

# Verification of distributed dataspace architectures <sup>\*</sup>

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**Abstract.** The *space calculus* is introduced as a language to model distributed dataspace systems, i.e. distributed applications that use a shared (but possibly distributed) dataspace to coordinate. The publish-subscribe and the global dataspace are particular instances of our model. We give the syntax and operational semantics of this language and provide tool support for functional and performance analysis of its expressions. Functional behaviour can be checked by an automatic translation to  $\mu$ CRL and the use of a model checker. Performance analysis can be done using an automatically generated distributed C prototype.

## 1 Introduction

A distributed system is generally seen as a number of single-threaded applications together with a distributed communication layer that coordinates them. Various shared dataspace models have been proposed to solve the task of coordination for parallel computing applications (Linda [8], Gamma [1]), network applications (Bonita [22], WCL [23]), command and control systems (Splice [3]), management of federations of devices (JavaSpaces [15]).

A shared dataspace (or tuple space) architecture is a distributed storage of information and/or resources, viewed as an abstract global store, where applications read/write/delete pieces of data. In this paper, we focus on the problem of designing, verifying and prototyping distributed shared dataspace systems. Building correct distributed systems is a difficult task. Typical required properties include transparent data distribution and fault-tolerance (by application replication and data replication), which are usually ensured at the price of giving up some performance. Many questions occur when deciding on the exact shape of the distributed dataspace. For instance: what data should be replicated (in order to prevent single points of failure or increase efficiency)? should the local storages be kept synchronized or should they be allowed to have different views on the global space? should the migration of data between local storages be on a subscription basis or rather “on demand”?

**The space calculus**, introduced in this paper, is an experimental framework in which verification and simulation techniques can be applied to the design of distributed systems that use a shared dataspace to coordinate their components.

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We provide a tool that translates a space calculus specification into a  $\mu$ CRL specification [16]. From this code a state space graph can be generated and analyzed by means of the model checker CADP [13]. A second tool generates distributed C code to simulate the system. System designers may use the automatic verification and simulation possibilities provided by the  $\mu$ CRL toolset [2] to verify properties of their architecture. Complementary, the distributed C prototype can be used for testing purposes, and to get an indication about the performance (e.g. number of messages, used bandwidth, bottlenecks). Several design choices can be rapidly investigated in this way. Ultimately, the prototype C implementation could even be used as a starting point for building a production version.

The operational semantics of our space calculus provides the formal ground for algebraic reasoning on architectures. Despite its apparent simplicity, our calculus is highly expressive, capable of modeling various destructive/non destructive, global/local primitives. By restricting the allowed space connectives and the allowed coordination primitives, we obtain well known instances, such as the kernels of Splice (a publish-subscribe architecture), and JavaSpaces (global space). Some specific features, like the transactions in JavaSpaces or dynamic publish/subscribe in Splice are out of our scope. Our goal is a uniform framework where core characteristics of various dataspace architectures should be present, in order to allow for studies and comparisons. The verification tool will help getting fast insights in the replication and distribution behaviour of certain architectures, for instance. The simulation tool can help identifying the classes of applications appropriate to each architecture.

**Related work.** An overview of shared dataspace coordination models is given in [24]. Some work that studies different semantics has been done in [4, 5, 7, 6], on which we based the style of our operational semantics. [7] compares the publish/subscribe with the shared dataspace architectural style by giving a formal operational semantics to each of them. We also aim at being able to compare the two paradigms, but we take a more unifying perspective: we consider both as being particular instances of the more general distributed dataspace model and express them in the same framework. [10] was the first attempt to use a Unity-like logic to reason on a shared dataspace coordination model (Swarm). [19] has goals similar to ours. It provides a framework for describing software architectures in the theorem prover PVS. However, it seems that the verification of functional behaviour is out of the scope of that paper. In [9], a language for specification and reasoning (with TLA) about software architectures based on hierarchical multiple spaces is presented. The focus there is on the design of the coordination infrastructure, rather than on the behaviour of systems using it. In [17], a translator from the design language VPL to distributed C++ code is presented. VPL specifications can be verified using the CWB-NC toolset. Compared to that approach, our work is more specific. We concentrate on shared dataspace architectures and define a “library” of carefully chosen set of primitives that are both handy and expressive. In [12], scenario-based verification is introduced as a useful technique in between verification and testing. Our language also supports that.

In section 2, we present the syntax and semantics of the space calculus and we comment its main characteristics. Then (section 3) we introduce the supporting tools. Section 4 contains two examples. We end with some concluding remarks (section 5).

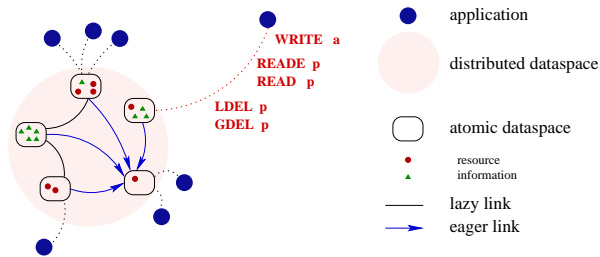


Fig. 1. A distributed dataspace architecture

## 2 The Space Calculus

### 2.1 Informal view

We model the shared space as a graph with atomic spaces as nodes (see Figure 1). We consider two types of links between spaces: *eager* and *lazy*. When elements are written in a local space, they are asynchronously transferred over all eager links that start in this local space. Eager links can be used to model subscription and notification mechanisms. Lazy links, on the other hand, are demand driven. Only when a data item is requested in some atomic space, it is transferred via a lazy link from one of the neighbouring spaces. Besides modeling the shared space, the space calculus provides a set of coordination primitives for *applications*: write, blocking and non-blocking read, local and global delete operations. Applications are loosely coupled in the sense that they cannot directly address other applications.

With so many existing shared dataspace models, it is difficult to decide what features are the most representative. Some choices that we are faced to are: atomic spaces can be sets or multisets; when transferring data items between different spaces, they could be replicated or moved; the primitives can be location-aware or location-independent; the retrieve operation can be destructive or non-destructive, etc. The answers depend of course on the specific application or on the purpose of the architecture. In order to allow the modeling of as many situations as possible, we let the user make the distinction between data items that should be treated as *information* (e.g. data from a sensor), for which multiplicity is not relevant, and data items that should be treated as *resource* (e.g. numbers to be added, jobs to be executed), for which multiplicity is essential. When handling elements, the space takes into account their type. The transfer between spaces means “copy” for information items and “move” for resources. Similarly, the lookups requested by applications are destructive for resources and non-destructive for information items.

The atomic spaces are multisets in which the elements tagged as information are allowed to randomly increase their multiplicity. As for the question whether to give to applications the possibility to directly address atomic spaces by using handles, like for instance in [22], we have chosen not to, in order to keep the application layer and the coordination layer as separated as possible. The advantage of a clear separation is that the exact distribution of the space is transparent to the applications.

### 2.2 Syntax and semantics

As mentioned before, a system description consists in our view from a number of program applications and a number of connected atomic spaces. We refer to the topol-

ogy of the distributed space by giving atomic spaces (abstract) locations, denoted by  $i, j, \dots$ . The data items come from a set  $\mathcal{D}$  of values, ranged over by  $a, b, \dots$ . Furthermore, we assume a set of patterns  $\mathcal{Pat}(\mathcal{D})$ , i.e. properties that describe subsets of  $\mathcal{D}$ . ( $\mathcal{D} \subseteq \mathcal{Pat}(\mathcal{D})$ ).  $p, q, \dots$  denote patterns. We also postulate two predicates on patterns:  $\text{match} : \mathcal{Pat}(\mathcal{D}) \times \mathcal{D} \rightarrow \{\top, \perp\}$  to test if a given pattern matches a given value, and  $\text{inf} : \mathcal{Pat}(\mathcal{D}) \rightarrow \{\top, \perp\}$  to specify whether a given pattern should be treated as information or as resource. The predicate  $\text{res} : \mathcal{Pat}(\mathcal{D})$  will be used as the complementary of  $\text{inf}$ .

A process  $([P]^i)$  is a program expression  $P$  residing at a location  $i$ . A program expression is a sequence of coordination primitives: write, read, “read if exists”, local delete, global delete. These primitives are explained later in this section. Formal parameters in programs are denoted by  $x, y, \dots$ , the empty program is denoted by  $\varepsilon$ , and  $\perp$  denotes a special error value.

$$\begin{aligned} \text{Prim} &::= \text{write}(a) \mid \text{read}(p, x) \mid \text{read}\exists(p, x) \mid \text{l}del(p) \mid \text{g}del(p) \\ P &::= \varepsilon \mid \text{Prim}.P \\ \text{Proc} &::= [P]^i \end{aligned}$$

The lazy and eager behaviours of the connections are specified as special marks:  $\downarrow_p^i$  (meaning that atomic space  $i$  published data matching  $p$ ),  $\uparrow_p^i$  ( $i$  subscribes to data matching  $p$ ),  $\uparrow_i^j$  ( $i$  and  $j$  can reach each other’s elements). If  $\downarrow_p^i$  and  $\uparrow_q^j$  are present, then all data matching  $p \wedge q$  written in the space  $i$  by an application will be asynchronously forwarded to the space  $j$ . We say then that there is an eager link from  $i$  to  $j$ . The presence of  $\uparrow_i^j$  indicates that there is a (symmetric) lazy link from space  $i$  to  $j$ . That is, all data items of  $i$  are visible for retrieve operations addressed to  $j$  by an application.

For administrative reasons, the set of data items ( $a$ ) is extended with buffered items that have to be sent ( $!a^j$ ), pending request patterns ( $?p$ ) and subscription policies ( $\circ p, k, \circ p, k, t$ ). Subscription policies are inspired by Splice and their function is to filter the data coming into a space as consequence of a subscription. Based on *keys* and *timestamps*, some of the data in the space will be replaced (overwritten) by the newly arrived element. The parameters  $k, t$  are functions on data  $k : \mathcal{D} \rightarrow \text{Keys}$ ,  $t : \mathcal{D} \rightarrow \mathbb{N}$  that dictate how the keys and the timestamps should be extracted from data items. If the newly arrived element  $a$ , matching  $p$ , meets the filter  $\circ p, k$ , then  $a$  will overwrite all data matching  $p$  with the key equal to that of  $a$ . If it meets the filter  $\circ p, k, t$ , it will overwrite all data matching  $p$  with the key equal to that of  $a$  and timestamp not fresher than that of  $a$ . With this second filter, it is also possible that  $a$  drops out, if its timestamp is too big. A configuration ( $\mathcal{C}$ ) then consists of a number of atomic dataspaces and applications, each bound to a location, and a number of links. The parallel composition operator  $\parallel$  is associative and commutative.

$$\begin{aligned} \text{data} &::= a \mid !a^j \mid ?p \mid \circ p, k \mid \circ p, k, t \\ \text{Data} &::= \langle D \rangle^i, \text{ where } D \text{ is a finite set over data} \\ \text{Link} &::= \uparrow_i^j \mid \uparrow_p^i \mid \downarrow_p^i \\ \text{Conf} &::= \text{Data} \mid \text{Proc} \mid \text{Link} \mid \text{Conf} \parallel \text{Conf} \end{aligned}$$

The operational semantics of the space calculus is defined as a labeled transition relation  $\mathcal{C} \xrightarrow{a} \mathcal{C}'$ , meaning that if the system is in configuration  $\mathcal{C}$  then it can do an  $a$ -step to



$$\begin{array}{l}
D \uplus_p a = D - [b \in D \mid \text{match}(p, b)] + [a] \\
D \uplus_{p,k} a = D - [b \in D \mid \text{match}(p, b) \wedge k(b) = k(a)] + [a] \\
D \uplus_{p,k,t} a = D - [b \in D \mid \text{match}(p, b) \wedge k(b) = k(a) \wedge t(b) \leq t(a)] + [a] \\
D \uplus_d d = D - [d \in D] + [d] \\
\\
D \oplus d = \begin{cases} D \uplus_d d & \text{if } \text{inf}(d) \\ D + [d] & \text{if } \text{res}(d) \end{cases} & D \ominus a = \begin{cases} D & \text{if } \text{inf}(a) \\ D - [a] & \text{if } \text{res}(a) \end{cases} \\
\\
D \oplus [d_1 \dots d_n] = D \oplus d_1 \oplus \dots \oplus d_n & B \boxplus d = \begin{cases} B \uplus_d d & \text{if } \text{inf}(d) \\ [d] & \text{if } \text{res}(d) \end{cases} \\
\\
D \uplus a = \begin{cases} D \uplus_{p,k} a & \text{if } \bigcirc p, k \in D \wedge \text{match}(p, a) \\ D \uplus_{p,k,t} a & \text{if } \bigcirc p, k, t \in D \wedge \text{match}(p, a) \\ D \oplus a & \text{if } \nexists \bigcirc p, k, \bigcirc p, k, t \in D \text{ s.t. } \text{match}(p, a) \end{cases}
\end{array}$$

**Fig. 3.** Auxiliary operators on multisets.

the pattern occurring in  $d$ . That is,  $\text{inf}(!a^j) = \text{inf}(a)$  and  $\text{inf}(?p) = \text{inf}(\bigcirc p, k) = \text{inf}(\bigcirc p, k, t) = \text{inf}(p)$ . The same holds for  $\text{res}$ .

We now explain the intuitive semantics of the coordination primitives.

*write(a)*: write data item  $a$  into the local dataspace, to be automatically forwarded to all subscribed spaces.  $a$  is added to the local dataspace (W1) and an auxiliary  $w(i, a)$  step is introduced. When pushing  $w(i, a)$  to the top level, if  $a$  matches a pattern published by  $i$ , then  $!a^j$  items are introduced for all subscriptions  $\uparrow_p^j$  matching  $a$  (rules W2, W3). At top level, the auxiliary  $w(i, a)$  step gets promoted to a *write(a)* step (W4). Finally, the  $a$  items are sent to the subscribed spaces asynchronously (W5). The operator  $\uplus$  in the right hand side of rule (W5) states that the freshly received data item should be added to the local database taking into account the local subscription policies.

*read(p, x)*: blocking test for presence, in the local space and its lazy linked neighbouring spaces, of some item  $a$  matching  $p$   $x$  will be bound to  $a$ . This results in generating a  $?p$  request, keeping the application blocked (R $\tau$ ). If a matching  $a$  has been found, it is returned and the application is unblocked (R). Meanwhile, the lazy linked neighbours of the local space asynchronously respond to the request  $?p$ , if they have an item matching  $p$  (TAU). *read* $\exists(p, x)$ : non-blocking test for presence in the local space. If some item  $a$  matching  $p$  exists in the local space, it is bound to  $x$ ; otherwise a special error value  $\perp$  is returned. Delivers a matching  $a$  from the local space, if it exists (R $\exists$ 1). Otherwise an error value is returned (R $\exists$ 2). *ldel(p)*: atomically removes all elements matching  $p$  from the local space (LD). *gdel(p)*: this is the global remove primitive. It atomically deletes all items matching  $p$ , in all atomic spaces. Note that due to its global synchronous nature, *gdel* can not be lifted over atomic spaces (GD2). Finally,

the general parallel rule (act) defines parallelism by interleaving, except for *write* and *gdel* which have their own parallel rules to ensure synchronization.

### 2.3 Modeling some dataspace paradigms

The kernels of some well-known dataspace paradigms can be obtained by restricting the allowed configurations and primitives.

Splice [3] implements a publish-subscribe paradigm. It has a loose semantics, reflecting the unstable nature of a distributed network. Applications announce themselves as publishers or subscribers of data sorts. Publishers may write data items to their local agents, which are automatically forwarded to the interested subscribers. Typically, the Splice primitives are optimized for real-time performance, and don't guarantee global consistency. The space calculus fragment without lazy links and restricted to the coordination primitives *write*, *read*, *ldel* corresponds to the reliable kernel of Splice. Network searches (modeled by the lazy links) and global deletion (*gdel*) are typically absent. In Splice, data sorts have keys, and data elements with the same key may overwrite each other – namely at the subscriber's location, the “fresh” data overwrites the “old” one. The order is given by implicit timestamps that elements get in the moment when they are published. The overwriting is expressible in our calculus, by using the eager links with subscribe policies. The Splice's timestamps mechanism is not present, but some timestamping behaviour can be mimicked by explicitly writing and modifying an extra field in the tuples that models the data.

JavaSpaces [15] on the contrary can be viewed as a *global* dataspace. It typically has a centralized implementation, and provides a strongly consistent view to the applications, that can write, read, and take elements from the shared dataspace. The space calculus fragment restricted to a single atomic space to which all coordination primitives are attached, and with the primitives *write*, *read*, *read* $\exists$  forms a fragment of JavaSpaces. Transactions and leasing are not dealt with in our model. Note that with the mechanism of marking the data “information” or “resource”, we get the behaviour of both destructive and non-destructive JavaSpaces lookup primitives: our *read*, *read* $\exists$  works, when used for information, like *read* and *readIfExists* from JavaSpaces, and like *take* and *takeIfExists* when called for resources.

So, interesting parts of different shared dataspace models are expressible in this framework.

## 3 The verification and prototyping tools

We defined a mapping from every configuration in the operational semantics to a process in the  $\mu$ CRL specification language [16]. An incomplete description of this mapping is given later in this section. The generation of the  $\mu$ CRL specification following this mapping is automated. Therefore, the  $\mu$ CRL toolset [2] can be immediately used to simulate the behaviour of a configuration. This toolset is connected to the CADP [13] model checker, so that temporal properties on systems in the space calculus can be automatically verified by model checking. Typical verified properties are deadlock freeness, soundness, weak completeness, equivalence of different specifications.

The state of a  $\mu$ CRL system is the parallel composition of a number of processes. A process is built from atomic actions by sequential composition ( $\cdot$ ), choice ( $+$ , sum),

```

Atomic (id:Nat, D:TupleSet, Req: TupleSet,
        ToSend: NatTupleSet, todel:Tuple,
        LL: NatSet, PL: TupleSet, SL: SubscriptionList) =

% W
sum(v:Tuple,
    W(v). sum(NewToSend: NatTupleSet,
              sum(NewD: TupleSet,
                  getToSend(v, ToSend, NewToSend, D, NewD).
                  Atomic(id, NewD, Req, NewToSend,
                        todel, LL, PL, SL)))
    <| and(isData(v), match(v, PL)) |> delta )
+ sum(v:Tuple,
    W(v).
    Atomic(id, a(v,D), Req, ToSend, todel, LL, PL, SL)
    <| and(isData(v), not(match(v,PL))) |> delta )

% async send
+ sum(x:Nat, sum(y:Tuple,
    el_send(x,y).
    Atomic(id, D, Req, r(x,y,ToSend), todel, LL, PL, SL)
    <| in(x,y,ToSend) |> delta ))

% async receive
+ sum(x:Tuple,
    el_rcv(id,x).
    Atomic(id, add(x,D,SL), Req, ToSend, todel, LL, PL, SL))
...

```

**Fig. 4.** Fragment from a  $\mu$ CRL specification of an atomic space

conditionals ( $\cdot \triangleleft \cdot \triangleright \cdot$ ) and recursive definitions. For our purpose, we introduce processes for each atomic space and for each application. An additional process, called the *TokenManager*, has to ensure that operations requiring global synchronization (*gdel*) don't block each other, thus don't introduce deadlocks. Before initiating a global delete operation, a space has to first request and get the token from the manager. When it has finished, it has to return the token to the manager, therefore allowing other spaces to execute their *gdels*. A second additional process, *SubscriptionsManager*, manages the list (multiset) of current subscriptions. When an item  $a$  is written to an atomic space, that space synchronizes with the *SubscriptionsManager* in order to get the list of the other atomic spaces where the new item should be replicated or moved.

For simplicity, we model the data items as tuples of natural numbers – fields are modeled by the  $\mu$ CRL datasort `Nat`, tuples by `Tuple`.

An atomic space has two interfaces: one to the application processes, and one to the other atomic spaces. In  $\mu$ CRL calls between processes are modeled as synchronization between atomic actions. The primitives of the space calculus correspond to the following atomic actions of `Atomic`:  $\{W, R, RE, Ldel, Togdel, Gdel\}$ . The interface to the

other atomic processes is used to send/receive data items and patterns for read requests. In Figure 4, the  $\mu$ CRL specification of a space's *write* behaviour is shown.

The application programs are also mapped to  $\mu$ CRL processes. Execution of coordination primitives is modeled as atomic actions, that synchronize with the corresponding local space's pair actions. This synchronization with the space is described by a communication function.

Another tool translates space calculus specifications to a distributed implementation in C that uses MPI (Message-Passing Interface) [14] for process communication. Different machines can be specified for different locations, thus getting a real distribution of spaces and applications. By instrumenting this code, relevant performance measures for a particular system under design can be computed. The result of the translation is more than a software simulation. It is actually a prototype, that can be tested in real-time conditions, in a distributed environment.

### 3.1 The space calculus tool language

In order to make the space calculus usable as specification language, the tools supporting it work with a concrete syntax. The data universe considered is *tuples of naturals* and the patterns are incomplete tuples (e.g.  $\langle 1, *, 2 \rangle, \langle * \rangle$ ). Apart from the syntactical constructions already defined, we allow external actions (e.g. EXTping), *assignment* of data variables, assignment of tuple variables and *if* and *while* with standard semantics. We give now a brief description of this language, including a precise syntax written in a slightly simplified YACC format.

Since we allow exactly one space per location, it is nice to give *names* to spaces and to say, instead of saying that a program stays at location  $i$ , that the program runs at the space  $\langle \text{name} \rangle$ . A specification of a configuration consists of:

- (optional) fixing the tuple size (**nfields**) and the first natural value strictly greater than any field of any tuple (**upbound**). The default values are nfields=1, upbound=2.
- (optional) define the inf/res predicates, by mentioning the patterns for which res should be  $\top$ . Any pattern  $p$  not included by the **res** declaration has  $\text{inf}(p) = \top$ .
- describing the spaces, by giving each space a name and, optionally, the machine where it's supposed to live. The default machine is "localhost".
- describing the applications, by specifying for each application its name, the name of the space with which it shares the location (the physical location as well) and its program.

Apart from the primitives **read**, **readE**, **write**, **ldel**, **gdel**, the actual language includes some extra constructions to provide easy data manipulation and control possibilities: natural variable names and expressions, projection of a tuple on a field, assignments, if, while, external actions that can be specified as strings.

The condition of *if* and *while* is very simple: a standard boolean value or a variable name that gets tested for correctness. Namely, "*if x*" means "if  $x$  is not *error*". Extending the conditions is further work.

The key and timestamp functions needed in the subscription policies are considered to be projections on the fields of the tuples – one field for the timestamp, possibly more for the key. Therefore, key functions are represented as lists of field indices and timestamp functions as one field index.

EXTCOMMAND	means $[E][X][T][a - zA - Z]^+$
INTID	means $[i][a - zA - Z0 - 9]^*$
ID	means $[a - zA - Z][a - zA - Z0 - 9]^*$ (that is not INTID)
INT	means $[0 - 9]^+$
configuration	: settings declarations
settings	:   setting settings
setting	: <b>nfields</b> = INT   <b>upbound</b> = INT   <b>res</b> pattern
declarations	: space declarations   link declarations   application declarations
space	: <b>space</b> ID ( ID )   <b>space</b> ID
link	: <b>LL</b> ( ID , ID )   ID - > pattern   ID < - pattern   ID < - pattern intlist   ID < - pattern intlist INT
pattern	: < tuple >
tuple	: datum   tuple , datum
datum	: *   INT   INTID
intlist	: INT   intlist , INT
intexpression	: INT   INTID   projection   intexpression + intexpression
projection	: pattern / INTID   ID / INTID
application	: <b>app</b> ID @ ID { program }
program	:   command ; program
command	: <b>write</b> pattern   <b>write</b> ID   <b>read</b> pattern ID   <b>readE</b> pattern ID   ID := pattern   INTID := intexpression   <b>ldel</b> pattern   <b>gdel</b> pattern   <b>publish</b> pattern   <b>subscribe</b> pattern   <b>subscribe</b> pattern intlist   <b>subscribe</b> pattern intlist INT   <b>if</b> condition { program }   <b>while</b> condition { program }   EXTCOMMAND
condition	: ID   not(ID)   true   false

Fig. 5. The YACC style syntax definition

## 4 Examples

We use the new language to specify two small existing applications, studied in [21] and [18], respectively. The goal of these examples is to show that our language is very simple to use and to illustrate the typical kind of problems that space calculus is meant for: transparent distribution of data and transparent replication of applications.

### 4.1 Toward distributed JavaSpaces

One of the initial motivations of our work was to model a distributed implementation of JavaSpaces, still providing the same strongly consistent view to the applications. When restricting the primitives as discussed in section 2.3, the expression  $\langle \emptyset \rangle^0$  represents the kernel of JavaSpaces and the expression  $\langle \emptyset \rangle^0 \parallel \langle \emptyset \rangle^1 \parallel \downarrow_\star^0 \parallel \uparrow_\star^0 \parallel \downarrow_\star^1 \parallel \uparrow_\star^1$  models a distributed implementation of it, consisting of two spaces eagerly linked by subscriptions matching any item.

Two rounds of the Ping-Pong game ([15], [21]) can be written in the space calculus as follows:

$$\begin{aligned} Ping &= write(1).read(0, x).write(1).read(0, x) \\ Pong &= read(1, x).write(0).read(1, x).write(0) \end{aligned}$$

(with  $\mathcal{D} = \{0, 1\}$  and  $\text{inf}(x) = \perp, \forall x$ ). We wish that the distribution of the space should be completely transparent to the applications, i.e. that they run on one space exactly the same that they run on two:

$$[Ping]^0 \parallel [Pong]^0 \parallel \langle \emptyset \rangle^0 = [Ping]^0 \parallel [Pong]^1 \parallel \langle \emptyset \rangle^0 \parallel \langle \emptyset \rangle^1 \parallel \downarrow_\star^0 \parallel \uparrow_\star^0 \parallel \downarrow_\star^1 \parallel \uparrow_\star^1$$

We have checked this equivalence by writing the two specifications of the Ping-Pong game (with a single, respectively replicated space) in the “tool syntax” (Figure 6(a)), generating the two statespaces and using the model checker to verify that they satisfy the *safety equivalence* relation.

### 4.2 Transparent replication of some Splice applications

Some of the most interesting problems in a system with components are associated with *replication*: what components can be replicated, and at what costs? We claim that the space calculus is a good framework for studying this type of questions. In the sequel we give an example of how our space calculus can be used to rapidly check transparent replication of some applications on Splice.

Consider a simple Splice system, composed of three applications: a *Producer* that writes data to the Splice network, based on observations that it makes on the environment; a *Transformer* that reads the data, applies some transformations on it and writes it back; and a *Consumer* that reads the transformed data items and uses it further, for instance by displaying it on a screen. The producer and the consumer are the components that interact with the environment, while the transformer works “under water”. Therefore it is reasonable to ask whether it is possible to replicate the transformer without affecting the (external) behaviour of the system.

This producer-transformer-consumer example illustrates a specific pattern in Splice systems. The transparent replication of the middle component was extensively studied in [18], using both  $\mu\text{CRL}$  and PVS. We show how to model the problem in space calculus (Figure 6(b)), for the specific instance when two data items are produced, with values 0 and 1. The  $\text{it}_{\text{sp}}$  variable models the local clock. The two specifications are proved safety equivalent by the model checker.

<pre> nfields = 1 upbound = 2 res &lt;*&gt;  space JS (mik.sen.cwi.nl)  app Ping@JS {   write &lt;1&gt;; EXTping; read &lt;0&gt; x;   write &lt;1&gt;; EXTping; read &lt;0&gt; x; } app Pong@JS {   read &lt;1&gt; x; write &lt;0&gt;; EXTpong;   read &lt;1&gt; x; write &lt;0&gt;; EXTpong; } </pre>	<pre> nfields = 1 upbound = 2 res &lt;*&gt;  space JS (mik.sen.cwi.nl) space JSbis (boeg.sen.cwi.nl) JS -&gt; &lt;*&gt; JS &lt;- &lt;*&gt; JSbis -&gt; &lt;*&gt; JSbis &lt;- &lt;*&gt;  app Ping@JS {   write &lt;1&gt;; EXTping; read &lt;0&gt; x;   write &lt;1&gt;; EXTping; read &lt;0&gt; x; } app Pong@JSbis {   read &lt;1&gt; x; write &lt;0&gt;; EXTpong;   read &lt;1&gt; x; write &lt;0&gt;; EXTpong; } </pre>
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(a) A Ping-Pong application on one JavaSpace (left) and on two (right)

<pre> nfields = 3 upbound = 3 space A1 space A2 space A3 A1 -&gt; &lt;1,*,&gt; A2 -&gt; &lt;2,*,&gt; A2 &lt;- &lt;1,*,&gt; 1 3 A3 &lt;- &lt;2,*,&gt; 1 3  app Producer@A1 {   itsp := 0; EXTin;   write &lt;1,0,itsp&gt;;   itsp := itsp + 1;   write &lt;1,1,itsp&gt;; } app Transformer@A2 {   while (true) {     read &lt;1,*,&gt; x;     ivx := x/2+1; itx := x/3;     write &lt;2,ivx,itx&gt;;   }; } app Consumer@A3 {   while (true) {     read &lt;2,*,&gt; x;     EXTout;   }; } </pre>	<pre> nfields = 3 upbound = 3 space A1 space A2 space A3 space A4 A1 -&gt; &lt;1,*,&gt; A2 -&gt; &lt;2,*,&gt; A2 &lt;- &lt;1,*,&gt; 1 3 A3 &lt;- &lt;2,*,&gt; 1 3 A4 -&gt; &lt;2,*,&gt; A4 &lt;- &lt;1,*,&gt; 1 3  app Producer@A1 {   itsp := 0; EXTin;   write &lt;1,0,itsp&gt;;   itsp := itsp + 1;   write &lt;1,1,itsp&gt;; } app Transformer@A2 {   while (true) {     read &lt;1,*,&gt; x;     ivx := x/2+1; itx := x/3;     write &lt;2,ivx,itx&gt;;   }; } app Transformer@A4 {   while (true) {     read &lt;1,*,&gt; x;     ivx := x/2+1;     itx := x/3;     write &lt;2,ivx,itx&gt;;   }; } app Consumer@A3 {   while (true) {     read &lt;2,*,&gt; x;     EXTout;   }; } </pre>
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(b) The Producer/Consumer/Transformer application with one (left) and two (right) transformers

**Fig. 6.** Space calculus program examples

## 5 Conclusions

This paper presents our initial research in a unifying framework for the design and analysis of various distributed dataspace systems. We introduced the space calculus, in which basic concepts of some dataspace paradigms can be modeled. A formal syntax and operational semantics provides a rigorous basis to this calculus.

We aim at two goals: comparing the various paradigms with respect to their meta-properties and facilitating the analysis of individual systems based on heterogeneous shared dataspace architectures.

For the first goal, we view a particular dataspace paradigm as a fragment of the space calculus and we address questions like: does a fragment admit transparent replication of data/processes, what are the costs of a distributed implementation, what are the typical applications for a certain fragment. An answer to the last question would facilitate early architectural design decisions. Some of these questions have been answered for Splice already [11, 18, 20].

The second goal is supported by automatic translations to  $\mu$ CRL and to C. The  $\mu$ CRL specifications can be used as an input to a model checker, thus formally establishing the functional correctness of a system. The approach follows a previous successful attempt for JavaSpaces [21]. The distributed C simulator can be used to find performance bottlenecks in the high-level architecture. These could be solved by transforming the space calculus expression to a functionally equivalent one with a better performance.

As future work, we plan to investigate meta-properties for (fragments) of the space calculus and identify behaviour-preserving transformation rules. Also, we intend to study more examples, in order to establish the validity of our framework and to improve it where necessary. An interesting extension might be to allow dynamic creation of spaces and applications or the dynamic change of the link structure.

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