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# A formal specification and prototyping language for multi-core system management.

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**Abstract**—We relate the experience of a defining a formal domain specific language (DSL) for the construction and reasoning about OS-level management logic of multi-core systems. The approach is based on a novel, iterative development principle where results of prototyping studies feed back into the next language revision. We illustrate the DSL with several examples of executable scripts.

## I. INTRODUCTION

With the end of processor frequency growth, scaling into the direction of multi- and many-core systems has become the primary way to translate the advances in manufacturing technology into increased computer performance. This change has already started to affect operating systems and application software design. Clearly, we have to plan now for the many-core systems of tomorrow.

In this paper we propose a method and toolkit that has grown around an effort to build a database of knowledge related to multi-core systems. The exercise started with a simple compilation of fact using a natural language. The result turned out to be inadequate due to many ambiguities, hidden assumptions and subtle inter-relations between concepts. A decision was made to transfer into a formal specification language. We have chosen Event-B modelling language [1] mainly for our previous experience with it. The formal specification of the knowledge database resolved the known ambiguities and exposed many omissions and inconsistencies. The effort took several months and resulted in a substantial model with hundreds of properties. It was expected that the model is consulted when designing or writing OS-level software for multi-core system. It turned out that a formal model of this scale is a difficult read and not a practical blueprint to relate with actual software. Moreover, there was a lack of confidence in the model completeness and liveness properties - conditions that are impractical to address with static theorem proving and turned out impossible to delegate to a model checker due to model complexity and scale.

To address these weaknesses while staying in the formal domain, it was decided to build a control flow extension to Event-B that would "drive" (to take the term from a conceptual inspiration - the CSP||B technique [2]) the original Event-B model. The pursuit of this goal led to the development of a novel approach to the construction of formal imperative-style domain-specific languages where the heart of the language (its domain-specific part such as commands, variables and constants) is designed in a separate formal notation (Event-B) while the "glue" part (control flow constructs like *if* and *while*) are generic and reused by all DSL instances. The method and the accompanying tools are referred to as DSL-Kit.

The most profound implication is the facilitation of iterative design: design prototypes realised in a DSL would often highlight deficiencies in the DSL itself (lack of progress,

missing concepts); this may be addressed by going back to Event-B specification of the domain, doing necessary changes, proving them correct and *automatically*, with the help of a translation tool, transferring the result into DSL-Kit to construct a new instance of domain DSL. DSL-Kit itself offers fairly reach reasoning facilities. On top of this, it can execute (with many limitations) DSL specifications and a custom state visualisation may be plugged in to facilitate debugging and design comprehension.

The first part of the paper (Section III) is a narrative build around an Event-B model that attempts to capture the essence of multi-core systems from the mostly software and OS perspective. The model does not include any notions of control logic. It rather gives a concrete definition of the subjects of the prospective control system and attempts to explain their inter-relationships; it also describes the general life cycle of a multi-core system.

The second part (Section IV) describes the transition in the DSL-Kit environment and relates the initial experience with the design of an adaptive run-time controller. We show how the domain model may be turned into a formal virtual machine over which control logic executes. The reasoning style also changes from refinement proofs and inductive verification of safety invariant of Event-B to the verification of imperative programs in the style of Floyd-Hoare logic [3].

## II. EVENT-B

### A. Event-B

The basis of our discussion is a formalism called Event-B [1]. It belongs to a family of state-based modelling languages that represent a design as a combination of state (a vector of variables) and state transformations (computations updating variables).

An Event-B development starts with the creation of a very abstract specification. A cornerstone of the Event-B method is the stepwise development that facilitates a gradual design of a system implementation through a number of correctness-preserving *refinement* steps. The general form of an Event-B model (or *machine*) is shown in Figure 1. Such a model encapsulates a local state (program variables) and provides operations on the state. The actions (called *events*) are characterised by a list of local variables (parameters)  $vl$ , a state predicate  $g$  called *event guard*, and a next-state relation  $S$  called *substitution* or event *action*.

Event guard  $g$  defines the condition when an event is *enabled*. Relation  $S$  is given as a generalised substitution statement [4] and is either deterministic ( $x := 2$ ) or non-deterministic update of model variables. The latter kind comes in two notations: selection of a value from a set, written as

```

MACHINE M
SEES Context
VARIABLES v
INVARIANT  $I(c, s, v)$ 
INITIALISATION  $R(c, s, v')$ 
EVENTS
   $E_1 \triangleq$  any  $vl$  where  $g(c, s, vl, v)$  then  $S(c, s, vl, v, v')$  end
  ...
END

```

Fig. 1. Event-B machine structure.

$x : \in \{2, 3\}$ ; and a relational constraint on the next state  $v'$ , e.g.,  $x : | x' \in \{2, 3\}$ .

### III. DOMAIN MODEL

#### A. Event-B domain model specification

a) *Cores*: A system contains a number of cores; at any moment a core may be operating or switched off. The set of all cores in a system is defined by a finite and non-empty set **CORES**. Current core status (on or off) is given by function variable *status*:

$$status \in \mathbf{CORES} \rightarrow \mathbf{STATUS}$$

where **STATUS** = {**ON**, **OFF**}. At this level, we may observe a core being switched on or off, as captured by the following two events:

```

on  $\triangleq$  any  $c$  where  $status(c) = \mathbf{OFF}$  then  $status(c) := \mathbf{ON}$  end
off  $\triangleq$  any  $c$  where  $status(c) = \mathbf{ON}$  then  $status(c) := \mathbf{OFF}$  end

```

b) *Frequency and voltage*: Two essential characteristics of a running core are the voltage of its power supply and the clock frequency. These are the principal attributes used to control core performance, power and reliability. The following partial functions define frequency and voltage of a core; these are undefined for cores switched off.

$$\begin{aligned}
freq &\in \mathbf{CORES} \leftrightarrow \mathbb{N} & vdd &\in \mathbf{CORES} \leftrightarrow \mathbb{N} \\
\text{dom}(freq) &= status^{-1}\{\{\mathbf{ON}\}\} & \text{dom}(vdd) &= status^{-1}\{\{\mathbf{ON}\}\} \\
freq &:= \emptyset & vdd &:= \emptyset
\end{aligned}$$

Since all the cores are initially off, voltage and frequency functions are initially empty. Voltage and frequency attributes are changed dynamically by event *core\_dvfs*:

```

core_dvfs  $\triangleq$ 
any  $c, f, v$  where
   $status(c) = \mathbf{ON} \wedge \dots$ 
then
   $freq(c) := f \parallel vdd(c) := v \parallel \dots$ 
end

```

The dots stand for omitted clauses; in a refined version, the event checks whether frequency/voltage pairs are correct for a given core, and also whether power budget and temperature constraints are satisfied for the new settings. Switching a core on also necessitates setting some initial voltage and frequency values..

#### B. Heat generation

Heat exchange happens between a core and the environment and a core and its neighbouring cores. The environment is assumed to possess infinite heat capacity so that its temperature is not affected by heat exchange.

For some core  $c$ , the amount of heat generated per unit of time is determined by core frequency  $f$  and core voltage  $V$ , according to the following law:

$$\frac{\partial H_c}{\partial t} = C_c f c V_c^2$$

where  $C_c$  is a core-specific constant. Heat rate  $\frac{\partial H}{\partial t}$  and heat exchange laws define core temperature delta over a time period. For any given core, given its heat capacitance, we can determine the time duration necessary to increase core temperature by one Kelvin. Alternatively, considering some fixed time period, we can determine temperature change caused by the heat rate during the period.

The following event controls core temperature change. At this step, it is un-timed and time will be added in a later refinement step.

```

core_temp  $\triangleq$ 
any  $c$  where
   $status(c) = \mathbf{ON}$ 
then
   $temp(c) := \max(\{temp(c) +$ 
     $\mathbf{HRATE}(c \mapsto vdd(c) \mapsto freq(c)) +$ 
     $\sum_{n \in \mathbf{NHB}(c)} \mathbf{NRATE}(n \mapsto c \mapsto vdd(c) \mapsto freq(c)) -$ 
     $\mathbf{ERATE}(c), 0\})$ 
end

```

In the above **HRATE**(...), **NRATE**(...) and **ERATE**(...) are the heat rates times time delta for the heat gained through resistive heating, heat exchange with neighbour cores and heat loss to the environment. When a core overheats, the system immediately (we shall clarify the meaning of immediacy with the introduction of time) switches off the core.

```

core_shutdown  $\triangleq$ 
any  $c, t$  where
   $status(c) = \mathbf{ON}$ 
   $temp(c) > \mathbf{CORE\_TEMP\_CRIT}(c) \wedge \dots$ 
then
   $status(c) := \mathbf{OFF} \parallel \dots$ 
end

```

We shall later see that shutting down a core cancels all the jobs and stops all the running threads. All the computation progress is irrecoverably lost.

c) *Threads*: A thread is a basic concurrency and program structuring unit in our model. By a thread we understand a potentially infinite sequence of commands that are continuously or intermittently executed by a core, perhaps switching between cores during its lifetime. At any given time, there is some number (potentially zero) of threads in the system:

$$threads \subseteq \mathbf{THREADS}$$

where **THREADS** is the universe of threads. A thread may be assigned to a core and then it is said to be running or *scheduled*. The thread/core association is functional and partial in the domain:

$$affinity \in threads \rightarrow \mathbf{CORES}$$

It is only possible to schedule a thread on a running core:

$$\text{ran}(affinity) \subseteq status\{\{\mathbf{ON}\}\}$$

Note that *affinity* is functional in one direction only (i.e., it is not injective); indeed, it is possible to map several threads onto the same core, for instance,

$$t, h \in threads \wedge t \neq h \wedge affinity(t) = affinity(h)$$

describes a situation where distinct threads  $t, h$  are running on a same core. In our model we choose to fix the lowest time resolution at the scale of  $\sim 1$  millisecond; thus for a sequence of events with overall time smaller than this limit, we see all

the events happening at the same time or concurrently. This is a standard abstraction technique and it allows one to conduct a formal refinement to a higher time resolution.

A new thread may be added to the system and, at some point, it may be destroyed:

$th\_start \triangleq$  any  $t$  where  $t \notin threads$  then  $threads := threads \cup \{t\}$  end  
 $th\_stop \triangleq$  any  $t$  where  $t \in threads$  then  $threads := threads \setminus \{t\}$  end

An existing thread may be scheduled to run on some operating system core  $c$ . An already running thread may be unscheduled and be left in a dormant state to be scheduled again:

$thread\_schedule \triangleq$ any $t, c$ where $t \in threads \setminus \text{dom}(affinity)$ $status(c) = \text{ON}$ then $affinity(t) := c$ end	$thread\_unschedule \triangleq$ any $t, c$ where $t \in \text{dom}(affinity)$ $affinity(t) = c$ then $affinity := \{t\} \triangleleft affinity$ end
---	---

d) *Application*: Several threads are grouped into an application. The primary role of an application is to define workload type shared by a number of threads. Applications may be created and destroyed during the system lifetime. The set of current applications is defined by set

$$apps \subseteq \text{APPS}.$$

Each thread belongs to an application and each application owns at least one thread. This is captured by the following (surjective and total) relation

$$app\_threads \in apps \leftrightarrow threads.$$

For each thread there is just one owning application:

$$app\_threads^{-1} \in threads \rightarrow apps.$$

When an application is created, it appears together with one thread of its own (this even *refines* event  $thread\_start$ ):

$app\_start \triangleq$   
any  $a, t$  where  
 $a \notin apps \wedge t \notin threads$   
then  
 $apps := apps \cup \{a\}$   
 $threads := threads \cup \{t\}$   
 $app\_threads := app\_threads \cup \{a \mapsto t\}$   
end

Destroying an application cancels all the running application threads and removes them from the system:

$app\_stop \triangleq$   
any  $t, a$  where  
 $t \in threads \wedge a \mapsto t \in app\_threads \wedge app\_threads[\{a\}] = \{t\}$   
then  
 $threads := threads \setminus \{t\} \parallel affinity := \{t\} \triangleleft affinity$   
 $apps := apps \setminus \{a\} \parallel app\_threads := app\_threads \setminus \{a \mapsto t\}$   
end

e) *Workload*: The purpose of an application is to provide a computation service. This it accomplishes by assigning incoming *workload* to application threads. The unit of a workload is a *job*. Every job is specific to an application so that only threads of a certain application may process a given job.

The pending and executing jobs of a systems are defined by variable  $jobs$ :

$$jobs \in \text{JOBS} \leftrightarrow apps$$

where **JOBS** is the universe of all jobs. To run a job, it must be allocated to a thread. At any given time a thread executes at most one job. The following partial injection captures this relationship:

$$job\_alloc \in \text{dom}(job\_alloc) \mapsto threads$$

A job allocation must agree with the ownership of a thread to which the job is assigned:

$$jobs^{-1}; job\_alloc \subseteq app\_threads$$

Here  $jobs^{-1}$  is the converse of function  $jobs$  and  $f;h$  denotes forward functional composition.

To reason about computational complexity of a job, we define the number of *steps* comprising a given job. A step is a normalised complexity measure independent of core properties and shared by all the jobs:

$$job\_steps \in \text{dom}(jobs) \rightarrow \mathbb{N}$$

A step execution time varies from core to core and with core frequency. A new job may appear at any moment; it must be assigned to an existing application.

$job\_create \triangleq$   
any  $j, a, w$  where  
 $j \notin \text{dom}(jobs) \wedge a \in apps \wedge w > 0$   
then  
 $jobs := jobs \cup \{j \mapsto a\} \parallel job\_steps(j) := w$   
end

In the above,  $w$  defines the job complexity in the terms of steps. An existing but yet unassigned job may be allocated to a thread of the job application. The thread in question must be mapped to a core but not already executing any other job:

$job\_allocate \triangleq$   
any  $j, t$ , where  
 $j \in \text{dom}(jobs) \setminus \text{dom}(job\_alloc)$   
 $t \in (app\_threads[\{jobs(j)\}] \setminus \text{ran}(job\_alloc)) \cap \text{dom}(affinity)$   
then  
 $job\_alloc := job\_alloc \cup \{j \mapsto t\}$   
end

When a job is assigned to a scheduled thread, the job “runs” going through a predefined number of timed steps. At some point, a job finishes and vanishes from the system.

f) *Timing*: A number of already defined phenomena require some form of timing. There is no native support for clocks and timers in Event-B but there are a number of established approaches to time modelling. The one we use is closely related to timed automata with integer clocks. To keep track of time progress and time various activities of a system we introduce a number discrete *timers*. A timer functions like a stopwatch - it counts down to zero from the initial integer value. All the system timers count in synchrony (that is, driven by one global clock) and there is no limit on the number of timers in the system. Once any timer reaches zero, time freezes to allow for a sub-system interested in this timer to react to a deadline; to enable further progress, the sub-system must also reset the timer or delete it. The set of all clocks is defined by function  $\tau$ :

$$\tau \in \text{TA} \rightarrow \text{TIME}, \quad \text{TIME} = \mathbb{N} \cup \{\text{DISABLED}\}$$

$\tau(x)$  gives the current value of clock  $x$  which is  $> 0$  before deadline, 0 on the deadline and **DISABLED** if the clock is not used. To avoid complicated progress arguments, we assume there is a plentiful supply of clocks. At this level of abstraction it suffices to require that set **TA** is infinite.

All the clocks are synchronously updated by an event *time*:

```

time  $\triangleq$ 
any p1, p2 where
  0  $\notin$  ran( $\tau$ )
  p1 =  $\tau^{-1}$ [[DISABLED, 0]]
  p2 = dom( $\tau$ ) \ p1
   $\forall c \cdot c \in status$  [[ON]]  $\Rightarrow temp(c) \leq CORE\_TEMP\_CRIT(c)$ 
then
   $\tau := (p1 \triangleleft \tau) \cup \{x \cdot x \in p2 | x \mapsto \tau(x) - 1\}$ 
end

```

Notice the disabling condition  $0 \in \text{ran}(\tau)$  which stops the timer until the deadline of a clock is processed. The last guard prevents clock progress when there is an overheated core: *it makes reaction to core overheating immediate*.

g) *Job deadlines*: One application of timing is the definition of the job deadlines and, to make the concept meaningful, the notion of job execution time. The latter is linked to the notion of jobs steps defined above.

A user deadline is set at the point of job allocation as opposed to the point of job creation. We do not yet model job queues (it is rather a set in this model) so there is no fairness property for job allocation. An extended version of the *job\_allocate* event sets a user deadline for a job.

```

job_allocate  $\triangleq$ 
any j, t, ta, ta_user, udl where
  ...
   $\tau(ta) = DISABLED \wedge \tau(ta\_user) = DISABLED$ 
   $ta \neq ta\_user \wedge udl \in \mathbb{N}$ 
then
  ...
  job_step_time(j) := ta || user_deadline(j) := ta_user
   $\tau := \tau \triangleleft \{$ 
     $ta \mapsto F\_STEP\_TIME(affinity(t) \mapsto freq(affinity(t))),$ 
     $ta\_user \mapsto udl \}$ 
end

```

The user deadline clock is stored in *user\_deadline(j)*; another clock times the execution of job steps and is stored in *job\_step\_time(j)*. The assignment to  $\tau$  initialises these two clocks. Note that the step running time is defined by the kind and the frequency of a core on which a thread executing the job is scheduled. User deadline clock *user\_deadline(j)* runs down without pauses from the point of job allocation irrespectively of whether a thread processing the job is running or not.

Job steps are timed according to the respective core performance and are affected by a change of a core frequency. Should a job fail to meet user deadline, it is cancelled even before all the job steps are done. Finally, if a job completes before a user deadline expires, it is accepted as a successful job execution (event *job\_finish*).

h) *Scheduling*: As defined above, a core runs several threads at the same time. In our model, we do not go into the minute details of scheduling within a group of threads assigned to a core but rather assume that core indeed runs all the mapped threads in parallel (that is, the *time band* in which the core model is given does not allow us to distinguish between instances of individual thread executions and the whole picture is blurred to give an illusion of a multi-threaded core). There are limits to a number and kind of threads that may be run on single core before the execution times of individual threads are affected. Intuitively, if a thread is computationally intensive and never has to wait for data, it cannot share a core with another thread without sacrificing performance. In practice,

many applications require data retrieval or do blocking system calls which make a thread idle and hence free a core to run another thread. The purpose of the scheduling refinement step is to bundle threads from various applications into thread groups that may be assigned to the same core without hindering thread performance.

Variable *thread\_load* characterises a thread in terms of lower bound of operations per second (or a comparable normalised measure) necessary to run the thread at full speed.

$$thread\_load \in threads \mapsto LOAD$$

Thread load must be defined for all the running threads,  $\text{dom}(affinity) \subseteq \text{dom}(thread\_load)$ . The crucial property is that the overall load of threads assigned to a core does not exceed the core computational capability at the current core frequency:

$$\forall c \cdot c \in status^{-1}[\{ON\}] \Rightarrow \text{sum}(affinity^{-1}[\{c\}] \triangleleft thread\_load) \leq CORE\_LOAD\_MAX(c \mapsto freq(c))$$

Thread load affects the way threads are mapped to cores:

```

thread_schedule  $\triangleq$ 
any t, c where
  ...
   $\text{sum}(affinity^{-1}[\{c\}] \triangleleft thread\_load) +$ 
   $thread\_load(t) \leq CORE\_LOAD\_MAX(c \mapsto freq(c))$ 
then
  affinity(t) := c
end

```

For a short while, the power budget may be lower than the power already drawn. To emphasize the immediacy of a reaction, the system timer is stopped until the power budget constraint is resolved either by shutting cores or adjusting their frequency and voltage.

$$time \triangleq \text{any } \dots \text{ where } \dots \wedge \text{sum}(c\_pwr) \leq pwr\_b \text{ then } \dots \text{ end}$$

#### IV. DSL-KIT

Ability to do design prototyping with the developed formal domain model was perceived to be one of the more interesting directions. To this end, we have developed an extension to the Event-B Rodin [6] modelling framework that adds a capability of rapidly defining a custom DSL on the basis of an Event-B specification. The extension, called DSL-Kit, works by combining a general-purpose imperative specification language with a custom set of 'commands' defined by an Event-B specification. The imperative layer provides control flow construct that allow one to write scenarios or imperative programs in the terms of Event-B events. From the viewpoint of Event-B, the imperative layer is an extra refinement step strengthening event guards and declaring new hidden state.

The core of DSL-Kit is a formal specification language based on the following principal structuring units:

- system - the top-level unit defined as a parallel composition of several *actors*;
- module - a self-contained unit providing definitions of *actors* and *actions*;
- actor - a unit of concurrency; its body is sequential code guarded by rely/guarantee conditions;
- action - a function-like entity;

```

actor dvfs(c : CORES)
rely ...
{
  if (status(c)! = OFF){
    core_threads : set(THREADS) = affinity-1[[c]];
    ml : int = CORE_LOAD_MAX(c, freq(c));
    if (core_threads! = {}){
      total_load : int = sum(thread_load[core_threads]);
      if (ml < total_load + STEP)
        if (enabled core_dvfs_running(...))
          core_dvfs_running(c, vdd(c) + CH, freq(c) + CH);
        else if (ml > total_load + STEP and ml > MIN_LOAD)
          if (enabled core_dvfs_running(...))
            core_dvfs_running(c, vdd(c) - CH, freq(c) - CH);
      } else
        #nothing is running, run down frequency gradually
        if (enabled core_dvfs(...) - 1) and ml > MIN_LOAD)
          core_dvfs(c, vdd(c) - 1, freq(c) - 1);
    }
  }
}
guar ...

```

Fig. 2. Cores-DSL script for a simple on-demand frequency/voltage governor.

- control flow statements - sequential composition, *if*, *while*, *for* and auxiliary variable declarations;
- expressions and predicates, expressed in the Event-B mathematical language; this makes possible to relate DSL-Kit to Event-B without logic mapping.

Not all scenarios defined on top of the Event-B domain model correspond to behaviour permitted by the domain model. For instance, the following defines a scenario made of a sequence of two events:

$$on(c, v, f); off(c);$$

The first event switches a core on and the second switched off the same core (assuming all identifiers are locally bound). If, instead, we write

$$on(c, v, f); on(c, v, f);$$

it appears that the second event cannot be enabled and the scenario must halt prematurely. In fact, even in the first example, we could not know that  $on(c, v, f)$  starts in state where the guard of even  $on$  instantiated with arguments  $c, v, f$  is enabled. To collect all such conditions automatically, we need to know the kind of states in which an event (or, for generality, any imperative statement) is enabled and also the kind of states it produces upon termination.

The module containing the Cores domain model is made available for new actor and system definitions. Module variables (translation of Event-B variables) are accessible read-only. Their state may only be modified via module actions. Event-B invariants thus become module axioms (Event-B has already established that events preserve invariants). This makes high-level design protected from small changes in the domain model specification, re-generated mechanically from an Event-B model,

DSL-Kit uses a variant of Floyd-Hoare logic to compute the strongest post-condition given some current state and a statement. On the basis of strongest post-conditions, the verification conditions generation procedure constructs a number of theorems. Among them are conditions establishing that every scenario runs until completion. To facilitate reasoning about design prototypes, DSL-Kit offers invariant, auxiliary variables, iteration and recursion variants, and rely/guarantee conditions for reasoning about concurrency and shared state.

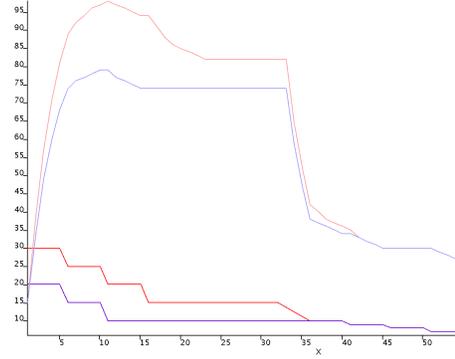


Fig. 3. Time graph of core frequencies (dark red/blue, lower) and temperatures (light red/blue, higher); the graph shows the performance of an on-demand DVFS governor (see script in Fig. 2) in a scenario where two jobs are scheduled simultaneously for both cores and cancelled after running for 33 seconds. The thread mapper script ramps up initial frequencies and voltages and these are gradually lowered by the DVFS script. Idling produces rapid cooling starting at 33 second mark.

The script in Fig. 2 shows a simple yet functional on-demand frequency governor. The performance of the script is shown in Fig. 3. The plotted graph is discrete and it depicts time series of domain model variables  $freq$  and  $core\_temp$ .

There is an important relation between DSL specifications and the original Event-B domain model: *for every DSL specification there exists an Event-B model refining the Event-B specification of the domain model*. A formal proof consists in exhibiting a translation procedure converting a DSL specification to Event-B model. Such a translation is trivial in principle but there does not seem to be a practical reason to implement it.

## V. CONCLUSIONS

We have presented a formal DSL for multi-core systems. It is a completely proven Event-B specification with 9 refinement steps and over 600 hundred proof obligations. Much of the final detalisation and 'debugging' (that is, fixing omissions and progress problems for safety and correctness are proven statically) was accomplished by building design prototypes in the developed DSL. We have done 15 revision cycles over the course of five months since the initial Event-B domain model was developed.

## ACKNOWLEDGEMENT

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