A module is (and it is denoted by the letter \mathcal{M}):

a) a computation module or

1.3

- b) a communication module or
- c) $[\mathcal{M}_1 \ OP \ \mathcal{M}_2]$, which is the composition of two modules \mathcal{M}_1 and \mathcal{M}_2 to be executed sequentially, returning an output depending on the nature of the operator OP; or
- d) $[M_1 \ OP \ M_2]$, which is the composition of two modules M_1 and M_2 to be executed, returning an output depending on the nature of the operator OP. These two modules will be executed in parallel if and only if OP supports parallelism, (i.e. some modules will be executed sequentially although they were grouped this way); or sequentially otherwise.

We denote the space of the modules by M and call compound modules to the composition of modules described in c) and d).

For a better understanding of Definition 4, Figure 1.4 shows graphically the structure of a compound module.

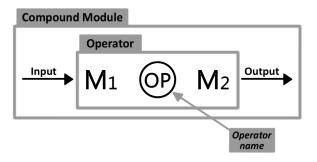


Figure 1.4: A compound module

As mentioned before, the abstract solver is independent from the computation modules and communication modules used in the solver. It means that one abstract solver can be used to construct many different solvers, by implementing it using different modules (see below the related concept of abstract solver instantiation). This is the reason why the abstract solver is defined only using abstract modules. Formally, we define an abstract solver as follows:

Definition 5 (Abstract Solver) An Abstract Solver AS is a triple $(M, \mathcal{L}^m, \mathcal{L}^c)$, where: **M** is a compound module (also called root compound module), \mathcal{L}^m a list of abstract computation modules appearing in \mathcal{M} , and \mathcal{L}^c a list of communication modules appearing in \mathcal{M} .

The root compound module can be defined also as a free-context grammar as follows:

Definition 6 (root compound module's grammar) $G_{POSL} = (\mathbf{V}, \Sigma, \mathbf{S}, \mathbf{R})$, where:

- a) $\mathbf{V} = \{CM, OP\}$ is the set of variables,
- $b) \ \ \Sigma = \left\{\alpha,\beta,be,[,],\llbracket,\rrbracket_p\,,(,),\{,\},\llbracket,\rrbracket^m,\rrbracket^o,\stackrel{?}{\longleftrightarrow},\stackrel{?}{\circlearrowleft},\circlearrowleft,\stackrel{?}{\o},\stackrel{?}{\smile},\stackrel$
- c) $S = \{CM\}$ is the set of start variables,
- d) and $\mathbf{R} =$

is a set of rules

In the following I explain some of the concepts in Definition 6:

- The variables CM and OP are two very important entities in the language, as it can be seen in the grammar. We name them *compound module* and *operator*, respectively.
- The terminals α and β represent a computation module and a communication module, respectively.
- \bullet The terminal be is a boolean expression.
- The terminals $[\]$, $[\]$ _p are symbols for grouping and defining the way the involved compound modules are executed. Depending on the nature of the operator, this can be either sequentially or in parallel:
 - a) [OP]: The involved operator is executed sequentially.
 - b) $[\![OP]\!]_p$: The involved operator is executed in parallel if and only if OP supports parallelism. Otherwise, an exception is thrown.
- The terminals (and) are symbols for grouping the boolean expression in some operators.
- The terminals { and } are symbols for grouping compound modules in some operators.
- The terminals $(.)^m$, $(.)^o$, are operators to send information to other solvers (explained bellow).
- The rest of terminals are POSL operators.

In the following we define POSL operators. In order to group modules, like in Definition 4(c)) and 4(d)), we will use |.| as generic grouper.

Definition 7 (Operator Sequential Execution) Let

a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and

b) $\mathcal{M}_2:\mathcal{D}_2\to\mathcal{I}_2$,

be modules, where $\mathcal{I}_1 \subseteq \mathcal{D}_2$. Then the operation $\left| \mathcal{M}_1 \middle{\mapsto} \mathcal{M}_2 \right|$ defines the compound module \mathcal{M}_{seq} as the result of executing \mathcal{M}_1 followed by executing \mathcal{M}_2 :

$$\mathcal{M}_{sea}:\mathcal{D}_1 o \mathcal{I}_2$$

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This is an example of an operator that does not support the execution of its involved compound modules in parallel, because the input of the second compound module is the output of the first one.

The following operator is very useful to execute modules sequentially creating bifurcations, subject to some boolean condition:

Definition 8 (Operator Conditional Execution) Let

- a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and
- b) $\mathcal{M}_2:\mathcal{D}_2\to\mathcal{I}_2$,

be modules, where $\mathcal{D}_1 \subseteq \mathcal{D}_2$. Then the operation $\left|\mathcal{M}_1\right|_{< cond} \mathcal{M}_2$ defines the compound module \mathcal{M}_{cond} as result of the sequential execution of \mathcal{M}_1 if < cond > is true or \mathcal{M}_2 , otherwise:

$$\mathcal{M}_{cond}: \mathcal{D}_1 \cap \mathcal{D}_2 \to \mathcal{I}_1 \cup \mathcal{I}_2$$

We can execute modules sequentially creating also cycles.

Definition 9 (Operator Cyclic Execution) Let $\mathcal{M}: \mathcal{D} \to \mathcal{I}$ be a module, where $\mathcal{I} \subseteq \mathcal{D}$. Then, the operation $|\circlearrowleft_{< cond>} \mathcal{M}|$ defines the compound module \mathcal{M}_{cyc} as result of the sequential execution of \mathcal{M} repeated while < cond> remains **true**:

$$\mathcal{M}_{cuc}:\mathcal{D}
ightarrow\mathcal{I}$$

Definition 10 (Operator Random Choice) Let

- a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and
- b) $\mathcal{M}_2: \mathcal{D}_2 \to \mathcal{I}_2$,

be modules, where $\mathcal{D}_1 \subset \mathcal{D}_2$ and a real value ρ . Then the operation $|M_1(\rho)\mathcal{M}_2|$ defines the compound module \mathcal{M}_{rho} that executes and returns the output of \mathcal{M}_1 with probability ρ , or executes and returns the output of \mathcal{M}_2 with probability $(1 - \rho)$:

$$\mathcal{M}_{rho}: \mathcal{D}_1 \cap \mathcal{D}_2 \to \mathcal{I}_1 \cup \mathcal{I}_2$$

The following operator is very useful if the user needs to use a *communication module* inside an *abstract solver*. As explained before, if a *communication module* does not receive any information from another solver, it returns *NULL*. This may cause the undesired termination of the solver if this case is not considered correctly. Next, I introduce the operator **Operator Not** *NULL* **Execution** and illustrate how to use it in practice with an example.

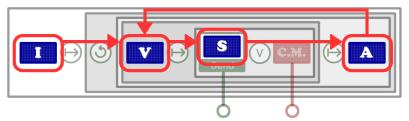
Definition 11 (Operator Not NULL Execution) Let

- a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and
- b) $\mathcal{M}_2: \mathcal{D}_2 \to \mathcal{I}_2$,

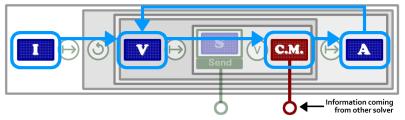
be modules, where $\mathcal{D}_1 \subseteq \mathcal{D}_2$. Then, the operation $|\mathcal{M}_1 \bigcup \mathcal{M}_2|$ defines the compound module \mathcal{M}_{non} that executes \mathcal{M}_1 and returns its output if it is not NULL, or executes \mathcal{M}_2 and returns its output otherwise:

$$\mathcal{M}_{non}: \mathcal{D}_1 \cap \mathcal{D}_2 \to \mathcal{I}_1 \cup \mathcal{I}_2$$

Figure 1.5 shows how to combine a connection module with the *computation module* S through the operator \bigcirc . Here, the *computation module* S will be executed as long as the *communication module* remains NULL, i.e., there is no information coming from outside. This behavior is represented in Figure 1.5a by the orange lines. If some data has been received through the *communication module*, the later is executed instead of the module S, represented in Figure 1.5b by blue lines.



(a) The solver executes the computation module ${\bf S}$ if no information is received through the connection module



(b) The solver uses the information coming from an external solver

Figure 1.5: Two different behaviors within the same solver

This is *short-circuit* operator. It means that if the first argument (module) does not return *NULL*, the second will not be executed. POSL provides another operator with the same functionality but not *short-circuit*:

Definition 12 (Operator BOTH Execution) Let

- a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and
- b) $\mathcal{M}_2:\mathcal{D}_2\to\mathcal{I}_2$,

be modules, where $\mathcal{D}_1 \subseteq \mathcal{D}_2$. Then the operation $|\mathcal{M}_1 \cap \mathcal{M}_2|$ defines the compound module \mathcal{M}_{both} that executes both \mathcal{M}_1 and \mathcal{M}_2 , then returns the output of \mathcal{M}_1 if it is not NULL, or the output of \mathcal{M}_2 otherwise:

$$\mathcal{M}_{both}: \mathcal{D}_1 \cap \mathcal{D}_2 \to \mathcal{I}_1 \cup \mathcal{I}_2$$

In the following definitions, the concepts of *cooperative parallelism* and *competitive parallelism* are implicitly included. We say that cooperative parallelism exists when two or more processes are running separately, they are independent, and the general result will be some combination of the results of all the involved processes (e.g. Definitions 13 and 14). On the other hand, competitive parallelism arise when the general result is the result of the process ending first (e.g. Definition 15).

Definition 13 (Operator Minimum) Let

- a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and
- b) $\mathcal{M}_2: \mathcal{D}_2 \to \mathcal{I}_2$,

be modules, where $\mathcal{D}_1 \subseteq \mathcal{D}_2$. Let also o_1 and o_2 be the outputs of \mathcal{M}_1 and \mathcal{M}_2 , respectively. Assume that there exists some order criteria between them. Then the operation $\left|\mathcal{M}_1(m)\mathcal{M}_2\right|$ defines the compound module \mathcal{M}_{min} that executes \mathcal{M}_1 and returns $\min \{o_1, o_2\}$:

$$\mathcal{M}_{min}: \mathcal{D}_1 \cap \mathcal{D}_2 \to \mathcal{I}_1 \cup \mathcal{I}_2$$

Similarly we define the operator **Maximum**:

Definition 14 (Operator Maximum) Let

- a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and
- b) $\mathcal{M}_2: \mathcal{D}_2 \to \mathcal{I}_2$,

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be modules, where $\mathcal{D}_1 \subseteq \mathcal{D}_2$. Let also o_1 and o_2 be the outputs of \mathcal{M}_1 and \mathcal{M}_2 , respectively. Assume that there exists some order criteria between them. Then the operation $\left|\mathcal{M}_1(M)\mathcal{M}_2\right|$ defines the compound module \mathcal{M}_{max} that executes \mathcal{M}_1 and returns $\max\{o_1,o_2\}$:

$$\mathcal{M}_{max}: \mathcal{D}_1 \cap \mathcal{D}_2 \to \mathcal{I}_1 \cup \mathcal{I}_2$$

Definition 15 (Operator Race) Let

- a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and
- b) $\mathcal{M}_2: \mathcal{D}_2 \to \mathcal{I}_2$,

be modules, where $\mathcal{D}_1 \subseteq \mathcal{D}_2$ and $\mathcal{I}_1 \subset \mathcal{I}_2$. Then the operation $\left|\mathcal{M}_1 \downarrow \mathcal{M}_2\right|$ defines the compound module \mathcal{M}_{race} that executes both modules \mathcal{M}_1 and \mathcal{M}_2 , and returns the output of the module ending first:

$$\mathcal{M}_{race}: \mathcal{D}_1 \cap \mathcal{D}_2 \to \mathcal{I}_1 \cup \mathcal{I}_2$$

The operators presented in Definitions 13, 14 and 15 are very useful in terms of sharing not only information between solvers, but also behaviors. If one of the operands is a communication module it can receive an external solver's computation module, providing the opportunity to instantiate it in the current solver. The operator either instantiates that module if it is not null, and executes it; or executes the other operand module otherwise.

Some others operators can be useful when dealing with sets.

Definition 16 (Operator Union) Let

- a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and
- b) $\mathcal{M}_2:\mathcal{D}_2\to\mathcal{I}_2$,

be modules, where $\mathcal{D}_1 \subseteq \mathcal{D}_2$. Let also V_1 and V_2 be the outputs of \mathcal{M}_1 and \mathcal{M}_2 , respectively. Then the operation $\left|\mathcal{M}_1 \bigcup \mathcal{M}_2\right|$ defines the compound module \mathcal{M}_{\cup} that executes both modules \mathcal{M}_1 and \mathcal{M}_2 , and returns $V_1 \cup V_2$:

$$\mathcal{M}_{\cup}: \mathcal{D}_1 \cap \mathcal{D}_2 \to \mathcal{I}_1 \cup \mathcal{I}_2$$

Similarly we define the operators **Intersection** and **Subtraction**:

Definition 17 (Operator Intersection) Let

a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and

b)
$$\mathcal{M}_2: \mathcal{D}_2 \to \mathcal{I}_2$$
,

be modules, where $\mathcal{D}_1 \subseteq \mathcal{D}_2$. Let also V_1 and V_2 be the outputs of \mathcal{M}_1 and \mathcal{M}_2 , respectively. Then the operation $|\mathcal{M}_1 \cap \mathcal{M}_2|$ defines the compound module \mathcal{M}_{\cap} that executes both modules \mathcal{M}_1 and \mathcal{M}_2 , and returns $V_1 \cap V_2$:

$$\mathcal{M}_{\cap}:\mathcal{D}_1\cap\mathcal{D}_2\to\mathcal{I}_1\cup\mathcal{I}_2$$

Definition 18 (Operator Subtraction) Let

a) $\mathcal{M}_1: \mathcal{D}_1 \to \mathcal{I}_1$ and

b) $\mathcal{M}_2: \mathcal{D}_2 \to \mathcal{I}_2$,

be modules, where $\mathcal{D}_1 \subseteq \mathcal{D}_2$. Let also V_1 and V_2 be the outputs of \mathcal{M}_1 and \mathcal{M}_2 , respectively. Then the operation $\left|\mathcal{M}_1 \supseteq \mathcal{M}_2\right|$ defines the compound module \mathcal{M}_- that executes both modules \mathcal{M}_1 and \mathcal{M}_2 , and returns $V_1 - V_2$:

$$\mathcal{M}_{-}:\mathcal{D}_{1}\cap\mathcal{D}_{2}\to\mathcal{I}_{1}\cup\mathcal{I}_{2}$$

Now, I define the operators which allows to send information to other solvers. Two types of information can be sent: i) the output of the *computation module* and send its output, or ii) the *computation module* itself. This utility is very useful in terms of sharing behaviors between solvers.

Definition 19 (Sending Data Operator) Let $\mathcal{M}: \mathcal{D} \to \mathcal{I}$ be a module. Then the operation $|(\mathcal{M})^o|$ defines the compound module \mathcal{M}_{sendD} that executes the module \mathcal{M} and sends its output outside:

$$\mathcal{M}_{sendD}: \mathcal{D}
ightarrow \mathcal{I}$$

Similarly we define the operator **Send Module**:

Definition 20 (Sending Module Operator) Let $\mathcal{M}: \mathcal{D} \to \mathcal{I}$ be a module. Then the operation $|\langle \mathcal{M} \rangle^m|$ defines the compound module \mathcal{M}_{sendM} that executes the module \mathcal{M} , then returns its output and sends the module itself outside:

$$\mathcal{M}_{sendM}: \mathcal{D}
ightarrow \mathcal{I}$$

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Once all desired abstract modules are linked together with operators, we obtain the *root* compound module, an important part of an abstract solver. To implement a concrete solver from an abstract solver, one must instantiate each abstract module with a concrete one respecting the required signature. From the same abstract solver, one can implement many different concrete solvers simply by instantiating abstract modules with different concrete modules.

An abstract solver is defined as follows: after declaring the **abstract solver**'s name, the first line defines the list of abstract computation modules, the second one the list of abstract communication modules, then the algorithm of the solver is defined as the solver's body (the root compound module), between **begin** and **end**.

An abstract solver in the following simple regular expression:

abstract solver name computation: L^m (communication: L^c)? begin \mathcal{M} end

where:

- name is the identifier of the abstract solver,
- ullet L^m is the list of abstract computation modules,
- L^c is the list of abstract communication modules, and
- \mathcal{M} is the root compound module.

For instance, Algorithm 1 illustrates the abstract solver corresponding to Figure 1.1b.

Algorithm 1: POSL pseudo-code for the abstract solver presented in Figure 1.1b

1.4 Third stage: creating POSL solvers

With computation and communication modules composing an abstract solver, one can create solvers by instantiating modules. This is simply done by specifying that a given solver must implements a given abstract solver, followed by the list of computation then communication modules. These modules must match signatures required by the abstract solver. Algorithm 2 implements Algorithm 1 by instantiating modules shown in Figure 1.5.

Algorithm 2: An instantiation of the abstract solver presented in Algorithm 1

solver solver_01 implements as_01 computation : $I_{rand}, V_{std}, S_{best}, A_{alw}$

connection: CM_{last}

Algorithm 2 is just an example of a solver instantiation, using some *computation modules* provided by POSL, that are used and explained in details in the Chapter ?? of this document:

- I_{rand} creates a random configuration.
- \bullet V_{std} creates a neighborhood of a given configuration, changing one element at a time.
- S_{best} selects the configuration of a neighborhood with the lowest cost.
- A_{alw} always accepts the incoming configuration.
- CM_{last} returns the last configuration arrived, if at the time of its execution, there is more than one configuration waiting to be received.

1.5 Forth stage: connecting the solvers

We call *solver set* to the pool of (concrete) solvers that we plan to use in parallel to solve a problem. Once we have our solvers set, the last stage is to connect the solvers to each other. Up to this point, solvers are disconnected, but they are ready to establish the communication. POSL provides a platform to the user such that cooperative strategies can be easily defined.

In the following we present two important concepts necessary to formalize the *communication* operators.

Definition 21 (Communication Jack) Let S be a solver. Then the operation $S \cdot M$ opens an outgoing connection from the solver S, sending to the outside either a) the output of M, if it is affected by a sending data operator as presented in Definition 19, or b) M itself, if it is affected by a sending module operator as presented in Definition 20.

Definition 22 (Communication Outlet) Let S be a solver. Then, the operation $S \cdot \mathcal{CM}$ opens an ingoing connection to the solver S, receiving from the outside either a) the output of some computation module, if \mathcal{CM} is a data communication module, or b) a computation module, if \mathcal{CM} is an object communication module.

The communication is established by following the following rules guideline:

- a) Each time a solver sends any kind of information by using a *sending* operator, it creates a *communication jack*.
- b) Each time a solver defines a communication module, it creates a communication outlet.
- c) Solvers can be connected to each other by linking communication jacks to communication outlets.

Following, we define the *connection operators* that POSL provides.

Definition 23 (Connection Operator One-to-One) Let

- a) $\mathcal{J} = [\mathcal{S}_0 \cdot \mathcal{M}_0, \mathcal{S}_1 \cdot \mathcal{M}_1, \dots, \mathcal{S}_{N-1} \cdot \mathcal{M}_{N-1}]$ be the list of communication jacks, and
- b) $\mathcal{O} = [\mathcal{Z}_0 \cdot \mathcal{CM}_0, \mathcal{Z}_1 \cdot \mathcal{CM}_1, \dots, \mathcal{Z}_{N-1} \cdot \mathcal{CM}_{N-1}]$ be the list of communication outlets

Then the operation

$$\mathcal{J} \bigoplus \mathcal{O}$$

connects each communication jack $S_i \cdot M_i \in \mathcal{J}$ with the corresponding communication outlet $Z_i \cdot \mathcal{CM}_i \in \mathcal{O}, \forall 0 \leq i \leq N-1$ (see Figure 1.6a).

Definition 24 (Connection Operator One-to-N) Let

- a) $\mathcal{J} = [S_0 \cdot \mathcal{M}_0, S_1 \cdot \mathcal{M}_1, \dots, S_{N-1} \cdot \mathcal{M}_{N-1}]$ be the list of communication jacks, and
- b) $\mathcal{O} = [\mathcal{Z}_0 \cdot \mathcal{CM}_0, \mathcal{Z}_1 \cdot \mathcal{CM}_1, \dots, \mathcal{Z}_{M-1} \cdot \mathcal{CM}_{M-1}]$ be the list of communication outlets

Then the operation

$$\mathcal{J} \stackrel{(\leadsto)}{(\leadsto)} \mathcal{O}$$

connects each communication jack $S_i \cdot M_i \in \mathcal{J}$ with every communication outlet $Z_j \cdot \mathcal{C}M_j \in \mathcal{O}$, $\forall 0 \leq i \leq N-1$ and $0 \leq j \leq M-1$ (see Figure 1.6b).

Definition 25 (Connection Operator Ring) Let

- a) $\mathcal{J} = [\mathcal{S}_0 \cdot \mathcal{M}_0, \mathcal{S}_1 \cdot \mathcal{M}_1, \dots, \mathcal{S}_{N-1} \cdot \mathcal{M}_{N-1}]$ be the list of communication jacks, and
- b) $\mathcal{O} = [\mathcal{S}_0 \cdot \mathcal{CM}_0, \mathcal{S}_1 \cdot \mathcal{CM}_1, \dots, \mathcal{S}_{N-1} \cdot \mathcal{CM}_{N-1}]$ be the list of communication outlets

Then the operation

$$\mathcal{J} \left(\longleftrightarrow \right) \mathcal{O}$$

connects each communication jack $S_i \cdot M_i \in \mathcal{J}$ with the corresponding communication outlet $\mathcal{Z}_{(i+1)\%N} \cdot \mathcal{CM}_{(i+1)\%N} \in \mathcal{O}, \ \forall 0 \leq i \leq N-1 \ (\text{see Figure 1.6c}).$

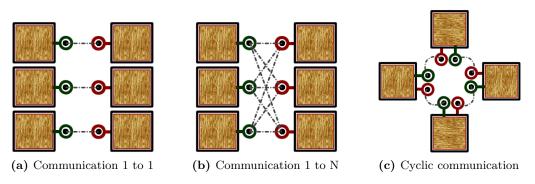


Figure 1.6: Graphic representation of communication operators

POSL also allows to declare non-communicating solvers to be executed in parallel, declaring only the list of solver names:

$$[S_0, S_1, \ldots, S_{N-1}]$$

When we apply a connection operator op between a communication jacks list \mathcal{J} and a communication outlets list \mathcal{O} , internally we are assigning an abstract computation unit (typically a thread) to each solver that we declare in each list. This assignment receives the name of Solver Scheduling. Before running the solver set, this abstract unit of computation is just an integer $\tau \in [0..N]$ identifying uniquely each of the solvers. When the solver set is launched, the solver with the identifier τ runs into the computation unit τ . This identifier assignation remains independent of the real availability of resources of computation. It just takes into account the user declaration. This means that, if the user declares 30 solvers (15 senders and 15 receivers) and the solver set is launched using 20 cores, only the first 20 solvers will be executed, and in consequence, there will be 10 solvers sending information to nowhere. Users should take this into account when declaring the solver set.

The connection process depends on the applied conection operator. In each case the goal is to assign, to the sending operator $((.)^o)$ or $(.)^m$ inside the abstract solver, the identifier of

the solver (or solvers, depending on the connection operator) where the information will be sent. Algorithm 3 presents the connection process.

Algorithm 3: Scheduling and connection main algorithm

```
input : \mathcal{J} list of communication jacks,
                \mathcal{O} list of communication outlets
1 while no available jacks or outlets do
        S_{jack} \leftarrow \texttt{GetNext}(\mathcal{J})
        R_{oulet} \leftarrow \texttt{GetNext}(\mathcal{O})
3
        S \leftarrow \texttt{GetSolverFromConnector}(S_{iack})
4
        R \leftarrow \texttt{GetSolverFromConnector}(R_{oulet})
5
        Schedule(S)
6
        R_{id} \leftarrow \texttt{Schedule}(R)
7
        Connect(root(S), S_{jack}, R_{id})
9 end
```

In Algorithm 3:

- GetNext(...) returns the next available solver-jack (or solver-outlet) in the list, depending on the connection operator, e.g., for the connection operator One-to-N, each communication jack in \mathcal{J} must be connected with each communication outlet in \mathcal{O} .
- GetSolverFromConnector(...) returns the solver name given a connector declaration.
- Schedule(...) schedules a solver and returns its identifier.
- Root(...) returns the root compound module of a solver.
- Connect(...) searches the computation module S_{jack} recursively inside the root compound module of S and places the identifier R_{id} into its list of destination solvers.

1.5.1 Solver name space expansion

One of the goals of POSL is to provide a way to declare sets of solvers to be executed in parallel fast and easily. For that reason, POSL provides two forms of namespace expansion, in order to create sets of solvers using already declared ones:

Solver name expansion - Uses an integer K to denote how many times the solver name S will appear in the declaration. $[\ldots S_i \cdot \mathcal{M}(K), \ldots]$ expands as $[\ldots S_i \cdot \mathcal{M}, S_i^2 \cdot \mathcal{M}, \ldots S_i^K \cdot \mathcal{M} \ldots]$

and all new solvers S_i^j , $j \in [2..K]$ are created using the same solver declaration of solver S_i .

Connection declaration expansion - Uses an integer K to denote how many times the connection will be repeated in the declaration. Let a) $[S_1 \cdot M_1, \dots, S_N \cdot M_N]$ and

b) $[\mathcal{R}_1 \cdot \mathcal{CM}_1, \dots, \mathcal{R}_M \cdot \mathcal{CM}_M]$ be the list of *communication jacks* and *communication outlets*, respectively, and c) (op) a connection operator. Then

$$[S_1 \cdot \mathcal{M}_1, \dots, S_N \cdot \mathcal{M}_N]$$
 (op) $[R_1 \cdot \mathcal{C}\mathcal{M}_1, \dots, R_M \cdot \mathcal{C}\mathcal{M}_M]$ K

expands as

1.6

$$[\mathcal{S}_{1} \cdot \mathcal{M}_{1}, \dots, \mathcal{S}_{N} \cdot \mathcal{M}_{N}] \underbrace{op} [\mathcal{R}_{1} \cdot \mathcal{C}\mathcal{M}_{1}, \dots, \mathcal{R}_{N} \cdot \mathcal{C}\mathcal{M}_{N}]$$

$$[\mathcal{S}_{1}^{2} \cdot \mathcal{M}_{1}, \dots, \mathcal{S}_{N}^{2} \cdot \mathcal{M}_{N}] \underbrace{op} [\mathcal{R}_{1}^{2} \cdot \mathcal{C}\mathcal{M}_{1}, \dots, \mathcal{R}_{N}^{2} \cdot \mathcal{C}\mathcal{M}_{N}]$$

$$\dots$$

$$[\mathcal{S}_{1}^{K} \cdot \mathcal{M}_{1}, \dots, \mathcal{S}_{N}^{K} \cdot \mathcal{M}_{N}] \underbrace{op} [\mathcal{R}_{1}^{K} \cdot \mathcal{C}\mathcal{M}_{1}, \dots, \mathcal{R}_{N}^{K} \cdot \mathcal{C}\mathcal{M}_{N}]$$

and all new solvers S_i^k , $i \in [1..N]$ and R_j^k , $j \in [1..M]$, $k \in [2..K]$, are created using the same solver declaration of solvers S_i and R_j , respectively.

Step-by-step POSL code example

In this section we explain how to create a solver using POSL through an example. POSL creates solvers based on local search meta-heuristics algorithms. These algorithms have a common structure: 1. they start by initializing some data structures (e.g. a tabu list for Tabu Search [34], a temperature for Simulated Annealing [32], etc.). 2. An initial configuration s is generated. 3. A new configuration s' is selected from the neighborhood $\mathcal{V}(s)$. 4. If s' is a solution for the problem P, then the process stops, and s' is returned. If not, the data structures are updated, and s' is accepted or not for the next iteration, depending on some criterion. An example of such data structure can be the penalizing features of local optima defined by Boussaïd et al [31] in their algorithm Guided Local Search.

Abstract computation modules composing local search meta-heuristics are the following:

Abstract Computation module - 1 I: Generating a configuration s

 $\overline{Abstract}$ Computation $module-2 \mid V$: Defining the neighborhood $\mathcal{V}\left(s\right)$

Abstract Computation module $-3 \mid S$: Selecting $s' \in \mathcal{V}(s)$

Abstract Computation module $-4 \mid A$: Evaluating an acceptance criteria for s'

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To be more specific in our example, we describe come concrete *computation modules* that we can use:

Computation module $-1 \mid I_{rand}$: Generates a random configuration s

Computation module $-2 \mid V_{1ch}$: Defines the neighborhood $\mathcal{V}(s)$ changing only one element

Computation module -3 S_{best} : Selects the best configuration $s' \in \mathcal{V}(s)$ improving the current cost.

Computation module - 4 $A_{a.i.}$: Evaluates an acceptance criteria for s'. We have chosen the classical module selecting the configuration with the lowest global cost, *i.e.*, the one which is likely the closest to a solution.

We can combine modules to create more complex ones. For example, in the example we use in this section, we want to apply some selection criteria, but if there is not improvement in the current configuration, we select a configuration randomly. To do so, we use the operator ? to combine the *computation module* 3 with another:

Computation module -3.1 S_{rand} : Selects a random configuration $s' \in \mathcal{V}(s)$.

Let's make this solver a little bit more complex: in the time of applying the acceptance criteria, suppose that we want to receive a configuration from other solver, to compare it with ours and take it if its cost is lower. We can do that applying the operator (m) to combine the computation module 4 with a communication module:

Communication module - 1: C.M.: Receiving a configuration.

Figure 1.7 shows a graphic representation of the *abstract solver* that we create for our example. In this Figure all the modules are abstract, so we can instantiate them afterwards to create concrete solvers. Note that we also want to send the output of the *computation module A*, in order to share the current configuration with other solvers.

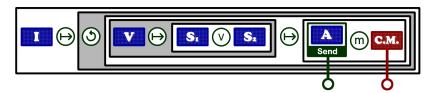


Figure 1.7: Graphic representation of an abstract solver

Algorithm 4 shows the POSL pseudo-code for the *abstract solver* described above, using predefined operators, and Algorithm 5 shows the concrete solver definition using the concrete *computation* and *communication modules* already presented. In the later, we have used the

concrete communication module CM_{last} , that receives all the configuration sent to it, and selects the last arrived.

In practice, POSL provides information regarding the execution process, such as number of ITERATIONS, solver execution TIME, BEST found solution, among others.

Algorithm 4: POSL pseudo-code for abstract solver presented in Figure 1.7

```
1 abstract solver as\_example
2 computation : I, V, S_1, S_2, A
3 connection: C.M.
4 begin
5 I \hookrightarrow
6 [\circlearrowleft (ITR \% K_2) \text{ begin}
7 [V \hookrightarrow [S_1 \lor S_2]] \hookrightarrow [(A)^o \circlearrowleft C.M.]
8 end]
9 end
```

Algorithm 5: An instantiation of the abstract solver in Algorithm 4

```
1 solver S implements as_example
2 computation : I_{rand}, V_{1ch}, S_{best}, A_{a.i.}
3 connection: CM_{last}
```

Now, supose that we want to create another solver Z using a different neighborhood function, that we will call V_{2ch} , performing two changes instead of only one. As a communication strategy, we want also connect them through the operator 1 to N, using Z as sender and S as receiver. Then, we need to declare, using name spare expansions, how many solvers we want to connect each others. Algorithm 6 shows the desired communication strategy.

Algorithm 6: A communication strategy

```
1 [ Z \cdot A (3) ] \longleftrightarrow [ S \cdot CM_{last} (3) ] 2;
```

Notice that the connection operation is affected also by the number 2 at the end, to expand also the topology. In that sense, supposing that we have 12 units of computation available, we obtain a *solver set* working on parallel following the topology described in Figure 1.8.

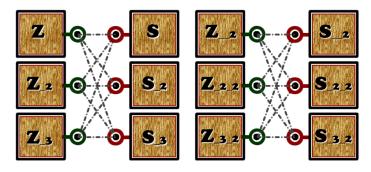


Figure 1.8: An example of connection strategy for 12 units of computation

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