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A Refinement-based compiler development for synchronous languages

Jean-Paul Bodeveix
IRIT Université Paul Sabatier
Toulouse, France
bodeveix@irit.fr

Mamoun Filali-Amine
IRIT CNRS
Toulouse, France
filali@irit.fr

Shuanglong Kan
Nanjing University of Aeronautics and Astronautics
Nanjing, China
kanshuanglong@outlook.com

ABSTRACT

In this paper, we are concerned by the elaboration of generic development steps for the code generation for synchronous languages. Our aim is to provide a correct by construction solution. For that purpose, we adopt a refinement-based approach where proof obligations for each step guarantee properties preservation. We use the Event-B formal method. We start with a big step semantics specified by an Event-B machine. Through a sequence of refinements, expressed as Event-B refinement machines, we end up with a code generation step which implements a small step semantics preserving the properties of the big step semantics.

KEYWORDS

synchronous languages, refinement, code generation, semantics, verification

1 INTRODUCTION

Our study concerns system development. We are interested by the development of reactive systems. For such systems, for instance, avionic systems and interactive systems, synchronous languages have revealed well suited. From the beginning, synchronous languages have put forward a strong mathematical framework underlying their basic principles: synchrony and deterministic concurrency [5]. Thanks to that, synchronous languages have received a significant audience, especially within the critical embedded systems community where high levels of certification are required. However, one must acknowledge that although significant efforts have been devoted to explain the software architecture for implementing frameworks for the development of such languages and even to mechanize them [6, 10, 19], the production of a certified compiler remains a challenge.

In this paper, we are concerned by the elaboration of generic development steps for the code generation for synchronous languages. Our aim is to provide a correct by construction solution. For that purpose, we adopt a refinement-based approach where proof obligations for each step guarantee the preservation of some classes of properties [7]. We use the Event-B formal method. We start with a big step semantics specified by an Event-B machine. Through a sequence of refinements, expressed as Event-B refinement machines, we end up with a code generator and an interpreter which implements a small step semantics preserving the properties of the initial big step semantics, i.e. the sequences of reactions observed when executing the concrete machine were allowed by the big step semantics.

The rest of the paper is organized as follows: in the next section, we give an overview of the Event-B language. In Section 3, we present the development of a generic compiler for synchronous languages. Finally, in Sections 4 and 5, we discuss related and future work.

2 EVENT-B

The Event-B method allows the development of correct by construction systems and software [1]. To achieve this, it supports natively a formal development process based on a refinement mechanism with mathematical proofs. Figure 1 illustrates a refinement step where a machine M0 using a context C0 (the sees edge) is refined (the refines edge) by a machine M1 using an extension C1 of C0 (the extends edge). Contexts define abstract data types through sets, constants and axioms while machines define symbolic labelled transition systems. The state of a transition system is defined as the value of machine variables. Labelled transitions are defined by events specifying the new value of variables while preserving invariant properties. Proof obligations for wellformedness and invariant preservation are automatically generated by the Rodin tool [18].

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They can be discharged thanks to the available automatic proof engines (CVC4, Z3, ...) or through human-assisted proofs.

2.1 Notations

For the most part, Event B uses standard set theory and its usual set notation. As a matter a fact, in Event-B, an array and a function are both considered as set of couples. Some notations are specific to Event B:

- **pair construction**: pairs are constructed using the maplet operator \( \leadsto \). A pair is thus denoted \( a \leadsto b \) instead of \( (a,b) \). The set of pairs \( a \leadsto b \) where \( a \in A \) and \( b \in B \) is denoted \( A \times B \).
- **sets A X B**: is a relation. A relation \( r \) has a domain \( \text{dom}(r) \) and a codomain \( \text{ran}(r) \). When a relation \( r \) relates an element of \( \text{dom}(r) \) with at most one element, it is called a function. The set of partial functions from \( A \) to \( B \) is denoted \( A \rightharpoonup B \), the set of total functions is denoted \( A \rightarrow B \). The image of a set \( A \) by a relation \( r \) is denoted \( r[A] \). Last, the relational composition is denoted by the ; infix operator. Suppose \( f \in A \rightarrow B \) and \( g \in B 

2.2 Preliminary example

As a preliminary example for section 3.1, Figure 2 illustrates basic labelled transitions systems [4]. The static description of transition systems is given through the context \( \text{cLTS} \). \( \text{cLTS} \) introduces State and Label as abstract sets. Then, the set of initial states is given through the constant \( I \), and the transition relation through the constant Trans. The description of the dynamics of a transition system is given by the machine \( \text{mLTS} \). First, we represent the state space by the variable \( \text{st} \in \text{State} \). The dynamics is specified through two events: the \( \text{INITIALISATION} \) event which chooses non-deterministically a state among the elements of the set \( I \) and the next event which also chooses non-deterministically a label and a transition having such a label.

Figure 3 illustrates an Event-B refinement of the preceding machine \( \text{mLTS} \). Intuitively, we model the introduction of an internal event that will take control for some time after which the next event becomes enabled again. For such a purpose, the machine \( \text{mLTS}_r \) refines \( \text{mLTS} \) and introduces a new variable \( \text{cpt} \) : a counter that will be initialized non deterministically. When the counter reaches the zero value, next is enabled. The refined machine refines the events \( \text{INITIALISATION} \) and \( \text{next} \) and introduces the new event \( \text{silent} \) which is similar to \( r \) events of CCS [12]. We notice that in the new (refining) next event, the enabled guard is strengthened by the conjunct \( \text{cpt} = 0 \). Moreover, this new next event provides a witness of the parameter of the old (refined) next event: the parameter \( lr \) is said to be the witness of the parameter \( l \). When the new next event has been executed, \( \text{cpt} \) is again assigned a positive value. In Event-B, the proof obligations enforce a weak refinement semantics [12]. The purpose of the \( \text{variant} \) clause is to enforce that introduced events will not take control forever. Usually, the variant belongs to a well ordered domain and is decreased by newly introduced (internal) events declared to be convergent.

3 A GENERIC COMPILER DEVELOPMENT FOR SYNCHRONOUS LANGUAGES

In this section, we elaborate a generic refinement-based approach for the compilation of synchronous languages. Starting from the so-called labelled synchronous transition systems, we propose a sequence of refinement steps, illustrated by Figure 4, leading to code generation. Thanks to the refinement-based approach, the generated code is known to preserve by construction the initial semantics provided all the proof obligations have been discharged.

3.1 Labelled synchronous transition systems

In the synchronous context, executions are seen as sequences of reactions [17]. A reaction \( r \) assigns values to the subset of variables said to be present: undefinedness of \( r \) on a variable \( v \) means absence of \( v \) in \( r \). We suppose a fixed domain \( \mathcal{D} \) for the values and define a reaction over the set of variables \( V \) as a partial function from \( V \) to the domain \( \mathcal{D} \), i.e., \( \text{Reaction}(V) = V \rightarrow \mathcal{D} \). We remark that the usual definition of a reaction is a total function to \( \mathcal{D} \cup \perp \) where \( \perp \) becomes either a specific value or undefinedness.
With respect to the Signal language, we essentially do not consider
The semantics of Signal can be defined as an LSTS where a state
Definition 3.1 (LSTS). A Labelled Synchronous Transition System
3.1.1 Signal and its LSTS semantics. In order to illustrate the
definition each variable at most once:
• \( x_1 = x_2 \) when \( x_3 \) undersampling
• \( x_1 = x_2 \) default \( x_3 \) deterministic merge
• \( x_1 = f(x_2, x_3) \) instantaneous function
• \( x_1 = x_2 \) if initial delay
With respect to the Signal language, we essentially do not consider
local variables. This aspect is not relevant for our presentation.
the semantics of Signal can be defined as an LSTS where a state
is a mapping from memorized (or persistent) variables to values.
A variable is memorized if it is introduced by a delay equation.
In order to define the denotational semantics of Signal, we introduce
the function \( R_m \) where \( m \) is a memory state. It takes as parameter
a Signal equation and returns the set of allowed reactions. We also
suppose the existence of a constant \( T \in \mathcal{D} \) representing true.
• \( R_m[x_1 = x_2 \text{ default } x_3] = \{
  \{ r \in V \mapsto \mathcal{D} \} \mid \forall u \cdot x_1 \mapsto u \mapsto r \Rightarrow \ 
  (x_2 \mapsto u \mapsto r \land x_3 \mapsto T \mapsto r) \} \)
3.1.3 Example. Consider the Signal program which increments \( x \) and restarts it from the signal \( r \) each time it occurs:

\[
\text{process INCR = } (\{\text{integer } r; \text{integer } x;\} \\
| \{ x : = x + 1 \} \\
| \{ x : = r \text{ default } nx \} \\
| )/\text{px, nx}
\]

If we keep local variables visible, the associated rLSTS is defined as follows:

- \( \mathcal{V} = \{ r, x, px, nx \} \)
- \( \mathcal{M} = \{ px \} \)
- \( \mathcal{I} = \{ px \rightarrow 0 \} \)
- \( \mathcal{N} = \{ px \rightarrow x \} \)
- \( \mathcal{C}(m) = R_m[px = x \text{ init } 0] \) \( \cap \) \( R_m[nx = px + 1] \) \( \cap \) \( R_m[x = r \text{ default } nx] \)

Figure 5 is a partial run of this example. A run is an alternating sequence of states (ovals) and transitions labelled by reactions (vectors). In the first reaction, \( r \) is absent and \( x \) is one more than the memorized value. The second transition is labelled by a reaction where \( r \) is present. Thus \( x \) takes the value of \( r \). In the third transition, \( r \) is chosen to be absent and \( x \) is incremented. The reaction could also be empty (all variables being absent). However, this will be forbidden in the remaining of the paper to avoid Zeno behaviors.

3.1.4 Event-B representation. First of all, we define the context \( \text{cMeta} \) (Fig. 6) containing the various fields of the rLSTS we will consider. With respect to definition (3.2), we have added one parameter of type \( \mathcal{P}(\mathcal{V}) \) to the function \( \mathcal{C} \) denoting the constraint problem of which reactions should be solutions. It contains the set of variables that are constrained and is now the full set \( \mathcal{V} \). This parameter will be modified to specify the incremental computation of the reaction in the next refinement.

Second, we define the semantics of the given rLSTS through an Event-B machine which specifies the Labelled transition system associated to the rLSTS through the state variable \( m \) initialized using \( I \). The state evolves through the \text{bigStep} event parameterized by a reaction \( r \). The reaction is accepted if it is a solution of the constraint system.

This \text{bigStep} event is intended to be split into small steps that compute reactions incrementally.

```plaintext
Figure 6: The Event-B context \( \text{cMeta} \)
```

```
context \text{cMeta} 
sets \( \mathcal{V} \mathcal{D} \)
constants \( M I N C \) 
axioms 
@\text{fVat finite}(\mathcal{V}) 
@I \in M \rightarrow \mathcal{D} 
@N \in M \rightarrow \mathcal{V} 
@C \in (M \rightarrow \mathcal{D}) \rightarrow (\mathcal{P}(\mathcal{V}) \rightarrow \mathcal{P}(\mathcal{V} \rightarrow \mathcal{D}))
```

end

```plaintext
machine \text{mBigStep} sees \text{cMeta} 
variables \( m \) // memory 
invariants 
@m_\text{ty} m \in M \rightarrow \mathcal{D} 
events 
\text{event \text{INITIALISATION}} then 
@m_\text{init} m = I 
end 
\text{event \text{bigStep}} any \( r \) where 
@r \in \mathcal{P}(\mathcal{V}) \rightarrow \mathcal{D} // r is a reaction 
@r_\text{ty} r \in \mathcal{C}(m)(\mathcal{V}) // r is a solution of the constraint system 
@n_\text{ty} r \neq 0 // to avoid Zeno behaviors 
then 
@m m = m \text{ end } (\mathcal{N} x) // memory update 
end 
end
```

3.2 Solved form (context \text{cMeta SF}) and small step semantics (machine \text{mSmallStep})

This refinement aims at splitting the big step which specifies what are valid reactions. The small steps should compute one variable at a time. To make it possible, the constraint system should be put in solved form.

3.2.1 Solved form. We define the constraint problem as a set of equations supposed to be in triangular form: a left hand side variable is said to depend on right hand side variables and this dependency relation should be acyclic. In fact, two informations must be computed: the presence of a variable, and if present, its value. The value of a present variable is either defined by a formula, or read from input (indeterminate), or read from memory. Thus, we first introduce the constant \text{isPresent} together with the dependency relation \text{dep}. The function \text{isPresent} takes as parameters a variable and a partially computed reaction and returns a set of
booleans which indicate if the variable is allowed to be present or absent, or both. isPresent returns either \( \text{TRUE} \), either \( \text{FALSE} \), or \( \text{[TRUE,FALSE]} \) to encode indeterminacy, as it is the case for master clocks of Signal programs [8]. Concerning dep, it maps each variable to the set of variables on which its presence depends. Moreover, a variable which does not depend on other variables can be chosen to be present or absent.

\[
\text{Figure 7: cMeta$_SF$ (part I)}
\]

Now, we introduce the eqns function for value computation, together with the dependency relation dep. Contrary to presence, values are specified in a functional way. Indeterminacy of values is inferred from the absence of equations for some variables. The function eqns takes as parameters the considered variable and returns its value for the currently computed reaction. This function is partial at several orders: it is defined for a subset of non memory variables. Accepted reactions (which are partial functions) should at most give a value to a variables on which the first argument depends. These reactions are those that allow the presence of the variable.

\[
\text{Figure 8: cMeta$_SF$ (part II)}
\]

Figure 9 introduces the link between presence and the existence of a value for a variable: a partial solution \( ps \) allows the presence of a variable \( v \) if and only if the partial solution \( ps \) is in the domain of eqns \((v)\).

\[
\text{Figure 9: cMeta$_SF$ (part III)}
\]

We suppose that for any variable \( v \), \( \text{dep}(v) \) is a subset of \( \text{dep}(v) \): the value of \( v \) (depending on variables of \( \text{dep}(v) \)) can be computed once we know its presence (depending on variables of \( \text{dep}(v) \)). Then, we add properties stating that dep is transitively closed and irreflexive, thus a strict partial order.

Now, we define the constraint problem \( C_{m,S} \) for a memory state \( m \) and a subset \( S \subseteq V \) of seen variables. A solution \( r \) is a partial function on \( S \), undefined meaning absence. It is constrained by three properties (Fig. 11):

- The presence or the absence of a variable \( v \) within a reaction \( r \) should be allowed by the set returned by the function isPresent where \( r \) is restricted to variables on which \( v \) depends.
- The restriction of \( r \) to memory variables is given by the memory state.
- The value of a present variable \( v \) associated to an equation is given by this equation taken as parameter \( v \) and a restriction of \( r \) to variables it depends on.

\[
\text{Figure 10: cMeta$_SF$ (part IV)}
\]

\[
\text{Figure 11: The constraint problem}
\]

3.2.2 Memory access. We introduce dedicated variables, from the set NextMem, used as markers for memory updates. The injective function write associates to each memory variable \( v \), its memory write event. This event should be fired after all reads of \( v \) and after the computation of the next value of \( v \) given by the variable \( N(v) \). The presence of the write event depends on the presence of the next value. It is actually present if the next value is present.

\[
\text{3.2.3 Small step semantics. Given a constraint system in triangular form, it is possible to compute reaction variables one by one through so-called small steps. In the same way, the memory can be updated incrementally once the current value of a variable becomes useless in the current cycle: it will not be read anymore. For this purpose, we introduce a refinement machine with three new variables: the partially updated memory \( pm \), the partial solution \( ps \) and the set of processed variables done.}
\]

\[
\text{machine mSmallStepRefines mBigStep sees cMeta$_SF$}
\]

variables \( pm \) // partially updated memory
\( ps \) // partial solution
done // processed signals

invariants
\( \text{@pm \in M \rightarrow } D \)
\( \text{@ps \in M \rightarrow } D \)
\( \text{@done \in M \rightarrow } D \)
These variables satisfy the following invariant properties:

- The restrictions of pm and ps to unprocessed memorized variables are identical functions.
- ps is a partial solution restricted to variables of done.
- If a variable is present, variables of which depends its value (given by dep) have been processed.
- If a variable has been processed, variables on which depends its presence (given by pdep) have also been processed.
- The partial solution is not empty as soon as one variable has been processed, i.e. the first processed variable should be marked as present.
- The memory is partially updated by the next value of memory variables of which the write event has been processed.

Initially, the partial solution and the set of processed variables are empty. Memory gets the rLSTS initial value.

Small steps determine the presence and compute the value of individual variables. These steps must be transparent from the point of view of the abstract machine: they do not modify previously introduced variables and should apply a finite number of times. This is why they are declared convergent, which means that they must decrease the variant. Several convergent events are defined. They respectively process absent variables, present variables defined by an equation, memory accesses or undetermined variables (considered as inputs or parameters). Due to space restrictions, two cases are presented: equational variables and absent variables. Other events are related to memory read and write variables, parameters (non memory and non equational variables).

The following listing manages variables defined by an equation. The event selects such a variable v which has not been processed before. Variables on which it depends should have been processed and v is allowed to be present. Then, its binding to the value given by the equation is inserted into the partial solution and v is added to done. It follows that the variant decreases.

The following listing manages absent variables. Contrary to present ones, they are processed by blocks. This will allow optimisations in the generated code. So, we consider a non empty set of variables V such as variables on which they depend are either already processed or in V. Absence should be allowed for all these variables. Note that this permission is checked by supposing variables of V are absent as unprocessed variables are seen as absent. Lastly, the partial solution should not be empty to ensure that the computed reaction will not be the empty one.

### Invariants

```plaintext
@psm ∀v ∈ M: done ⇒ pm(v) = m(v) // pm equals m on unprocessed variables
@psm ok ps ∈ C(m(done)) // ps is the restriction to done of a reaction
@seen_val ∀v ∈ done ⇒ dep(v) ⊆ done
@seen_pr ∀v ∈ done ⇒ pdep(v) ⊆ done
@done_ne done = ∅ ⇒ ps = 0 // forces non empty reactions
@pm_m pm m = (write⁻¹(done)m) // pm is a partial update of m
```

### Events

- **INITIALISATION**
  ```plaintext
  event INITIALISATION
  then
  @ist ps ps := ∅
  @inst done done := 0
  @inst pm pm := I
  end
  ```

- **smallStep**
  ```plaintext
  event smallStep // processing of a present variable
  any v where
  @v ∈ dom(eqns) \ done
  @dep(v) ⊆ done // thus pdep(v) ⊆ done
  @ps ps(v) := eqns(v) \ done
  then
  @done done := done ∪ {v}
  ```

- **smallStep_P**
  ```plaintext
  event smallStep_P // processing of a present variable
  any v where
  @v ∈ dom(eqns) \ done
  @dep(v) ⊆ done // thus pdep(v) ⊆ done
  @ps ps(v) := eqns(v) \ done
  then
  @done done := done ∪ {v}
  ```

- **convergent event**
  ```plaintext
  convergent event smallStep_m // processing of absent variables
  any V where
  @V \ done
  @m V \ done = 0 // ensures that the variant decreases
  @d ∀v ∈ V ⇒ pdep(v) ⊆ done ∪ V
  // variables of V are allowed to be absent
  @abs ∀v ∈ V ⇒ FALSE = isPresent(v) \ pdep(v) \ ps
  @ps ps # 0
  then
  @done done := done ∪ V
  end
  ```

When all the variables have been processed by the small steps, we can refine the big step of the abstract machine and prepare the next round by resetting the variable done.

### 3.2.4 Example

Given the Signal example previously introduced in Paragraph 3.1.3, we add the write event marker ps_w for the memory variable px. Figure 12 shows how the declared constants introduced in the solved form can be instantiated. The graph represents the functions pdep and dep. All the antecedents with plain arrows of a node x form its dependency for pdep. dep is obtained by adding the dashed arrows.

Concerning the isPresent function, written isP in the figure, it allows presence and absence for r. If r is actually present, px must be present; otherwise px presence does not care. Then the authorization of presence of variables nx, x and px_w is determined by the actual presence of px and x.

Concerning eqns, it associates a value to px and x by accessing to the value of variables they depend on.

![Figure 12: Solved Form level (example)](image-url)
3.3 Introduction of clocks (machine mSmallStep_CC)

We now associate clocks to variables. The role of a clock is to manage presence. Thus, a variable will be present if and only if its clock is present. For this purpose, the set Clock is declared as a subset of variables supposed to be neither memory variables nor defined by an equation: they are not valued. Each variable has a clock (defined by the function cP). The clock of a clock is itself.

A new refinement can now be defined. It differs from the previous machine by the fact that presence of clocks are tested through the isPresent function while presence of valued variables comes to the presence of their clocks. One invariant is added. It states that for a processed variable, its presence and the presence of its clock are equivalent. The following listing illustrates the transformation of the two events. For equational variables: cP(v) presence is now checked. For absence, constraints on V are strengthened to ensure clocks are also selected and that they are absent. The big step is unchanged.

3.4 Clock hierarchy (machine mSmallStep_SUB)

The goal of this refinement is to introduce the clock inclusion relation: if a clock is absent, sub-clocks are also absent and it is not useful to check them at runtime. For this purpose, we introduce the relation sub represented by a function associating to a clock the set of clocks it subsumes (or contains). This relation should be compatible with the pdep relation: a clock depends on clocks containing it. Subsumption declarations should be consistent with presence semantics as given by the isPresent function.

3.3.1 Example. Given our working example, we introduce the clocks \( \hat{r}, \hat{x}, \hat{x_n}, \hat{px}, \hat{px_w} \), thus defining the cP function by \( cP(v) = \hat{v} \).

We have chosen to take the same clock for \( x \) and \( nx \). This is allowed by the axioms since their respective presence are defined in the same way. The dependencies between the variables are now given by the following graph:

\[
\hat{r} \xrightarrow{r} \hat{px} \xrightarrow{px} \hat{x} = \hat{nx} \\
\hat{nx} \xrightarrow{x} \hat{px_w} \xrightarrow{px_w} \]

These new arrows define a subgraph of the dependency graph and should satisfy the axiom sub_present. For instance, let us consider the bendarrow between \( \hat{px} \) and \( \hat{x} \). Given a partial solution \( ps \) defined at least on \( pdep(px) \) and not on \( px \) we must check the implication between each presence property. This property is satisfied since the right side of the implication is True because \( px \notin dom(ps) \).
3.5 Sequential code generation (machine mSmallStepSeq)

This refinement step is the key point of the verified compiler generation. It consists in separating actions that can be performed at compile time (statically) from actions performed at run time, the link between them being the generated code. The principle is the following:

- At the abstract level, an event is fired if its parameters satisfy static properties (linked to the dependency relations) and dynamic properties (linked to presence / value) of variables in the partial reaction already built.
- To each such event is associated a convergent event that will select its parameter depending only on static properties and store it in the sequence of instructions built by the compilation phase. These events constitute the code of the compiler.
- The abstract event itself is refined by an event which simulates the created instruction by testing and updating the dynamic state.
- The correction of this separation relies on the fact that static properties should be sufficient to select event parameters. It is ensured if we suppose that, given parameters satisfying the static properties, there exists an event of which dynamic properties are also satisfied.

These principles are applied to small step events. Such an event may be fired on a variable if previous variables (for dep or pdep partial orders) have been processed and if some conditions depending on the partial reaction (presence of the clock for example) are satisfied. Building the reaction is left for the dynamic phase. However, variable selection, i.e., topological sort, can be performed at compile time. As a result, we get a sequence of statements specifying the event to fire and the variable to be selected. Statements are declared in the cMetaSeq context:

context cMetaSeq extends cMeta_SUB
sets Stmt
constants
GetPresence // check clock presence
GetP // read parameter statement
ReadM // read memory statement
WriteM // write memory statement
Assign // assignment statement
axioms
@get partition((Stmt,GetPresence), GetP, ReadM, WriteM, Assign)
end

3.5.1 The compilation events. The compiler encodes machine instructions numbered from 0 to N – 1 through two variables: stmt for the name of the statement and arg for the concerned variable. In order to take into account the clock hierarchy, we add a jump pointer (iff read if false) to each presence test. It indicates the continuation in case the test fails. So, we introduce the following variables:

machine mSmallStepSeq refines mSmallStep_SUB sees cMetaSeq
variables pm // memory dynamic, inherited
ps // partial reaction dynamic, inherited
N // size of code (static)
pc // program counter (dynamic)
stmt // statements (static)
arg // argument of statement (static)
iff // argument of statement (static)
\n\nconvergent event sta_smallStep_P
any v where
@pv v ∈ dom(eqns) \ ran(arg)
@ldep(v) \ ran(arg)
thend
@N N := N + 1
@arg arg(N) := v
@assign stmt(N) := Assign
@iff iff := iff + 1
\{i | i < stmt(\{GetPresence\}) ∧ iff(i) \ N ∧ \ cP(v) \ sub(arg(i)) \ \ cP(\{arg(i)\}) \ \ N+1\]
end

3.5.2 The interpretation of generated events. The second event recognizes the generated instruction and performs the dynamic part of the refined event: the presence of the argument of the instruction is checked and the reaction is updated. Note that presence check could be removed: within the bloc of code skipped by the iff pointer, the presence condition is known to be satisfied. Due to lack of space, this optimization is not detailed here.

event smallStep_P refines smallStep_P where
@pc pc < N // end of generated code reached

The invariant specifies the link between old and new dynamic variables (done and pc) and between generated instructions and expected properties of the variable selected by the old small step events:

- processed variables (the set done) are arguments of previous statements in the code sequence,
- dependencies are arguments of previous statements
- arguments of GetPresence are clocks
- arguments of GetParam are parameters
- arguments of ReadMem are memorized variables
- arguments of WriteMem are memory write events
- arguments of Assign are equational variables
- iff pointers are forward pointers
- iff pointers skip subsumed clocks

The principles are applied to small step events. Such an event may be fired on a variable if previous variables (for dep or pdep partial orders) have been processed and if some conditions depending on the partial reaction (presence of the clock for example) are satisfied. Building the reaction is left for the dynamic phase. However, variable selection, i.e., topological sort, can be performed at compile time. As a result, we get a sequence of statements specifying the event to fire and the variable to be selected. Statements are declared in the cMetaSeq context:
Absence being managed through the update of the iff pointers during the generation of all small step statement related to presence, only remains the dynamic part which consists in skipping a block of instructions in case a presence test fails. The new event declares the values of the parameters of the refined event.

```plaintext
@arg stmt(pc) = Assign // current statement is an Assign @pc stmt(pc) = GetPresence with @v = arg(pc) // the processed variable is the argument of the statement then @ps ps(arg(pc)) := exprs(arg(pc))(dep(arg(pc)) * ps) @npc pc := pc + 1

end
```

At last, the big step event is modified to take into account the new state variables: the compilation should have been performed before, the small steps should have been executed. It has to be noted that, except for this last event, compilation and execution of generated instructions can be interleaved during the computation of the first reaction of the run. This property is partially lost if presence test optimization is considered: if execution jumps to the end of the code as a consequence of a presence test failure, the compiler could insert an optimized statement (without presence check) at this point.

```plaintext
event smallStep_mA refines smallStep_mA when @pcN pc < N // end of generated code reached @pc stmt(pc) = GetPresence @pc pcaN > 0 // it is the first statement @abs FALSE ∈ isPresent[ arg(pc) = pdep[ arg(pc)] * ps] // absence allowed with @v = arg(pc) // the processed variable @V = arg(pc) .. iff(pc) // the set of variables to be skipped then @npc pc := iff(pc) // jump to next block

end
```

### 3.6 Development overview

All the preceding development has been performed under the Rodin framework for Event-B. For all the development (contexts and machines), the tool generated 17 proof obligations for contexts and 239 proof obligations for machines. Most of the proof obligations could be discharged automatically thanks to the SMT solvers (CVC3-4, Z3, veriT). Some proof obligations needed manual quantification instantiation and case analysis.

We remark that unlike usual assistant theorem provers [9, 14, 15], elementary proof obligations are generated automatically, e.g., the preservation of each conjunct of the invariant by each action of each event. However, concerning Rodin, strategy languages are still missing.

### 4 RELATED WORK

In this section we review some related works. In the following, we position our work with respect to them.

#### 4.1 Clock calculus

Here, we have adopted an axiomatic approach: in the context cMeta_SUB, we have specified a subsumption relation compatible with data dependency. The work of [3] can be seen as a concretization of this specification. As a future work, we intend to validate some of these constructions by taking into account the language to which they apply.

#### 4.2 Translation validation

Translation validation has been pioneered by A. Pnueli et al. [16]. Recently, [13] have pushed forward this approach for the Signal language. The basic idea is that each individual translation is followed by a validation phase which verifies that the produced code implements the source code. Unlike this approach, here we are interested in the compilation approach: which verifies the compiler once for all. It follows that any produced code is guaranteed to be correct.

#### 4.3 CompCert

The CompCert project [11] aims at the development of a formally verified C compiler. This work is far more complex than our one and could be used as a back end of a synchronous language to C compiler. However, one point is interesting to be discussed and is put forward in the CompCert project. It is related to what means...
for a compiler to be correct? It should preserve the semantics of the source code, i.e., preserve some observable behaviors, which is ultimately expressed as proving a forward simulation for safe programs property. In this paper, a safe program is a program that can be put in solved form. Forward simulation is a consequence of refinement. Observable behaviors are sequences of reactions. Thus, Event-B brings us a methodology to incrementally build a similar proof (for a much less complex language).

### 4.4 Constructive semantics

A so called constructive semantics has been proposed for various synchronous languages [21]. This semantics aims at specifying how to compute output variables from inputs for each reaction. As in our case, small steps are proposed to solve constraints upon presence and values of the variables of the program. Specific rules are provided for various synchronous languages. The main difficulty of this approach, which is linked to the considered languages, is to find a trigger, i.e. a signal of which presence can be enforced in order to determine the presence and value of all other signals without failing to find a global solution when solving incrementally the constraint system. The authors define several classes of processes (reactive, deterministic, isochronous, ...) ensuring convergence properties. In this paper, we do not address these aspects and suppose that constraints have been put in solved form, which guarantees the ability to incrementally build a solution and has the same consequences as isochrony [21], i.e., the ability to non deterministically select a non empty subset of triggers to be present. Then, we mainly focus on refinement-based methodology to extract an optimizing compiler and the interpreter of the compiled code from the small step semantics which is itself shown to conform with the big step semantics. Furthermore, our framework is independent from a specific synchronous language.

### 5 CONCLUSION

In this paper, we have presented the development of a generic compiler for synchronous languages. Our main contribution was to structure such a development as successive refinement steps. Through these steps, we have introduced progressively the main ingredients of such a compilation process: presence of variables, clock calculus, optimization and code generation. Properties preservation was ensured thanks to the semantics of refinement supported by the Event-B formal method. With respect to the expression of the properties, set theory has revealed well suited to express: as well atomicity in high level specifications thanks to the composition of relations, as well the progressive approach thanks to partial functions. The proof obligations have been generated by the Rodin platform. All the proofs have been done. These proofs were made easier by the automatic SMT provers.

We envision to carry on this work with respect to two directions: first we would like to instantiate this generic compiler by well known synchronous languages as [5, 20]. Second, also, we would like to adapt code generation for targeting concurrent architectures [2].

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**REFERENCES**