## ForSyDe-Atom

## User Mānual

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(c)(i)(0)

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## List of Case Studies

George Ungureanu and Ingo Sander (2017). "A layered formal framework for modeling of cyberphysical systems". In: 2017 Design, Automation \& Test in Europe Conference $\mathcal{E}$ Exhibition (DATE). IEEE, pp. 1715-1720
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George Ungureanu, José E. G. de Medeiros, and Ingo Sander (2018). "Bridging discrete and continuous time with Atoms". In: 2018 Design, Automation $\mathfrak{E B}^{2}$ Test in Europe Conference 8 Exhibition (DATE). IEEE
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## Introduction

In this chapter we introduce the purpose, organization and usage of this document, as well as brief instructions and references for helping to set up the ForSyDe-Atom libraries. The scope is to facilitate the reader's progression through this document.

### 1.1 Purpose \& organization

This book is a living document which gathers material related to ForSyDe-Atom and binds it in form of a user manual. Most of the text contained by this book originates from actual inline or literate source code documentation, in form of examples, tutorials, reports and even library API documentation. This means that this document evolves with the ForSyDe-Atom project itself and is periodically updated.

ForSyDe-Atom is a shallow-embedded DSL in the functional programming language Haskell for modeling cyber-physical and parallel systems. It enforces a disciplined way of modeling by separating the manifold concerns of systems into orthogonal layers. The ForSyDe-Atom formal framework aims to providing (where possible) a minimum set of primary common building blocks for each layer called atoms, capturing elementary semantics. Even so, the modeling framework provides library blocks and modules commonly used in CPS defined in terms of patterns of atoms. For more information on ForSyDe-Atom itself, please consult the associated scientific publications or the API extended documentation ${ }^{1}$.

This document is structured in two parts. The first part gathers examples, tutorials and experiments in a learning progression. Each chapter associated with (and actually generated from) a Haskell Cabal project included in the forsyde-atom-examples ${ }^{2}$ repository. This means that the first part of the book is meant to be read in parallel with running the associated example project which conveniently exports functions to test the listed code "on-the-fly". The second part of the book is the actual inline API documentation of ForSyDe-Atom generated with Haddock, which serves also as an extended library report and provides information both on theoretical and implementation issues.

If for any reason you have difficulties following the document or you encounter bugs or discrepancies in the code and text do not hesitate to contact the author(s) or maintainer(s) on GitHub or by email.

### 1.2 Getting ForSyDe-Atom

The ForSyDe-Atom EDSL can be downloaded from https://github.com/forsyde/forsyde-atom. The main page contains enough information for acquiring the dependencies and installing the library on your own. However, each example project from the forsyde-atom-examples repository

[^0]comes with a set of installation scripts written for user convenience, which should be enough for traversing this manual.


#### Abstract

Attention! Be advised that different chapters of the book have been written during different development stages of ForSyDe-Atom and are compatible with different library releases. The code listed throughout the book might be incompatible with the latest release. We strongly recommend using the installation scripts included in each project associated with a chapter, which acquire and install the right dependencies in a local sandbox.


Provided you have an OS installation where the minimum dependencies (GNU make, a Git CLI client, Haskell Platform and cabal-install) are working and accessible by your user profile, to run the getting-started example associated with chapter 2, you simply need to type in the terminal:

```
# download the examples
git clone https://github.com/forsyde/forsyde-atom-examples.git
# change directory to the desired project folder
cd forsyde-atom-examples/getting-started
# install the project (and dependencies) in a sandbox
make install
# open an interpreter session with the examples loaded
cabal repl
```


### 1.3 Using this document

## Disclaimer

The document assumes that the reader is familiar with the syntax of Haskell and the usage of a Haskell interpreter (e.g. ghci). Otherwise, we recommend consulting at least the introductory chapters of one of the following books by Lipovača, 2011 and Hutton, 2016 or other recent books in Haskell.

Most of this document has been created using literate programming. This means that all code shown in the listings is compatible with the ForSyDe-Atom library version mentioned at the beginning of each chapter. Throughout the document you will find two listing styles. This style

```
- / API documentation comment
myIdFunc : : a }->\textrm{a
myIdFunc = id
```

shows source code as it is found in the implementation files. We have taken the liberty to display some code characters as their literate equivalent (e.g. -> is shown as $\rightarrow$, $\backslash$ is shown as $\lambda$, and so on).

This style

```
Prelude> 1 + 1
```

2
suggests interactive commands given by the user in a terminal or an interpreter session. The listing above shows a typical ghci session, where the string after the prompter symbol > suggests the user input (e.g. $1+1$ ). Whenever relevant, the expected output is printed one row below that (e.g. 2).

The code examples are bundled as separate Cabal packages and is provided as libraries meant to be loaded in an interpreter session in parallel with reading this document. Detailed instructions
on how to install the packages can be found in the README.md file in each project. The best way to install the packages is within sandboxed environments with all dependencies taken care of, usually scripted within the make commands. After a successful installation, to open an interpreter session pre-loaded with the main sandboxed library, you just need to type in the following command in a terminal from the package root path (the one containing the .cabal file):

```
|# cabal repl
```

Each section of this document contains a small example written within a library module, like: |module $X$ where

One can access all functions in module X by importing it in the interpreter session, unless otherwise noted (e.g. library X is re-exported by Y ).
|*Y> import X
Now suppose that function myIdFunc above was defined in module $X$, then one would have direct access to it, e.g.:

```
*Y X> :t myIdFunc
myIdFunc :: a -> a
*Y X> myIdFunc 3
3
```

By all means, the code for myIdFunc or any source code for that matter can be copied/pasted in a custom .hs file and compiled or used in any relevant means. The current format was chosen because it is convenient to "get your hands dirty" quickly without thinking of issues associated with compiler suites.

A final tip: if you think that the full name of $X$ is polluting your prompter or is hard to use, then you can import it using an alias:
|*Y> import Extremely.Long.Full.Name.For.X as ShortAlias
|*Y ShortAlias>

### 1.4 References

Hutton, Graham (2016). Programming in Haskell. 2nd. New York, NY, USA: Cambridge University Press. ISBN: $1316626229,9781316626221$.
Lipovača, Miran (2011). Learn You a Haskell for Great Good!: A Beginner's Guide. 1st. San Francisco, CA, USA: No Starch Press. ISBN: 1593272839, 9781593272838.

## Part I

## Examples \& Reports



## Getting Started with ForSyDe-Atom

This chapter is meant to introduce the reader to using the basic features of FORSYDeAtom library as a Haskell EDSL for modeling and simulating embedded and cyberphysical systems. It is not meant to substitute the API documentation nor provide any detail on the theoretical foundation, and it references external documents whenever necessary. It starts from modeling basics, goes through a toy example seen from different perspectives, and into more advanced features like creating custom patterns.

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Info
Compatibility : FORSYDE-ATOM version 0.2.1
Repository : https://github.com/forsyde/forsyde-atom-examples
Path : <repo_root>/getting-started
Other deps. : none

### 2.1 Goals

The main goals of this chapter are:

- introduce the reader to basic modeling features such as: importing library modules, using a Haskell interpreter, using helper functions, composing functions, designing with layers, understanding type signatures, using basic input/output.
- provide a step-by-step guide for modeling a toy system expressing concerns from four layers: function, extended behavior, model of computation and recursive/parallel composition.
- describe the above system as executing with the semantics dictated by four MoCs: synchronous dataflow (SDF), synchronous (SY), discrete event (DE) and continuous time (CT). For this purpose the system will be first instantiated multiple times using specialized helpers, and then described as a network of patterns overloaded with MoC semantics by injecting the right data types, thus exposing the polymorphism of atoms.
- briefly introduce the concepts of atoms and patterns and their usage and guide through creating custom patterns and behaviors.


### 2.2 The basics

This section introduces some basic modeling features of ForSyDe-Atom, such as helpers and process constructors. The module is re-exported by AtomExamples.GettingStarted which is pre-loaded in a repl session, so there is no need to import it manually.
|module AtomExamples.GettingStarted.Basics where
We usually start a ForSyDe-Atom module by importing the ForSyDe.Atom library which provides some commonly used types and utilities.
| import ForSyDe.Atom
In this section we will only test synchronous processes as patterns defined in the MoC layer. An extensive library of types, utilities and helpers for SY process constructors can be used by importing the ForSyDe.Atom.MoC.SY module.
| import ForSyDe.Atom.MoC.SY
Next we import the Absent extended behavior, defined in the ExB layer, to get a glimpse of modeling using multiple layers. As with the previous, we need to specifically import the ForSyDe.Atom.ExB.Absent library to access the helpers and types.
|import ForSyDe.Atom.ExB (res11, res21)
import ForSyDe.Atom.ExB.Absent
The signal is the basic data type defined in the MoC layer, and it encodes a tag system which describes time, causality and other key properties of CPS. In the case of SY MoC, a signal defines a total order between events. There are several ways to instantiate a signal in ForSyDe-Atom. The most usual one is to create it from a list of values using the signal helper. By studying its type signature in the online API documentation, one can see that it needs a list of elements of type a as argument, so let us create a test signal testsig1:

```
|testsig1 = signal [1,2,3,4,5]
```

You can print or check the type of testsig1
*AtomExamples.GettingStarted> testsig1
$\{1,2,3,4,5\}$
*AtomExamples.GettingStarted> :t testsig1
testsig1 :: ForSyDe.Atom.MoC.SY.Core.Signal Integer
The type of testsig1 tells us that the signal helper created a SY Signal carrying Integer values. If you are curious, you can print some information about this mysterious type

```
*AtomExamples.GettingStarted> :info ForSyDe.Atom.MoC.SY.Core.Signal
type ForSyDe.Atom.MoC.SY.Core.Signal a =
    Stream (ForSyDe.Atom.MoC.SY.Core.SY a)
        -- Defined in ForSyDe.Atom.MoC.SY.Core
```

which shows that it is in fact a type alias for a Stream of SY events. If this became too confusing, please read the MoC layer overview in this online API documentation page. Unfortunately the names printed as interactive information are verbose and show their exact location in the structure of ForSyDe-Atom. We do not care about this in the source code, since we imported the SY library properly. To benefit from the same treatment in the interpreter session, we need to do the same:

```
*AtomExamples.GettingStarted> import ForSyDe.Atom.MoC.SY as SY
*AtomExamples.GettingStarted SY> :info Signal
type Signal a = Stream (SY a)
    -- Defined in ForSyDe.Atom.MoC.SY.Core
```

Another way of creating a SY signal is by means of a generate process, which generates an infinite signal from a kernel value. By studying the online API documentation, you can see that the SY library provides a number of helpers for this particular process constructor, the one generating one output signal being generate1. Let us first check the type signature for this helper function:

```
*AtomExamples.GettingStarted SY> :t generate1
```

generate1 :: (b1 -> b1) -> b1 -> Signal b1

So basically, as suggested in the online API documentation, this helper takes a "next state" function of type a -> a, a kernel value of type a, and it generates a signal of tyle Signal a. With this in mind, let us create testsig2:
|testsig2 = generate1 (+1) 0
Printing it would jam the terminal... we were serious when we said "infinite"! This is why you need to select a few events from the beginning to see whether the signal generator behaves correctly. To do so, we use the takeS utility:

```
*AtomExamples.GettingStarted SY> takeS 10 testsig2
{0,1,2,3,4,5,6,7,8,9}
*AtomExamples.GettingStarted SY> :t testsig2
testsig2 :: Signal Integer
```

generate was a process with no inputs. Now let us try a process that takes the two signals testsig1 and testsig2 and sums their synchronous events. For this we use the combinatorial process comb, and the SY constructor we need, with two inputs and one output is provided by comb21. Again, checking the type signature confirms that this is the helper we need:
*AtomExamples.GettingStarted SY> :t comb21
comb21 :: (a1 -> a2 -> b1) -> Signal a1 -> Signal a2 -> Signal b1
So let us instantiate testproc1:

```
|testproc1 = comb21 (+)
```

Calling it in the interpreter with testsig1 and testsig2 as arguments, it returns:

```
*AtomExamples.GettingStarted SY> testproc1 testsig2 testsig1
```

\{1,3,5, 7, 9$\}$
which is the expected output, as based on the definition of the SY MoC, all events following the sixth one from testsig2 are not synchronous to any event in testsig1.


Figure 2.1: Simple process network as composition of processes
Suppose we want to increment every event of testsig2 with 1 before summing the two signals. This particular behavior is described by the process network in fig. 2.1. There are multiple ways to instantiate this process network, mainly depending on the coding style of the user:

```
testpn1 = comb21 (+) . comb11 (+1)
testpn2 s2 s1 = testproc1 (comb11 (+1) s2) s1
testpn3 s2 = testproc1 (comb11 (+1) s2)
testpn4 s2 s1 = let s2' = comb11 (+1) s2
    in testproc1 s2' s1
testpn5 s2 = comb21 (+) s2'
    where s2' = comb11 (+1) s2
```

All of the above functions are equivalent. testpn1 uses the point-free notation, i.e. the function composition operator, between two partially-applied process constructors. testpn2 makes use of the previously-defined testproc1 to enforce a global hierarchy. testpn3 is practically the same, but it exposes the partial application mechanism, by not making the s1 argument explicit. testpn4 makes use of local hierarchy in form of a let-binding, while testpn5 does so through a where clause. Printing them only confirms their equivalence:

```
*AtomExamples.GettingStarted SY> testpn1 testsig2 testsig1
{2,4,6,8,10}
*AtomExamples.GettingStarted SY> testpn2 testsig2 testsig1
{2,4,6,8,10}
*AtomExamples.GettingStarted SY> testpn3 testsig2 testsig1
{2,4,6,8,10}
*AtomExamples.GettingStarted SY> testpn4 testsig2 testsig1
{2,4,6,8,10}
*AtomExamples.GettingStarted SY> testpn5 testsig2 testsig1
{2,4,6,8,10}
```

One key concept of ForSyDe-Atom is the ability to model different aspects of a system as orthogonal layers. Up until now we only experimented with two layers: the MoC layer, which concerns timing and synchronization issues, and the function layer, which concerns functional aspects, such as arithmetic and data computation. Let us rewind which DSL blocks we have used an group them by which layer they belong:

- the Integer values carried by the two test signals (i.e. $0,1, \ldots$ ) and the arithmetic functions (i.e. (+) and (+1)) belong to the function layer.
- the signal structures for testsig1 and testsig2 (i.e. Signal a) the utility (i.e. signal) and the process constructors (i.e. generate1, comb11 and comb21) belong to the MoC layer.

It is easy to grasp the concept of layers once you understand how higher order functions work, and accept that ForSyDe-Atom basically relies on the power of functional programming to define structured abstractions. In the previous case entities from the MoC layer "wrap around" entities from the function layer like in fig. 2.2, "lifting" them into the MoC domain. Unfortunately it is not that straightforward to see from the code syntax which entity belongs to which layer. For now, their membership can be determined solely by the user's knowledge of where each function is defined, i.e. in which module. Later in this guide we will make this apparent from the code syntax, but for now you will have to trust us.


Figure 2.2: Layered structure of the processes in fig. 2.1


Figure 2.3: Layered structure of the processes describing absent events

As a last exercise for this section we would like to extend the behavior of the system in fig. 2.1 in order to describe whether the events are happening or not (i.e. are absent or present) and act accordingly. For this, ForSyDe-Atom defines the Extended Behavior (ExB) layer. As suggested in fig. 2.3, this layer extends the pool of values with symbols denoting states which would be impossible to describe using normal values, and associates some default behaviors (e.g. protocols) over these symbols.

The two processes fig. 2.1 are now defined below as testAp1 and testAp2. This time, apart from the functional definition (name $=$ function) we specify the type signature as well (name : : type), which in the most general case can be considered a specification/contract of the interfaces of the newly instantiated component. Both type signatures and function definitions expose the layered structure suggested in fig. 2.3. As specific ExB type, we use AbstExt and as behavior pattern constructor we choose a default behavior expressing a resolution res.

```
testAp1 :: Num a
    => Signal (AbstExt a) - - input signal of absent-extended values
    Signal (AbstExt a) - output signal of absent-extended values
testAp1 = comb11 (res11 (+1))
testAp2 :: Num a
    => Signal (AbstExt a) - first input signal of absent-extended values
    Signal (AbstExt a) - second input signal of absent-extended values
    Signal (AbstExt a) - output signal of absent-extended values
testAp2 = comb21 (res21 (+))
```

Now all we need is to create some test signals of type Signal (AbstExt a). One way is to use the signal utility like for testAsig1, but this forces to make use of AbstExt's type constructors. Another way is to use the library-provided process constructor helpers, such as filter', like for testAsig2.

```
testAsig1 = signal [Prst 1, Prst 2, Abst, Prst 4, Abst]
testAsig2 = filter' (/=4) testsig2
```

Printing out the test signals in the interpreter session this is what we get:

```
*AtomExamples.GettingStarted SY> testAsig1
{1,2,\perp,4,\perp}
*AtomExamples.GettingStarted SY> takeS 10 testAsig2
{0,1,2,3,\perp,5,6,7,8,9}
```

Trying out testAp1 on testAsig1:

```
*AtomExamples.GettingStarted SY> testAp1 testAsig1
{2,3,\perp,5,\perp}
```

Everything seems all right. Now testing testAp2 on testAsig1 and testAsig2:

```
*AtomExamples.GettingStarted SY> takeS 10 $ testAp2 testAsig2 testAsig1
```

$\{1,3, * * *$ Exception: [ExB.Absent] Illegal occurrence of an absent and present event

Uh oh... Actually this is the correct behavior of a resolution function for absent events, as defined in synchronous reactive languages such as Lustre Halbwachs et al., 1991. Let us remedy the situation, but this time using another library-provided process constructor, when'.

```
testAsig2' = when' mask testsig2
    where
        mask = signal [True, True, False, True, False]
```

Now printing testAp2 looks better:

```
*AtomExamples.GettingStarted SY> testAsig2'
{0,1,\perp,3,\perp}
*AtomExamples.GettingStarted SY> testAp2 testAsig2' testAsig1
{1,3,\perp,7,\perp}
```

And recreating the process network from fig. 2.1 gives the expected result:

```
|testApn1 = testAp2 . testAp1
*AtomExamples.GettingStarted SY> testApn1 testAsig2' testAsig1
{2,4,\perp,8,\perp}
```

This section has provided a crash course in modeling with ForSyDe-Atom, with focus on a few practical matters, such as using library-provided helpers and constructors and understanding the role of layers. The following sections delve deeper into modeling concepts such as atoms and making use of ad-hoc polymorphism.

### 2.2.1 Visualizing your data

Up until now, we have made use of the Show instance of the ForSyDe-Atom data types to print out signals on the terminal screen. While this remains the main way to test if a model is working properly, there are alternative ways to plot data. This section introduces the reader to the ForSyDe.Atom.Utility. Plot library of utilities for visualizing signals or other data types.

The functions presented in this section are defined in the following module, which is exported by AtomExamples.GettingStarted.
|module AtomExamples.GettingStarted.Plot where
We will be using the signals defined in the previous section, so let us import the corresponding module:
| import AtomExamples.GettingStarted.Basics
And, as mentioned, we need to import the library with plotting utilities:
| import ForSyDe.Atom.Utility.Plot
Upon consulting the API documentation for this module, you might notice that most utilities input a so-called PlotData type, which is an alias for a complex structure carrying configuration parameters, type information and data samples. Using Haskell's type classes, ForSyDe-Atom is able to provide few polymorphic utilities for converting most of the useful types into PlotData.

For example, the prepare function takes a "plottable" data type (e.g. a signal of values), and a Config type, and returns PlotData. The Config type is merely a record of configuration parameters useful further down in the plotting pipeline. At the time of writing this report ${ }^{1}$, a configuration record looked like this:

```
config = 
    , file = "plot" - base name of the eventual files generated
    , rate = 0.01 - sampling rate if relevant (e.g. ignored by SY signals).
    - Useful just for e.g. explicit-timed signals.
    , xmax = 20 - maximum x coordinate. Necessary for infinite signals.
    , labels = ["s1","s2"] - labels for all signals passed to be plotted.
    , verbose = True - prints additional messages for each utility.
    , fire = True - if relevant, fires a plotting or compiling program.
    , mklatex = True - if relevant, dumps a LaTeX script loading the plot.
    , mkeps = True - if relevant. dumps a PostScript file with the plot.
    , mkpdf = True - if relevant, dumps a PDF file with the plot.
```

ForSyDe.Atom.Utility.Plot provides several of these pre-made configuration objects, which can be modified on-the-fly using Haskell's record syntax, as you will see further on.

Let us see again the signals testsig1 and testsig2 defined in the previous section:

```
\lambda> testsig1
{1,2,3,4,5}
\lambda> takeS 20 testsig2
{0,1,2,3,4,5,6,7,8,9,10,11,12,13,14,15,16,17,18,19}
```

The utility showDat prints out sampled data on the terminal, as pairs of X and Y coordinates:

```
show1 = showDat $ prepare config testsig1
show2 = let cfg = config {xmax=15, labels=["testsig1","testsig2"]}
    in showDat $ prepareL cfg [testsig1,testsig2]
```

| 0 | 1.0 |
| :---: | :---: |
| 1 | 2.0 |
| 2 | 3.0 |
| 3 | 4.0 |
| 4 | 5.0 |
| $<$ |  |
| $\lambda>$ show2 |  |

[^1]| testsig1 $=$ |  |
| ---: | :---: |
| 0 | 1.0 |
| 1 | 2.0 |
| 2 | 3.0 |
| 3 | 4.0 |
| 4 | 5.0 |
| testsig2 |  |
| 0 | 0.0 |
| 1 | 1.0 |
| 2 | 2.0 |
| 3 | 3.0 |
| 4 | 4.0 |
| 5 | 5.0 |
| 6 | 6.0 |
| 7 | 7.0 |
| 8 | 8.0 |
| 9 | 9.0 |
| 10 | 10.0 |
| 11 | 11.0 |
| 12 | 12.0 |
| 13 | 13.0 |
| 14 | 14.0 |

The function dumpDat dumps the data files in a path specified by the configuration object. Based on the config object instantiated earlier, after calling the following function you should see a new folder called fig in the current path, with two new . dat files.

```
dump2 = let cfg = config {xmax=15, labels=["testsig1","testsig2"]}
    in dumpDat $ prepareL cfg [testsig1,testsig2]
```

```
|> dump2
Dumped testsig1, testsig2 in ./fig
["./fig/testsig1.dat","./fig/testsig2.dat"]
```

The plot library also has a few functions which create and (optionally) fire Gnuplot scripts. In order to make use of them, you need to install the dependencies mentioned in the API documentation. For example, using the function plotGnu creates the following plot:

```
plot2 = let cfg = config {xmax=8, labels=["testsig1","testsig2"]}
    in plotGnu $ prepareL cfg [testsig1,testsig2]
```

Different input data creates different types of plots, as we will see in future sections.
One can also generate $\mathrm{IA}_{\mathrm{E}} \mathrm{X}$ code which is meant to be compiled with the ForSyDe-IATEX package, more specifically its signal plotting library. Check the user manual for more details on how to install the dependencies and how to use the library itself. Naturally, there is a function showLatex which prints out the command for a signals environment defined in FORSYDE-EATEX:

```
latex1 = let cfg = config {xmax=8, labels=["testsig1","testsig2"]}
    in showLatex $ prepareL cfg [testsig1,testsig2]
```

$\mid \lambda>$ latex 1
\begin\{signalsSY\} [] \{8.0\} }
\signalSY[]\{1.0:0,2.0:1,3.0:2,4.0:3,5.0:4\}
\signalSY[]\{0.0:0,1.0:1,2.0:2,3.0:3,4.0:4,5.0:5,6.0:6,7.0:7\}
\end\{signalsSY\} }
<

Also, there is a command plotLatex for generating a standalone $\mathrm{A}_{\mathrm{E}} \mathrm{EXX}_{\mathrm{E}}$ document and, if possible, compiling it with pdflatex. For example, calling the following function generates the image from section 2.2.1.

```
latex2 = let cfg = config {xmax=8, labels=["testsig1","testsig2"]}
    in plotLatex $ prepareL cfg [testsig1,testsig2]
```

The SY signal plot is nothing spectacular, but wraps the events in a matrix of nodes which can be embedded into a more complex TikZ figure. Other signals produce other plots. Most generated plots will need manual tweaking in order to look good. Check the user manual on how to customize each plot.


Figure 2.4: SY signal in Gnuplot as a impulse plot

| 1.0 | 2.0 | 3.0 | 4.0 | 5.0 |  |  |  |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| 0.0 | 1.0 | 2.0 | 3.0 | 4.0 | 5.0 | 6.0 | 7.0 |

Figure 2.5: SY signal plot in ForSyDe- ${ }^{\mathrm{A}} \mathrm{T}_{\mathrm{E}} \mathrm{X}$ as a matrix of nodes

### 2.3 Toy example: a focus on MoCs

This example has been used as a case study for introducing the new concepts of ForSyDe-Atom in the paper of Ungureanu and Sander, 2017. It describes the simple system from fig. 2.6b which exposes four layers, structured like in fig. 2.6a. This system is then fed vectors of signals describing different MoCs and its response is observed. In figs. 2.6 c to 2.6 f some possible projections on the different layers are depicted. For now they are used just as trivia, and you need not bother with theritibats msydhthetic example meant to introduce as many concepts as possible in a short amount of time and, among others, it highlights:

- the power of partial application for creating parameterized structures, such as the process network for stage1.
- alternative designs for the same toy system to show the effect of MoCs. First it is instantiated using different process constructor helpers defined for each MoC separately. Afterwards it is written as one single polymorphic instance using MoC layer patterns, overloaded with execution semantics in accordance with the tag system injected into the system.

For the sake of brevity, we also provide the functional description in the language introduced by Ungureanu and Sander, 2017 in eqs. (2.1)-(2.5) and table 2.1. Do not bother much about this notation either, as this exact definition will appear in the code in a more "human readable" form.

$$
\begin{align*}
& \text { toy : }\langle V\rangle \rightarrow\langle S\rangle \rightarrow S  \tag{2.1}\\
& \text { toy }\langle i\rangle\langle s\rangle=\operatorname{when}_{\mathrm{M}}\left(\Gamma_{w} \vdash w_{\mathrm{B}}\right) \circ \operatorname{reduce}_{\mathrm{S}}\left(r_{\mathrm{M}}\right) \circ \operatorname{map}_{\mathrm{S}}\left(p c_{\mathrm{M}}\right)\langle i\rangle\langle s\rangle
\end{align*}
$$



Figure 2.6. Views and projections for the toy system
where

$$
\begin{align*}
\operatorname{when}_{\mathrm{M}}\left(\Gamma_{w} \vdash w_{\mathrm{B}}\right)(s) & =\left((\perp \notin) \circ\left(\Gamma_{w} \vdash w_{\mathrm{B}}\right)\right) \oplus s  \tag{2.2}\\
r_{\mathrm{M}}(x, y) & =\Gamma_{r} \vdash r_{\mathrm{B}} \oplus(x, y)  \tag{2.3}\\
\operatorname{map}_{\mathrm{s}}\left(p c_{\mathrm{M}}\right)\langle v\rangle\langle s\rangle & =p c_{\mathrm{M}} \diamond\langle v\rangle *\langle s\rangle  \tag{2.4}\\
p c_{\mathrm{M}}(x, y) & =\operatorname{moore}_{\mathrm{M}}\left(\Gamma_{n s} \vdash n s_{\mathrm{B}}, \Gamma_{o d} \vdash o d_{\mathrm{B}}, x\right)(y) \tag{2.5}
\end{align*}
$$

Table 2.1: Contexts, Functions and Initial Tokens for the System in Eq. (2.1)

| MoC | $\Gamma_{w} \vdash$ | $w_{\text {B }}(x)$ | $\Gamma_{r} \quad \vdash r_{\mathrm{B}}(x, y)$ | $\Gamma_{n s} \vdash$ | $n s_{\mathrm{B}}(x, y)$ | $\Gamma_{o d} \vdash$ od ${ }_{\mathrm{B}}(x)$ | $\langle i\rangle=\langle(t, v)\rangle$ |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $\mathrm{SDF}^{2}$ | $2,2 \vdash\left(x_{1}\right.$ | $\left.<0, x_{2}<0\right)^{ \pm}$ | $(1,1), 1 \vdash\left(x_{1}+y_{1}\right)^{\triangle}$ | $\begin{aligned} &(1,2), 1 \vdash\left(x_{1}+y_{1}+y_{2}\right)^{\mathbb{A}} \\ & \vdash(x+y)^{\mathbb{A}} \\ & \vdash(x+y)^{\mathbb{®}} \\ & \vdash(x+y)^{\mathbb{®}} \end{aligned}$ |  | $\begin{aligned} 1,1 & \vdash x_{1} \mathbb{A} \\ & \vdash x \mathbb{A} \\ \vdash & \not \mathbb{A}_{\mathbb{A}} \\ & \vdash x \mathbb{A} \end{aligned}$ | $\begin{array}{\|ll} \langle & (,-1) \\ \langle & (,-1) \\ \langle & (0.5,-1) \\ \langle(1, \lambda t \rightarrow-1) \end{array}$ | $(, 1)$ | $(,-1)$ | $(, 1)$ | > |
| SY | $\vdash$ | $(x<0)^{\text {A }}$ | $\vdash(x+y)^{\mathbb{A}}$ |  |  | $(\mathrm{t})$ |  | $(,,-1)$ | $(\mathrm{l}, 1)$ | > |
| DE | $\vdash$ | $(x<0)^{\mathbb{A}}$ | $\vdash(x+y)^{\mathbb{A}}$ |  |  | $(1.4,1)$ |  | $(1.0,-1)$ | $(1.4,1)$ | ) |
| CT | $\vdash$ | $(x<0)^{\text {A }}$ | $\vdash(x+y)^{\mathbb{A}}$ |  |  | $(1.4, \lambda t \rightarrow$ |  | ) $(1, \lambda t \rightarrow-1)$ | $(1.4, \lambda t \rightarrow$ |  |

[^2]
### 2.3.1 Test input signals

In the following examples we will use a set of test signals defined in the following module, which is also re-exported by AtomExamples.GettingStarted (i.e. you don't need to import it):
|module AtomExamples.GettingStarted.TestSignals where
The test signals need to define tag systems belonging to different MoCs. Each MoC has an own dedicated module under ForSyDe. Atom. MoC which defines atoms, patterns, types and utilities. Just like in the previous section, we need to import the needed modules. This time we name them using short aliases, to disambiguate between the different DSL items, often sharing the same name, but defined in different places.

| import ForSyDe.Atom.ExB.Absent (AbstExt (..)) |  |
| :--- | :--- | :--- |
| import ForSyDe.Atom.MoC.SY | as SY |
| import ForSyDe.Atom.MoC.DE | as DE |
| import ForSyDe.Atom.MoC.CT | as CT |
| import ForSyDe.Atom.MoC.SDF | as SDF |
| import ForSyDe.Atom.MoC.Time | as T |
| import ForSyDe.Atom.MoC.TimeStamp | as Ts |
| import ForSyDe.Atom.Skeleton.Vector | as V |
| import ForSyDe.Atom.Utility.Plot |  |

Let the signals sdf1-sdf4 denote four SDF signals, i.e. sequences of events. Instead of using the signal utility, we use readSignal which reads a string, tokenizes it and converts it to a SDF signal. This utility function needs to be "steered" into deciding which data type to output so in order to specify the type signature we use the inline Hakell syntax name = definition :: type. All events, although extended, are present.

```
sdf1 = SDF.readSignal "{ 1, 1, 1, 1, 1, 1}" :: SDF.Signal (AbstExt Int)
sdf2 = SDF.readSignal "{-1, 1, -1, 1,-1, 1}" :: SDF.Signal (AbstExt Int)
sdf3 = SDF.readSignal "{ 0, 0, 1, 1, 0 }" :: SDF.Signal (AbstExt Int)
sdf4 = SDF.readSignal "{-1,-1,-1,-1,-1 }" :: SDF.Signal (AbstExt Int)
```

Similarly, let the signals sy1-sy4 denote four SY signals, i.e. all events are synchronized with each other. We use the SY version of readSignal, also "steered" by declaring the types inline, and all events are also present.

```
sy1 = SY.readSignal "{ 1, 1, 1, 1, 1, 1}" :: SY.Signal (AbstExt Int)
sy2 = SY.readSignal "{-1, 1, -1, 1, -1, 1}" :: SY.Signal (AbstExt Int)
sy3 = SY.readSignal "{ 0, 0, 1, 1, 0 }" :: SY.Signal (AbstExt Int)
sy4 = SY.readSignal "{-1,-1,-1,-1,-1 }" :: SY.Signal (AbstExt Int)
```

For the DE signals de1-de4 we need to specify for each event an explicit tag (i.e. timestamp), as required by the DE tag system. For this, the DE version of readSignal reads each event using the syntax value@timestamp. Needless to say, all events are also present.


Let ct1-ct4 denote four CT signals. As the events in a CT signal are themselves continuous functions of time, we cannot specify them as mere strings, thus we cannot use a readSignal utility any more. This time we will use the CT version of the signal utility, where each event is specified as a tuple (timestamp, $f(t)$ ), and can be considered as a continuous sub-signal. For representing time we use an alias Time for Rational, defined in the ForSyDe.Atom.MoC.Time. This module also contains utility functions of time, such as the constant function const or the sine sin. We define local functions to wrap the type returned by a CT subsignal into an AbstExt type.

```
pconst = T.const . Prst
ct1 = CT.signal [(0,pconst 1)] :: CT.Signal (AbstExt Time)
ct2 = CT.signal [(0,Prst . ( }\lambda\textrm{t}->\textrm{T}.\operatorname{sin (T.pi * t)))] :: CT.Signal (AbstExt Time)
ct3 = CT.signal [(0,pconst 0),(1.4,pconst 1),(2.8,pconst 0)] :: CT.Signal (AbstExt Time)
ct4 = CT.signal [(0,pconst (-1))] :: CT.Signal (AbstExt Time)
```

Finally, we need to bundle these signals into vectors of signals, to feed into the toy system from fig. 2.6b and eq. (2.1). For this purpose we pass the four signals of each MoC as a list to the vector utility.

```
vsdf = V.vector [sdf1, sdf2, sdf3, sdf4] :: V.Vector (SDF.Signal (AbstExt Int))
vsy = V.vector [ sy1, sy2, sy3, sy4] :: V.Vector ( SY.Signal (AbstExt Int))
vde = V.vector [ de1, de2, de3, de4] :: V.Vector ( DE.Signal (AbstExt Int))
vct = V.vector [ ct1, ct2, ct3, ct4] :: V.Vector ( CT.Signal (AbstExt Time))
```

Now we need to create for each MoC the vectors with the initial states for the Moore machines, i.e. $\langle i\rangle$ from table 2.1. For SDF, SY and DE we are also making use of the fact that the defined data types are Readable. The data types that we need within the vectors are in accordance to the moore process constructor helpers defined in each module, so be sure to check the online API documentation to understand why we need those particular types. For example, SDF requires a partition (list) of values as initial states whereas DE apart from a value it requires also the duration of the first event. For the CT vector, as before, we cannot read functions as string, so we use the vector utility. ict also shows the usage of the milisec utility which converts an integer into a timestamp.

```
isdf = read "<[-1],[ 1],[-1],[ 1]>" :: V.Vector [AbstExt Int]
isy = read "<(-1), 1, (-1), 1 >" :: V.Vector (AbstExt Int)
ide = read "< (1,-1),(1.4, 1),(1.0,-1),(1.4, 1)>"
:: V.Vector (TimeStamp, AbstExt Int)
ict = V.vector [
    (Ts.milisec 1000, pconst (-1)),
    (Ts.milisec 1400, pconst 1 ),
    (Ts.milisec 1000, pconst (-1)),
    (Ts.milisec 1400, pconst 1 )] :: V.Vector (TimeStamp, Time }->\mathrm{ AbstExt Time)
```

For the polymorphic instance in section 2.3 .6 we need to provide the initial states as wrapped in signals, so we create unit signals:

| sisdf $=$ SDF.signal | $<\$>$ isdf |
| :--- | :--- |
| sisy $=$ SY.unit | $<\$>$ isy |
| side $=$ DE.unit | $<\$>$ ide |
| sict $=$ CT.unit | $<\$>$ ict |

Finally, let us instantiate some plotting utilities, to test the different signals throughout the experiments:

```
plot until lbls = plotGnu . prepare defaultCfg {xmax=until, labels=lbls, rate=0.01}
latex until lbls = plotLatex . prepare defaultCfg {xmax=until, labels=lbls, rate=0.01}
plotV until lbls = plotGnu . prepareV defaultCfg {xmax=until, labels=lbls, rate=0.01}
latexV until lbls = plotLatex . prepareV defaultCfg {xmax=until, labels=lbls, rate=0.01}
```


### 2.3.2 SY instance

The SY instance of the toy system is created using process constructor helpers defined in ForSyDe.Atom.MoC.SY, and is defined in the following module (re-exported by AtomExamples.GettingStarted).
|module AtomExamples.GettingStarted.SY where
As in the previous examples we import the modules we need, and use aliases for referencing them in the code.

```
import ForSyDe.Atom
import ForSyDe.Atom.ExB.Absent (AbstExt)
import ForSyDe.Atom.ExB as ExB
import ForSyDe.Atom.MoC.SY as SY
import ForSyDe.Atom.Skeleton.Vector as V
```

Although Haskell's type engine can infer these type signatures, for the sake of documenting the interfaces for each stage, we will explicitly write their types. First, stage1 is defined as a farm network of moore processes, where the initial states are provided by a vector. Its definition makes use of partial application (i.e. arguments which are not explicitly written are supposed to be the same on the LHS as on the RHS). It is defined hierarchically, making use of local name bindings after the where keyword.

```
stage1SY :: V.Vector (AbstExt Int) - vector of initial states
    ->V.Vector (SY.Signal (AbstExt Int)) - - vector of input signals
    -> V.Vector (SY.Signal (AbstExt Int)) - ~ vector of output signals
```

```
stage1SY = V.farm21 pcSY
    where
        pcSY = SY.moore11 ns od
        ns = ExB.res21 (+)
        od = ExB.res11 id
```

Let us print and plot the inputs against the outputs, using the test signals and plotting functions latexV and plotV defined in section 2.3.1:

```
l> isy 
\lambda> vsy
<{1,1,1,1,1,1},{-1,1,-1,1,-1,1},{0,0,1,1,0},{-1,-1,-1,-1,-1}>
\lambda> stage1SY isy vsy
<{-1,0,1,2,3,4,5},{1,0,1,0,1,0,1},{-1,-1,-1,0,1,1},{1,0,-1,-2,-3,-4}>
\lambda> let latexIn = latexV 6 ["sy1","sy2","sy3","sy4"] vsy
\lambda> let latexS1 = latexV 7 ["sy1-1","sy2-1","sy3-1","sy4-1"] $ stage1SY isy vsy
\lambda> let gnuIn = plotV 6 ["sy1","sy2","sy3","sy4"] vsy
\lambda> let gnuS1 = plotV 7 ["sy1-1","sy2-1","sy3-1","sy4-1"] $ stage1SY isy vsy
```

| 1 | 1 | 1 | 1 | 1 | 1 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| -1 | 1 | -1 | 1 | -1 | 1 |
| 0 | 0 | 1 | 1 | 0 |  |
| -1 | -1 | -1 | -1 | -1 |  |

(a) latexIn

| -1 | 0 | 1 | 2 | 3 | 4 | 5 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | 0 | 1 | 0 | 1 | 0 | 1 |
| -1 | -1 | -1 | 0 | 1 | 1 |  |
| 1 | 0 | -1 | -2 | -3 | -4 |  |


(c) gnuIn
(b) latexS1


(d) gnuS1

The second stage of the toy system in fig. 2.6 b is defined as a reduce network of comb processes. As seen in its type signature, it inputs a vector of signals and it reduces it to a single signal.

```
stage2SY :: V.Vector (SY.Signal (AbstExt Int))
        SY.Signal (AbstExt Int)
stage2SY = V.reduce rSY
    where
        rSY = SY.comb21 (ExB.res21 (+))
```

Again, let us print and plot the output signals using the test inputs and utilities defined in section 2.3.1.

```
\lambda> let s2out = (stage2SY . stage1SY isy) vsy
\lambda> s2out
{0,-1,0,0,2,1}
\lambda> let latexS2 = latex 7 ["sy1-2","sy2-2","sy3-2","sy4-2"] s2out
\lambda> let gnuS2 = plot 7 ["sy1-2","sy2-2","sy3-2","sy4-2"] s2out
```

Finally, the last stage of the toy system applies a filter pattern on the reduced signal to mark all values less than 0 as absent.

```
stage3SY :: SY.Signal (AbstExt Int)
    -> SY.Signal (AbstExt Int)
```



```
stage3SY = SY.filter (>=0)
>
toySY :: V.Vector (AbstExt Int) - - initial states
    -> V.Vector (SY.Signal (AbstExt Int)) - - input
    -> SY.Signal (AbstExt Int) - - output
toySY i = stage3SY . stage2SY . stage1SY i
```

We print and plot the system response to the test signals defined in section 2.3.1.

```
\lambda> toySY isy vsy
{0, \perp,0,0,2,1}
\lambda> let latexS3 = latex 6 ["sy1-3"] $ toySY isy vsy
\lambda> let gnuS3 = plot 6 ["sy1-3"] $ toySY isy vsy
```



### 2.3.3 DE instance

The DE instance of the toy looks exactly the same as the SY instance in section 2.3.2, but is created using constructors from the ForSyDe.Atom.MoC.DE module. This is why we will skip most of the description, and jump straight to testing it. The following file, as you are used to by now, is re-exported by AtomExamples.GettingStarted.

## |module AtomExamples.GettingStarted.DE where

As previously, we use aliases for the imported modules.

```
import ForSyDe.Atom
import ForSyDe.Atom.ExB.Absent (AbstExt)
import ForSyDe.Atom.ExB as ExB
import ForSyDe.Atom.MoC.DE as DE
import ForSyDe.Atom.Skeleton.Vector as V
```

Again, we make the type signatures explicit for documentation purpose. For stage1 we use the same farm network but now using DE moore processes.

```
stage1DE :: V.Vector (TimeStamp, AbstExt Int) - ^ vector of initial states
    -> V.Vector (DE.Signal (AbstExt Int)) - - vector of input signals
    -> V.Vector (DE.Signal (AbstExt Int)) - ~ vector of output signals
```

```
stage1DE = V.farm21 pcDE
    where
        pcDE = DE.moore11 ns od
        ns = ExB.res21 (+)
        od = ExB.res11 id
```

When printing ide and vde we can see the effects of rounding the input floating point numbers to the nearest discrete timestamp. We also have to take into account that the DE version of the Moore machine produces infinite signals when we print them out. Notice that the generated graphs may need to be tweaked in order to show the information properly.

```
\lambda> ide
<(1s, -1),(1.399999999999s,1), (1s, -1),(1.399999999999s, 1)>
\lambda> vde
<{1 @0s},{ -1 @0s, 1 @0.699999999999s, -1 @1.399999999999s, 1 @2.1s, -1 @2.799999999999s, 1 @3.5s
    },{ 0 @0s, 1 @1.399999999999s, 0 @2.799999999999s},{ -1 @0s}>
\lambda> fmap (takeS 6) $ stage1DE ide vde
<{ -1 @0s, 0 @1s, 1 @2s, 2 @3s, 3 @4s, 4 @5s},{ 1 @0s, 0 @1.399999999999s, 2 @2.099999999998s, -1
    @2.799999999998s, 1 @3.499999999997s, 3 @3.499999999999s},{ -1 @0s, -1 @1s, -1 @2s, 0 @2
    .399999999999s, 0 @3s, 1 @3.399999999999s},{ 1 @0s, 0 @1.399999999999s, -1 @2.799999999998s, -2
        @4.199999999997s, -3 @5.599999999996s, -4 @6.999999999995s}>
\lambda> let latexIn = latexV 3.3 ["de1","de2","de3","de4"] vde
\lambda> let latexS1 = latexV 3.3 ["de1-1","de2-1","de3-1","de4-1"] $ stage1DE ide vde
\lambda> let gnuIn = plotV 3.3 ["de1","de2","de3","de4"] vde
\lambda> let gnuS1 = plotV 3.3 ["de1-1","de2-1","de3-1","de4-1"] $ stage1DE ide vde
```


(a) latexIn

(c) gnuIn

(b) latexS1
(d) gnuS1

The second stage according to fig. 2.6b is also defined as a reduce network but this time we use the DE process constructor for the comb processes.

```
stage2DE :: V.Vector (DE.Signal (AbstExt Int))
    DE.Signal (AbstExt Int)
stage2DE = V.reduce rDE
    where
        rDE = DE.comb21 (ExB.res21 (+))
```

```
\lambda> let s2out = (stage2DE . stage1DE ide) vde
\lambda> takeS 10 s2out
{ 0 @0s, 1 @1s, -1 @1.399999999999s, 0 @2s, 2 @2.099999999998s, 3 @2.399999999999s, -1 @2
    .799999999998s, 0 @3s, 1 @3.399999999999s, 3 @3.499999999997s}
\lambda> let latexS2 = latex 3.3 ["de1-2","de2-2","de3-2","de4-2"] s2out
\lambda> let gnuS2 = plot 3.3 ["de1-2","de2-2","de3-2","de4-2"] s2out
```

For the last stage of the toy system there is no DE process constructor in ForSyDe.Atom.MoC.DE so we need to create it ourselves.

(a) latexS2

(b) gnuS2

```
stage3DE :: DE.Signal (AbstExt Int)
    -> DE.Signal (AbstExt Int)
stage3DE = deFilter (>=0)
    where
        deFilter p s = DE.comb21 ExB.filter (predSig p s) s
        predSig p s = DE.comb11 (ExB.res11 p) s
toyDE :: V.Vector (TimeStamp, AbstExt Int) - ~ initial states
        ->V.Vector (DE.Signal (AbstExt Int)) - - input
        DE.Signal (AbstExt Int) - - output
toyDE i = stage3DE . stage2DE . stage1DE i
```

```
\lambda> takeS 10 $ toyDE ide vde
{ 0 @0s, 1 @1s, \perp@1.399999999999s, 0 @2s, 2 @2.099999999998s, 3 @2.399999999999s, \perp @2
    .799999999998s, 0 @3s, 1 @3.399999999999s, 3 @3.499999999997s}
\lambda> let latexS3 = latex 3.3 ["de1-3"] $ toyDE ide vde
\lambda> let gnuS3 = plot 3.3 ["de1-3"] $ toyDE ide vde
```



The CT instance of the toy looks exactly the same as the previous ones in section 2.3.2 and ??, but is created using constructors from the ForSyCt.Atom.MoC.CT module. The following file, as you are used to by now, is re-exported by AtomExamples.GettingStarted.

## |module AtomExamples.GettingStarted.CT where

As prevoiusly, we use aliases for the imported modules.

| import ForSyDe.Atom |  |
| :--- | ---: |
| import ForSyDe.Atom.ExB.Absent | (AbstExt) |
| import ForSyDe.Atom.ExB | as ExB |
| import ForSyDe.Atom.MoC.CT | as CT |
| import ForSyDe.Atom.Skeleton.Vector | as V |

Again, stage1 is a farm network of CT moore processes.

```
stage1CT :: V.Vector (TimeStamp, Time -> AbstExt Time) - - vector of initial states
    ->V.Vector (CT.Signal (AbstExt Time)) - - vector of input signals
    | V.Vector (CT.Signal (AbstExt Time)) - - vector of output signals
stage1CT = V.farm21 pcCT
    where
        pcCT = CT.moore11 ns od
        ns = ExB.res21 (+)
        od = ExB.res11 id
```

We cannot print ict nor vct any more, but we can plot them. Again, the generated plots might need to be tweaked.

```
\lambda> let latexIn = latexV 3.5 ["ct1","ct2","ct3","ct4"] vct
\lambda> let latexS1 = latexV 3.5 ["ct1-1","ct2-1","ct3-1","ct4-1"] $ stage1CT ict vct
\lambda> let gnuIn = plotV 3.5 ["ct1","ct2","ct3","ct4"] vct
\lambda> let gnuS1 = plotV 3.5 ["ct1-1","ct2-1","ct3-1","ct4-1"] $ stage1CT ict vct
```


(a) latexIn

(c) gnuIn

(b) latexS1

(d) gnuS1

The second stage is a reduce network of CT comb processes.

```
stage2CT :: V.Vector (CT.Signal (AbstExt Time))
    CT.Signal (AbstExt Time)
stage2CT = V.reduce rCT
    where
        rCT = CT.comb21 (ExB.res21 (+))
```

```
\lambda> let s2out = (stage2CT . stage1CT ict) vct
\lambda> let latexS2 = latex 3.3 ["ct1-2","ct2-2","ct3-2","ct4-2"] s2out
\lambda> let gnuS2 = plot 3.3 ["ct1-2","ct2-2","ct3-2","ct4-2"] s2out
```

For the last stage of the toy system there is no CT process constructor in ForSyDe.Atom.MoC.CT so we need to create it ourselves.

(a) latexS2

(b) gnuS2

```
stage3CT :: CT.Signal (AbstExt Time)
    ->CT.Signal (AbstExt Time)
stage3CT = ctFilter (>=0)
    where
        ctFilter p s = CT.comb21 ExB.filter (predSig p s) s
        predSig p s = CT.comb11 (ExB.res11 p) s
toyCT :: V.Vector (TimeStamp, Time -> AbstExt Time) - ~ initial states
        ->V.Vector (CT.Signal (AbstExt Time)) - - input
        CT.Signal (AbstExt Time) - - output
toyCT i = stage3CT . stage2CT . stage1CT i
```

| $\lambda>$ let latexS3 $=$ latex 3.3 ["ct1-3"] \$ toyCT ict vct |
| :--- |
| $\lambda>$ let gnuS3 $=$ plot $3.3[" c t 1-3 "] \$$ toyCT ict vct |

## ct1-3


(a) latexS3

(b) gnuS3

### 2.3.5 SDF instance

The SDF instance of the toy system is created using process constructor helpers defined in ForSdfDe.Atom.MoC. SDF, and can be found in the following module (re-exported by AtomExamples.GettingStarted).
|module AtomExamples.GettingStarted.SDF where

```
import ForSyDe.Atom
import ForSyDe.Atom.ExB.Absent (AbstExt)
import ForSyDe.Atom.ExB as ExB
import ForSyDe.Atom.MoC.SDF as SDF
import ForSyDe.Atom.Skeleton.Vector as V
```

stage1 is defined as a farm network of SDF moore processes. As SDF Moore processes are in principle graph loops, we take the initial tokens for each loop from a vector of lists of tokens. Also, both next state and output decoder functions are defined over lists of values instead of values, and they need to be provided within a context which describes the production and consumption rates. Read more about the particularities of SDF in the API documentation.

```
stage1SDF :: V.Vector ([AbstExt Int]) - vector of initial tokens
    -> V.Vector (SDF.Signal (AbstExt Int)) - ~ vector of input signals
```

```
     V.Vector (SDF.Signal (AbstExt Int)) - ~ vector of output signals
stage1SDF = V.farm21 pcSDF
    where
        pcSDF = SDF.moore11 ((1,2),1,ns) (1,1,od)
        ns [x1] [y1,y2] = [ExB.res31 (\lambdaa b c -> a + b + c) x1 y1 y2]
        od [x1] = [ExB.res11 id x1]
```

Let us print and plot the inputs against the outputs, using the test signals and plotting functions latexV and plotV defined in section 2.3.1:

```
\lambda> isdf
<[-1], [1], [-1], [1]>
\lambda> vsdf
<{1,1,1,1,1,1},{-1,1,-1,1,-1,1},{0,0,1,1,0},{-1,-1,-1,-1,-1}>
\lambda> stage1SDF isdf vsdf
<{-1,1,3,5},{1,1,1,1},{-1,-1,1},{1,-1,-3}>
\lambda> let latexIn = latexV 6 ["sdf1","sdf2","sdf3","sdf4"] vsdf
\lambda> let latexS1 = latexV 7 ["sdf1-1","sdf2-1","sdf3-1","sdf4-1"] $ stage1SDF isdf vsdf
\lambda> let gnuIn = plotV 6 ["sdf1","sdf2","sdf3","sdf4"] vsdf
\lambda> let gnuS1 = plotV 7 ["sdf1-1","sdf2-1","sdf3-1","sdf4-1"] $ stage1SDF isdf vsdf
```

| 1.0 | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| -1.0 | 1.0 | -1.0 | 1.0 | -1.0 | 1.0 |
| 0.0 | 0.0 | 1.0 | 1.0 | 0.0 |  |
| -1.0 | -1.0 | -1.0 | -1.0 | -1.0 |  |


(b) latexS1
(a) latexIn

(c) gnuIn

(d) gnuS1
stage2 is again a reduce network of comb processes. As with stage1, we need to provide the production and consumption rates.

```
stage2SDF :: V.Vector (SDF.Signal (AbstExt Int))
    -> SDF.Signal (AbstExt Int)
stage2SDF = V.reduce rSDF
    where
        rSDF = SDF.comb21 ((1,1),1,rF)
        rF [x1] [y1] = [ExB.res21 (+) x1 y1]
```

Again, let us print and plot the output signals using the test inputs and utilities defined in section 2.3.1.

```
\lambda> let s2out = (stage2SDF . stage1SDF isdf) vsdf
\lambda> s2out
{0,0,2}
\lambda> let latexS2 = latex 3 ["sdf1-2"] s2out
\lambda> let gnuS2 = plot 3 ["sdf1-2"] s2out
```

As for DE and CT instances, a SDF filter process does not really make sense in practice, but for the scope of this toy system, we need to instantiate one ourselves.

```
stage3SDF :: SDF.Signal (AbstExt Int)
    -> SDF.Signal (AbstExt Int)
stage3SDF = sdfFilter (>=0)
    where
```

$$
\begin{array}{lll}
0.0 & 0.0 & 2.0
\end{array}
$$

(a) latexS2

(b) gnuS2

```
sdfFilter p s = SDF.comb21 ((2,2),2,filterF) (predSig p s) s
filterF pl sl = zipWith ExB.filter pl sl
predSig p s = SDF.comb11 (1,1,fmap (ExB.res11 p)) s
```

```
|
    ->V.Vector (SDF.Signal (AbstExt Int)) - - input
    SDF.Signal (AbstExt Int) - - output
toySDF i = stage3SDF . stage2SDF . stage1SDF i
```


## $0.0 \quad 0.0$

(a) latexS3

(b) gnuS3

### 2.3.6 Polymorphic instance

In the previous section you've seen how to model systems in ForSyDe-Atom using the helper functions for instantiating process constructors in different MoCs. In this section we will be instantiating the "raw" polymorphic form of the same process constructors, not overloaded with any execution semantics. The execution semantics are deduced from the tag system of the input signals, i.e. their types. These process constructors are defined as patterns of MoC atoms in the ForSdfDe.Atom.MoC module. The code below is exported by AtomExamples.GettingStarted.

## |module AtomExamples.GettingStarted.Polymorphic where

Notice that apart from the polymorphic MoC patterns, we are also using "raw" extended behavior and skeleton patterns.

```
import ForSyDe.Atom
import ForSyDe.Atom.MoC.SDF (Prod, Cons)
import ForSyDe.Atom.ExB as ExB
import ForSyDe.Atom.MoC as MoC
import ForSyDe.Atom.Skeleton as Skel
```

stage1 is defined, like in all previous instances, as a farm network of moore processes.

```
stage1 :: (Skeleton s, MoC m, ExB b)
    => Fun m (b a) (Fun m (b a) (Ret m (b a))) - - next state function
    Fun m (b a) (Ret m (b a)) - output decoder function
    ->s(Stream (m (b a))) - signals with initial tokens
    -> s (Stream (m (b a))) - ~ vector of input signals
    ->s (Stream (m (b a))) - vector of output signals
stage1 ns od = Skel.farm21 (MoC.moore11 ns od)
```

We can immediately observe some main differences in the type signature. First, the Vector, Signal and AbstExt data types are not explicit any more, but suggested as type constraints. The first line in the type signature (Skel s, MoC m, ExB b) suggests that the type of should belong to the skeleton layer, the type of $m$ should belong to the MoC layer and the type of b should belong to the extended behavior layer. Another peculiarity is the presence of the first two structures, but there should be nothing frightening about them: e.g. a structure Fun ma(Fun mbet $m \mathrm{c}$ ) ) simply stands for the type of a function a -> b $->\mathrm{c}$, which was wrapped in a context specific to a MoC m. Read the MoC layer's API documentation for more on function contexts. This means that the functions for the next state decoder and the output decoder need to be provided as arguments for stage1, and they might differ depending on the MoCs. A third peculiarity is that the initial states are provided as signals and not through some specific structure any more. Indeed, the MoC atoms extract initial states from signals, and deal with them in different ways depending on the MoC they implement.

The main two classes of MoCs, based on their notion of tags, but also based on how they deal with events, are timed MoCs (e.g. SY, DE, CT) and untimed MoCs (e.g. SDF). Concerning the functions they lift from layers below, we can say that in ForSyDe-Atom timed MoCs lift functions on individual values, whereas untimed MoCs lift functions on lists of values (i.e. multiple tokens). Based on this observation, let us define the next state and output decoders for timed and untimed/SDF MoCs.

```
nsT :: (ExB b, Num a) => b a }->\textrm{b}a\textrm{a}->\textrm{b} \textrm{a
odT :: (ExB b, Num a) => b a }->\textrm{b}\mathrm{ a
nsT = ExB.res21 (+)
odT = ExB.res11 id
nsSDF :: (ExB b, Num a) => (Cons, [b a] -> (Cons, [b a] -> (Prod, [b a])))
odSDF :: (ExB b, Num a) => (Cons, [b a] -> (Prod, [b a]))
nsSDF = MoC.ctxt21 (1,2) 1 ( }\lambda[\textrm{x}1] [y1,y2] -> [ExB.res31 (\lambdaa b c -> a + b + c) x1 y1 y2]) 
odSDF = MoC.ctxt11 1 1 ( }\lambda[\textrm{x}1]\quad->\mathrm{ [ExB.res11 id x1])
```

OBS: for the sake of simplicity, the ExB component has been left as part of the nsT and odT, respectively nsSDF and odSDF, and not part of stage1. Describing all layers within the stage1 function would have rendered the type signature a bit more complicated and is left as an exercise for the reader.

We postpone plotting the input and output signals for later. Carrying on with instatiating stage 2 as a reduce network of comb processes:

```
stage2 :: (Skeleton s, MoC m, ExB b)
    => Fun m (b a) (Fun m (b a) (Ret m (b a))) - - reduce function
    s (Stream (m (b a))) - - vector of input signals
    Stream (m (b a)) - - output signal
stage2 r = Skel.reduce (MoC.comb21 r)
```

Again, the passed functions need to be specifically defined for each MoC, and for simplicity we include the ExB part as well:

```
rT :: (ExB b,Num a) => b a }->\textrm{b}a->\textrm{a}|\textrm{a
rT = ExB.res21 (+)
rSDF :: (ExB b, Num a) => (Cons, [b a] -> (Cons, [b a] -> (Prod, [b a])))
rSDF = MoC.ctxt21 (1,1) 1 ( }\lambda[\textrm{x}1] [y1] -> [ExB.res21 (+) x1 y1]),
```

Finally stage3, the filter pattern, we create it ourselves in terms of existing ones. This time we incorporate the extended behaviors in the definition of stage3, an we only ask for a context wrapper as input argument. Don't be alarmed by the scary type signature, the actual implementation is quite elegant.

```
stage3 :: (MoC m, ExB b, Ord a, Num a)
    =((b a }->\textrm{b}a
        Fun m (b a) (Ret m (b a))) - context wrapper for the filter behavior
    -> Stream (m (b a)) - - input signal
    ->Stream (m (b a)) - - output signal
stage3 fctx = MoC.comb11 (fctx filtF)
    where filtF a = ExB.filter (ExB.res11 (>=0) a) a
```

And now for the timed/untimed context wrappers:

```
fctxT = id
fctxSDF f = MoC.ctxt11 2 2 (fmap f)
```

The full definition of the toy system:

```
|toy ns od r fctx is = stage3 fctx . stage2 r . stage1 ns od is
```

And that's it! Let us plot now the test signals and the responses of the system for each stage. This time we will use only $\mathrm{ET}_{\mathrm{A}} \mathrm{X}$ plots. We can also plot initial states as they are wrapped as signals. The test results can be seen in fig. 2.7.

```
\lambda> let noLabel = ["","","",""]
\lambda> let iSDF = latexV 2 noLabel sisdf
\lambda> let iSY = latexV 2 noLabel sisy
\lambda> let iDE = latexV 2 noLabel side
\lambda> let iCT = latexV 2 noLabel sict
\lambda>
\lambda> let vSDF = latexV 6 noLabel vsdf
\lambda> let vSY = latexV 6 noLabel vsy
\lambda> let vDE = latexV 3.3 noLabel vde
\lambda> let vCT = latexV 3.3 noLabel vct
\lambda>
\lambda> let s1SDF = latexV 6 noLabel $ stage1 nsSDF odSDF sisdf vsdf
\lambda> let s1SY = latexV 6 noLabel $ stage1 nsT odT sisy vsy
\lambda> let s1DE = latexV 3.3 noLabel $ stage1 nsT odT side vde
\lambda> let s1CT = latexV 3.3 noLabel $ stage1 nsT odT sict vct
\lambda> let s2 ns od r is = stage2 r . stage1 ns od is
\lambda> let s2SDF = latex 6 noLabel $ s2 nsSDF odSDF rSDF sisdf vsdf
\lambda> let s2SY = latex 6 noLabel $ s2 nsT odT rT sisy vsy
\lambda> let s2DE = latex 3.3 noLabel $ s2 nsT odT rT side vde
\lambda> let s2CT = latex 3.3 noLabel $ s2 nsT odT rT sict vct
\lambda> let s3SDF = latex 6 noLabel $ toy nsSDF odSDF rSDF fctxSDF sisdf vsdf
\lambda> let s3SY = latex 6 noLabel $ toy nsT odT rT fctxT sisy vsy
\lambda> let s3DE = latex 3.3 noLabel $ toy nsT odT rT fctxT side vde
\lambda> let s3CT = latex 3.3 noLabel $ toy nsT odT rT fctxT sict vct
```

As expected, the results in fig. 2.7 are exactly the same as the ones presented in section 2.3.2 and ??????. In conclusion we have succesfully instantiated a MoC-agnostic system, whose execution semantics are inferred according to the input data types. This is possible thanks to the notion of type classes, inferred from the host language Haskell. In this section, instead of MoC-specific helpers, we have used the "raw" process constructors as defined in the ForSdfDe. Atom. MoC module as patterns of MoC-layer atoms.

This example, used as a case study by Ungureanu and Sander, 2017, has been focused on the MoC layer. A similar approach based on atom polymorphism could target other layers as well since, as you have seen, all layers are implemented as type classes. At the moment of writing this report the extended behavior layer was represented only by the AbstExt type, while the skeleton layer had only Vector. Nevertheless, future iterations of ForSyDe-Atom will describe more types.

### 2.4 Making your own patterns

The final section of this report introduces the reader to constructing custom patterns in ForSyDeAtom. Up until now we have been using patterns which were pre-defined as compositions of atoms. Atoms are primitive, indivizible building blocks capturing the most basic semantics in each layer.
| \{-\# LANGUAGE PostfixOperators \#-\}
The code for this section is found in the following module, which is not re-exported, i.e. needs to be manually imported.
| module AtomExamples. GettingStarted. CustomPattern where
For this exercise, we will create a custom comb pattern with 5 inputs and 3 outputs, as a process constructor in the MoC layer. We will test this pattern with a set of SY and a set of DE input signals, thus we need to import the following modules:




$$
\begin{aligned}
& \text { - } \\
& \begin{array}{c}
\text { yaszs ( } \mathrm{p} \text { ) } \\
0 \mathrm{Z} \\
0.0 \\
0.0
\end{array}
\end{aligned}
$$

```
import ForSyDe.Atom
import ForSyDe.Atom.MoC
import ForSyDe.Atom.MoC.SY as SY
import ForSyDe.Atom.MoC.DE as DE
```

The best way to start building your own patterns is to study the source code for the existing patterns and see how they are made. If you don't want to dig into the source code of ForSyDeAtom, there is a link in the API documentation for each exported element, as suggested in fig. 2.8.


Figure 2.8: Screenshot from the API documentation. The link to the source code is marked with a red rectangle

Studying the comb22 pattern, you can see that it is defined in terms of the lift and sync atoms which are represented by the infix operators -.- and $-*$ - respectively, and the unzip utility represented by the postfix operator $-*<$. lift and sync are atoms becuse they capture an interface for exeucution semantics, whereas unzip is just a utility because it is merely a type traversal which alters the structure of data types and rebuilds it to describe "signals of events carrying values".

Considering the applicative nature of the -.- and -*- atoms, the comb pattern with 5 inputs and 3 outputs can be written as the mathematical formula below. This one-liner tells that function f is "lifted" into the MoC domain, and applied to the five input signals which are synchronized. The $-*$ « postfix operator then "unzips" the resulting signal of triples into three synchronized signals of values. The applicative mechanism explained in the previous paragraph is depicted in fig. 2.9.

$$
\begin{aligned}
& \text { comb53 f s1 s2 s3 s4 s5 } \\
& \quad=(\mathrm{f}-.-\mathrm{s} 1-*-\mathrm{s} 2-*-\mathrm{s} 3-*-\mathrm{s} 4-*-\text { s5 }-* \ll)
\end{aligned}
$$



Figure 2.9: Composition of atoms forming the comb pattern
If we want to restrict the pattern to one specific MoC, then we must mention this in the type signature we associate it with, like in the example below.

```
comb53SY :: (a1 -> a2 }->\textrm{a}3->\textrm{a}4->\textrm{a}5->(\textrm{b}1,\textrm{b}2,\textrm{b}3)
    SY.Signal a1 - - input signal
    SY.Signal a2 - - input signal
    SY.Signal a3 - - input signal
    -> SY.Signal a4 - - input signal
    SY.Signal a5 - - input signal
    ( SY.Signal b1
        , SY.Signal b2
        , SY.Signal b3) - - 3 output signals
comb53SY f s1 s2 s3 s4 s5
=(f -.- s1 -*- s2 -*- s3 -*- s4 -*- s5 -*<<)
```

To test the output, let us create five signals and a function that needs to be lifted. For the example, the terminal printouts should suffice to test our simple pattern.

```
\lambda> import AtomExamples.GettingStarted.CustomPattern as CP
\lambda> let fun a b c d e = (a+c+e, d-b, a*e)
\lambda> let sy1 = SY.signal [1,2,3,4,5]
\lambda> let sy2 = SY.comb11 (+10) sy1
\lambda> let sy3 = SY.constant1 100
\lambda> let de1 = DE.signal [(0,1),(3,2),(7,3),(9,4),(11,5)]
\lambda> let de2 = DE.signal [ (0,11), (3,12), (5,13), (9,14), (11,15)]
\lambda> let de4 = DE.constant1 100
\lambda>
\lambda> let (o1,o2,o3) = CP.comb53 fun sy1 sy1 sy2 sy2 sy3
\lambda> o1
\lambda> {112,114,116,118,120}
\lambda> o2
\lambda> {10, 10, 10, 10, 10}
\lambda> o3
\lambda> {100, 200, 300,400,500}
\lambda>
\lambda> let (o1,o2,o3) = CP.comb53 fun de1 de1 de2 de2 de4
\lambda> o1
\lambda> { 112 @0s, 114 @3s, 115 @5s, 116 @7s, 118 @9s, 120 @11s}
\lambda> o2
\lambda> { 10 @0s, 10 @3s, 11 @5s, 10 @7s, 10 @9s, 10 @11s}
\lambda> o3
\lambda> { 100 @0s, 200 @3s, 200 @5s, 300 @7s, 400 @9s, 500 @11s}
\lambda>
\lambda> let (o1,o2,o3) = CP.comb53SY fun sy1 sy1 sy2 sy2 sy3
\lambda> o1
\lambda> {112,114,116,118,120}
\lambda> o2
\lambda> {10, 10, 10, 10, 10}
\lambda> o3
\lambda> {100, 200, 300,400,500}
\lambda>
\lambda> let (o1,o2,o3) = CP.comb53SY fun de1 de1 de2 de2 de4
<interactive>:71:56-58:
    Couldn't match type 'DE Integer' with 'SY b3'
    Expected type: SY.Signal b3
        Actual type: DE.Signal Integer
    Relevant bindings include
        it :: (SY.Signal b3, SY.Signal b2, SY.Signal b3)
            (bound at <interactive>:71:1)
    In the second argument of 'c0omb53SY', namely 'de1'
    In the expression: comb53SY fun de1 de1 de2 de2 de4
...
```


### 2.5 Conclusion

This report has introduced the reader to the basic features of ForSyDe-Atom a framework for modeling and testing of cyber-physical systems. It has covered basic usage such as instantiating systems and plotting signals. It briefly went through concepts such as layers, atoms and patterns, and has focused on their practical usage. A step-by-step tutorial has been presented, showing alternative ways to instantiate systems and demonstrating the polymorphism of layers.

The reader is recommended to further consult the API documentation which also acts as a manual for the library, as well as the related publications listed on the project web site. Future reports will assume familiarity with using and understanding the framework and will focus mainly on results.

### 2.6 References

Halbwachs, Nicholas et al. (1991). "The synchronous data flow programming language LUSTRE". In: Proceedings of the IEEE 79.9, pp. 1305-1320.

Ungureanu, George and Ingo Sander (2017). "A layered formal framework for modeling of cyberphysical systems". In: 2017 Design, Automation $\mathcal{B}$ Test in Europe Conference $\mathcal{E}$ Exhibition (DATE). IEEE, pp. 1715-1720.

## Hybrid CT/DT Models in ForSyDE-ATom

This chapter gathers examples and experiments which involve hybrid models focusing on the the semantics and implications of combining and translating between continuous and discrete domains. As such, the focus is mainly on exploring alternative models and analyzing the implications on fidelity, precision and performance. This chapter is also meant to support the associated scientific publications with experimental results.

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$$
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Info
Compatibility : FORSYDE-ATOM version 0.2.2
Repository : https://github.com/forsyde/forsyde-atom-examples
Path : <repo_root>/hybrid
Other deps. : none

### 3.1 Goals

The reader is assumed to have been familiarized with the ForSyDe-Atom modeling framework. A good resource for that is chapter 2. The main goals of this chapter are:

- explore alternative ForSyDe-Atom models for widely known hybrid (discrete and continuous time) systems, with the scope of analyzing the implications from the modeling and simulation perspective.
- train the reader into the decision-making process of modeling hybrid systems, and the tradeoffs involved. While as per writing this report the ForSyDe-Atom modeling framework is still limited, future directions are hinted.
- support the associated scientific publications with the complete experiments and results for the models used as case studies. These publications include: ungureanu18a


### 3.2 RC Oscillator

In this section we model the RC oscillator setup in fig. 3.1. Despite its apparent simplicity, shows an interesting problem in CPS: how continuous systems react to discrete stimuli. This example was used in Ungureanu, Medeiros, and Sander, 2018, and for a study on the theoretical implications
of the chosen models, we strongly recommend reading this paper before going further. The source code for this section is found in the following module:
|module AtomExamples.Hybrid.RCOsc where


Figure 3.1: RC oscillator setup
These are the dependencies that need to be included within this module:

```
import ForSyDe.Atom
import ForSyDe.Atom.MoC.CT
import ForSyDe.Atom.MoC.DE as DE - DE MoC library
import ForSyDe.Atom.MoC.SY as SY - SY MoC library
import ForSyDe.Atom.MoC.Time as T - utilities for functions of time
import ForSyDe.Atom.MoC.TimeStamp (milisec) - utilities for time stamps
import ForSyDe.Atom.Utility.Plot - plotting utilities
```

The circuit in fig. 3.1 is characterized by two states: 1) $s w_{1}$ is closed and $s w_{2}$ open, equivalent to the situation where the capacitor C charges with VDD; 2) $s w_{1}$ is open and $s w_{2}$ is closed, equivalent to the situation where the capacitor C discharges to the ground through R . For the scope of this example, to keep the model simple and deterministic, we assume that $s w_{1}$ and $s w_{2}$ open/close alternatively at the same discrete time instants, there are no intermediary transitions, and we ignore the effects of discharge through switches.

Let us analyze case 1) above. The circuit in fig. 3.1 becomes a RC integrator, where the input signal is applied to the resistance with the output taken across the capacitor. The amount of charge that is established across the plates of the capacitor is equal to the time domain integral of the current, like in eq. (3.1). The rate at which the capacitor charges is directly proportional to the amount of the resistance and capacitance giving the time constant of the circuit like in eq. (3.2). As the capacitors current can be expressed as the rate of change of charge, $Q$ with respect to time, we can express the charge at any instant of time like in eq. (3.3). Eq. (3.3) brings together the previous equations and uses the instant voltage charge formula to obtain the final integral formula for $V_{O}$.

$$
\begin{gather*}
i_{C}(t)=C \frac{\partial V_{C(t)}}{\partial t}=\frac{V_{D D}}{R}  \tag{3.1}\\
R C=R \frac{Q}{V}=R \frac{i \times T}{i \times R}=T  \tag{3.2}\\
i=\frac{\partial Q}{\partial t} \Rightarrow Q=\int i \partial t  \tag{3.3}\\
V_{O}=V_{C}=\frac{Q}{C} \stackrel{(3)}{=} \frac{1}{C} \int i \partial t \stackrel{(1)}{=} \frac{1}{C} \int \frac{V_{D D}}{R} \partial t=\frac{1}{R C} \int V_{D D} \partial t \tag{3.4}
\end{gather*}
$$

If an ideal step voltage pulse is applied, that is with the leading edge and trailing edge considered as being instantaneous, the voltage across the capacitor will increase for charging exponentially over time at a rate determined by eq. (3.5). Similarly, in case 2) the circuit becomes an RC bridge where, assuming that C has been fully charged, it now discharges to the ground at a rate
determined by eq. (3.6).

$$
\begin{gather*}
V_{O, \text { charge }}(t)=V_{C}(t)=V_{D D}\left(1-e^{-\frac{t}{R C}}\right)  \tag{3.5}\\
V_{O, \text { discharge }}(t)=V_{C}(t)=V_{D D}\left(e^{-\frac{t}{R C}}\right) \tag{3.6}
\end{gather*}
$$

Translating eqs. (3.5) and (3.6) into ForSyDe-Atom Haskell code, we get the following, assuming to is the moment where the switch occurred, and ignoring the factor of $V_{D D}$ (which is considered and scaled in the output decoder anyway).

```
- / RC time constant.
rc = 0.1
- I V_O(tO,t) during the discharging. to is the instant when the switch occurred.
vcDischarge to = \lambdat }->\mathrm{ T.exp (-(t-t0) / rc)
- I V_O(tO,t) during the charging. to is the instant when the switch occurred.
vcCharge to = \lambdat -> 1 - T.exp (-(t-t0) / rc)
```

Naturally, in ForSyDe-Atom we can model the oscillator in multiple ways, depending on what aspect we want to focus. Our first example models the RC oscillator as a Mealy finite state machine which embeds the continuous and discrete time semantics as defined in the ForSyCt. Atom. MoC.CT library, like in fig. 3.2. As such, we imply from the model that the switching of $s w_{1}$ and $s w_{2}$ is performed periodically after e certain $\tau$, inferred from the initial state of the CT.mealy process. As such, the state machine is loaded with the voltage charging rule in eq. (3.5), and a duration $\tau$. The configuration of the mealy pattern ensures that at each multiple of $\tau$ the next state function ns will be performed, negating the state rule (i.e. the rate for $V_{O}$ ), basically translatig back and forth between eqs. (3.5) and (3.6). As mentioned before, the output decoder od merely rescales $V_{O}$ with $V_{D D}$.


Figure 3.2: Internal pattern of the initial RC oscillator setup

```
|_ I RC oscillator model as CT FSM with discrete semantics inherent in the CT model.
osc1 :: CT.Signal Rational - VDD as input signal
    CT.Signal Rational - - V_O as output signal
osc1 = CT.mealy11 ns od (milisec 500, vcCharge 0)
    where
        ns v _ = 1 + (-1 * v)
        od = (*)
```

Plotting the output against an "arbitrary" $V_{D D}$, we get fig. 3.3.

```
- I plotting configuration that will be used throughout this section
```

$\mathrm{cfg}=$ defaultCfg $\{\mathrm{xmax}=3$, rate=0.01\}

```
-_ I example input for testing 'osc1'
vdd1 = CT.signal [(0,\mp@subsup{\lambda}{-}{}->2),(1,\mp@subsup{\lambda}{-}{}->1.5),(2,\mp@subsup{\lambda}{-}{}->1)] :: CT.Signal Rational
- I plotting the response of 'osc1'
plot1 = plotGnu $ prepareL cfg {labels=["V_{DD}","V_0"]} $ [vdd1, osc1 vdd1]
```

The model for osc1, although strikes out as simple and elegant, is by all means limited. In the following paragraphs we will address these limitations one at a time, and try find solutions that overcome them in order to include more realistic or at least a family of behaviors correctly covering classes of (increasing) non-determinism.

Lee, 2016 argues that determinism is an attribute of the model, and it is dependent on what we consider as inputs and outputs. As such, the model osc1 in fig. 3.2 is a deterministic model for the circuit in fig. 3.1 and the output in fig. 3.3 is the correct response in the following conditions:


Figure 3.3: Response of osc1

- the changes in $V_{D D}$ happen at multiples of $\tau$. If a changes occur at any other time the response will be shown, but it will not describe the RC circuit in a realistic manner, as we will see shortly.
- the time constant follows roughly the rule $5 R C \leq \tau$ where the chosen $\tau$ is a part of the initial state of the Mealy state machine. This says that the discharging occurrs when the capacitor is charged at least $99.3 \%$ and vice-versa, according to eqs. (3.5) and (3.6).
- $\tau>0 . \tau=0$ would render the system as non-causal and would manifest Zeno behavior. In practical terms, this is the equivalent of the simulation being stuck and not advancing time.

Another property (or limitation, depending on what your intentions are) is that the switching of $s w_{1}$ and $s w_{2}$ cannot be controlled and is a property of the Mealy machine. This switching is inferred by $\tau$ which sets the (discrete) period of these events. Let us change that by giving the possibility to control the state of the capacitor as charging/discharging through a signal of discrete events. A system such as the one we propose implies some subtle changes in the modelling which require a deeper understanding on the underlying MoCs. First of all, we still require a stateful process (a state machine) which captures the notion of working modes, but which reacts to a state change instantaneously. But this implies that $\tau=0$ which, as said above, would lead to Zeno behavior. On the other hand, the particular behavior required is described precisely by synchronous reactive MoC, where the response of a system is performed in "zero time".

Continuing to adhere to the school of thought of Lee, 2016, where the merit of the modeler lies in choosing the right modeling paradigm for the right problem ${ }^{1}$, we choose to model the above described system as a synchronous reactive (SY) state machine wrapped inside a DE/CT environment like in fig. 3.4. As such, we should note a few particularities of this model:

- in CT the carried values are implicit functions of time, i.e. they carry time semantics. In DE and SY, the time semantics make no sense, thus the same functions of time are explicit (where Time itself is just a data type $\in V$ ). Therefore when converting $\mathcal{S}_{C T}(\alpha) \mapsto \mathcal{S}_{D E}(t \rightarrow \alpha)$ and vice-versa $\mathcal{S}_{D E}(t \rightarrow \alpha) \mapsto \mathcal{S}_{C T}(\alpha)$.

[^3]- in SY tags are useless, therefore DE tags and values are split into two separate SY signals upon conversion. This way we can easily recreate the DE signal without loss of information. So upon conversion $\mathcal{S}_{D E}(\alpha) \mapsto \mathcal{S}_{S Y}(t) \times \mathcal{S}_{S Y}(\alpha)$ and vice-versa $\mathcal{S}_{S Y}(t) \times \mathcal{S}_{S Y}(\alpha) \mapsto \mathcal{S}_{D E}(\alpha)$.
- the discrete control signal $S_{\text {control }}$, carries nothing, and it is used only for the timestamps generated by its tags, i.e. for generating $\mathcal{S}_{S Y}(t)$. These timestamps are further used by the od function to determine the time when the switch occurred to in the formulas for $V_{O}$ in eqs. (3.5) and (3.6).


Figure 3.4: RC oscillator model which reacts to a discrete control signal. Shapes decorating signals suggest the type of carried tokens: circles $=$ scalars; squares $=$ functions

Like in the model from fig. 3.2, the model for osc2 in fig. 3.4 scales the output with $V_{D D}$ through a combinational CT process. The next state decoder function $n s$ does not manipulate the rule for $V_{C}$ anymore, but rather switches between the states Charge and Discharge, and the output decoder od selects the proper $V_{C}$ rule from eq. (3.5) or eq. (3.6) respectively, based on the current state.

```
- I Encodes the capacitor state
data CState = Charging | Discharging
```

We encode the capcitor state (i.e. the states of $s w_{1}$ and $s w_{2}$ ) through the CState data type. Thus the code for fig. 3.4 is:

```
- I RC oscillator model as FSM which reacts to discrete impulses
osc2 :: CT.Signal Rational - - VDD input signal
    -> DE.Signal () - control signal of discrete impulses
    CT.Signal Rational - - output signal
osc2 s = CT.comb21 (*) s . embed (SY.mealy11 ns od Charging)
    where
        - SY next state function
        ns Charging _ = Discharging
        ns Discharging _ = Charging
        - SY output decoder function
        od Charging to = vcCharge (time t0)
        od Discharging tO = vcDischarge (time t0)
        - wrapper that embeds a SY process into a mixed DE/CT environment
        embed p de = let (t, _) = DE.toSY de
                in DE.toCT $ SY.toDE t (p t)
```

Now let us create two situations to test osc2. The control signal sCtrl injects events at timestamps $0,0.6,1.4,2.3$ and 2.8 , causing the SY state machine to react each time changing its state from Charge to Discharge and vice-versa. The first $V_{D D}$ input vdd21 is mirroring an "ideal case", when changes occur at time instants where switching happens, meaning that it does not affect the evolution of the $V_{O}$ rule. On the contrary, vdd22 comes at arbitrary times, thus affecting the output in a way that is unrealistic, i.e. $V_{O}$ changes values abruptly and instantaneously, which strays away from the acceptable behavior of a capacitor. This can be seen in fig. 3.5.

```
- I Signal of discrete events. Carries nothing, it is used only for the time stamps.
sCtrl = DE.signal [(0,()),(0.6,()),(1.4,()),(2.3,()),(2.8,())] :: DE.Signal ()
- I example input for testing 'osc2'
vdd21 = CT.signal [(0,\lambda_ ->2),(1.4, \mp@subsup{\lambda}{-}{}->1.5),(2.8,\mp@subsup{\lambda}{-}{}->1)] :: CT.Signal Rational
vdd22 = CT.signal [(0,\lambda_->1),(0.8,\mp@subsup{\lambda}{-}{}->2),(1.6,\mp@subsup{\lambda}{-}{}->1.5)] :: CT.Signal Rational
- | plotting the example responses of 'osc2'
plot21 = plotGnu $ prepareL cfg {labels=["V_{DD}","V_O"]} $ [vdd21, osc2 vdd21 sCtrl]
plot22 = plotGnu $ prepareL cfg {labels=["V_{DD}","V_O"]} $ [vdd22, osc2 vdd22 sCtrl]
```



Figure 3.5: Response of osc2

The second case in fig. 3.5 can be explained easily if we consider the fact that $V_{O}$ itself is the result of a statically pre-calculated function of time. There is no notion of feedback response to the continuous inputs, the only feedback is synchronous reactive for osc2 and even osc1. In other words any change on the input will affect how the output "looks like", but it will not affect the continuous behavior which, as said, is pre-calculated. In order to influence the continuous behavior, some sort of continuous feedback is necessary, which is usually non-causal, thus uncomputable. Without going too much into detail we can say that in order to model a realistic behavior of the capacitor response in $V_{O}$, we need to embed an ordinary differential equation (ODE) solver within a synchronous reactive state machine ${ }^{2}$, which models precisely eq. (3.4).


Figure 3.6: RC bridge: circuit (left); block diagram based on eq. (3.4) (right)

$$
\begin{equation*}
\int_{t_{0}}^{t_{0}+h} f(t, y(t)) \partial t \approx h f\left(t_{0}, y\left(t_{0}\right)\right) \tag{3.7}
\end{equation*}
$$

$$
\begin{equation*}
\dot{y}_{n}=\frac{h R C}{h+R C}\left(\frac{1}{R C} y_{n}+\frac{1}{h} y_{n-1}\right) \tag{3.8}
\end{equation*}
$$

First let us model a simple RC bridge like in fig. 3.6 which acts like a low-pass filter, and acts according to eq. (3.4). As mentioned, we model this circuit as a SY state machine embedded within a CT environment. Why a state machine? Because the circuit itself, due to the capacitor, is a memory system, i.e. its output is dependent on its history as well as inputs. The previous "history" at the start of simulation is encoded inside the initial state. The next state decoder itself needs to "feed-through" the input and mirror the behavior of the block diagram fig. 3.6. On the other hand, this block diagram shows a non-causal system, which as such is uncomputable, and needs to be transformed into a solvable form. As per the writing of this report, ForSyDe-Atom ${ }^{3}$ did not provide generic ODE solvers, thus for didactic purpose we write the following numerical solver for our RC circuit in fig. 3.6 using Euler's trapezoidal method eq. (3.7), by substituting by hand eq. (3.4) with its feed-forward version in eq. (3.8):

```
- I ODE solver for a simple RC low-pass filter using Euler's method
euler :: TimeStamp _ time step for the solver precision
    (Time -> Rational) - the input function being integrated
```

[^4]```
    Rational - the "history" of the integral at to
    TimeStamp - - to
    \rightarrow \text { Time } \rightarrow \text { Rational - - a function of time}
euler step f p to t = iterate p t0
    where
        h = time step
        - by-hand substitution of eq. (4) using the trapezoidal rule
        calc vp v = (h * rc)/(h + rc) * (1/rc * v + 1/h * vp)
        _ loop which calculates the integral step-wise from to to t
        iterate st ti
            | t < time ti = st
            | otherwise = iterate (calc st $ f (time ti)) (ti + step)
```

As said earlier, we model the RC circuit in fig. 3.6 by including the euler solver into a feedthrough state machine (see the state pattern) like in fig. 3.7. Like for osc2, the timestamps resulted after splitting the tags from values when interfacing between DE and SY are used for determining the $t_{0} \mathrm{~s}$ for each new incoming event. The solver inputs functions of time and outputs functions of time (i.e. the solver itself is the output).


Figure 3.7: ForSyDe-Atom model of the RC bridge in fig. 3.6

```
rcfilter :: CT.Signal Rational }->\mathrm{ CT.Signal Rational
rcfilter s
    = let (ts, sy) = DE.toSY $ CT.toDE s
        out = SY.state21 ns ( }\mp@subsup{\lambda}{-}{}->0) ts sy
        ns p t s = euler 0.01 s (p (time t)) t
    in DE.toCT $ SY.toDE ts out
```

Testing the filter against a square wave signal, we get the response plotted in fig. 3.8. As can be clearly seen, the behavior is the right one and it reacts correctly to the inputs, taking unto account the current state of the system, i.e. its history, and it is a continuous signal in the true sense of the word.

```
/ square wave signal for testing 'rcfilter'
```



```
    Signal Rational
- I plotting the example response of 'rcfilter'
plot31 = plotGnu $ prepareL cfg {labels=["V_I","V_O"]} $ [vi3, rcfilter vi3]
```

Now let us modify osc2 from fig. 3.4 to mirror the correct behavior of the circuit in fig. 3.1. Seems to it that there is not much to do, as rcfilter already responds correctly to any input. On the other hand the RC oscillator as represented in fig. 3.1 admits as inputs both $V_{D D}$ and the control signal for $s w_{1}$ and $s w_{2}$. To correctly model that in ForSyDe-Atom we need to separate the FSM which controls the state of the capacitor as charging or discharging, which in turn will generate $V_{I}$. We do that like in fig. 3.9.

```
- / ODE-based RC oscillator model
osc4 :: CT.Signal Rational - - VDD input signal
    DE.Signal () - control signal of discrete impulses
    CT.Signal Rational - - output signal
osc4 vdd ctl
    = let - generator for the switch state variable
    swState = DE.embedSY11 (SY.stated11 swF Charging) ctl
```



Figure 3.8: Response of rcfilter

```
swF Charging _ = Discharging
```

swF Charging _ = Discharging
swF Discharging _ = Charging
swF Discharging _ = Charging
_ transforms VDD into VI for the RC filter
_ transforms VDD into VI for the RC filter
vddSwitched = DE.comb21 vddF swState \$ CT.toDE vdd
vddSwitched = DE.comb21 vddF swState \$ CT.toDE vdd
vddF Charging v = v

```
vddF Charging v = v
```




```
- state machine with ODE solver modeling an RC filter
```

- state machine with ODE solver modeling an RC filter
vOut = embed (SY.state21 nsEU ( }\mp@subsup{\lambda}{-}{}->0)\mathrm{ ) vddSwitched
vOut = embed (SY.state21 nsEU ( }\mp@subsup{\lambda}{-}{}->0)\mathrm{ ) vddSwitched
nsEU p t s = euler 0.01 s (p \$ time t) t
nsEU p t s = euler 0.01 s (p \$ time t) t
_ custom wrapper that embeds a SY tag-aware process into a DE environment
_ custom wrapper that embeds a SY tag-aware process into a DE environment
embed p de = let (t, v) = DE.toSY de
embed p de = let (t, v) = DE.toSY de
in SY.toDE t (p t v)
in SY.toDE t (p t v)
in DE.toCT vOut

```


Figure 3.9: ForSyDe-Atom model of the RC oscillator described by osc4
Testing osc4 against the same inputs as osc2 in fig. 3.5, we get the response plotted in fig. 3.10, which is now the correct behavior of the RC circuit with respect to its inputs. As expected, the output describes correctly the dynamics of the system based on the state and history of the capacitor.
|plot41 = plotGnu \$ prepareL cfg \{labels=["V_\{DD\}","V_O"]\} \$ [vdd22, osc4 vdd22 sCtrl]
We have shown three different models of an RC oscillator circuit represented in fig. 3.1 at different levels of complexity, and a model of an RC bridge represented in fig. 3.6. The fact that


Figure 3.10: Response of osc4
they are different does not mean that either is "more correct/incorrect" than the other. Either of them might very well be treated as "correct" depending on what we consider or not the acceptable inputs (note that we have not defined them on purpose). As expected, the more we consider the inputs as sources of non-determinism, the more complex our model needs to be in order to cover "special" cases. As you might guess already, this comes at a severe cost of run-time performance.

Let us briefly measure this cost in performance between the four presented models. For this, we consider two situations:

Experiment 1 we need to find out \(V_{O}\) at \(t=2.8\) seconds.
Experiment 2 we need to sample \(V_{O}\) for the whole period \(t=[0,3]\) seconds, with a precision of 50 milliseconds.

We consider the same input scenarios for all 4 models:
```

- I plotting/sampling configuration for performance testing
cfgTest = defaultCfg {xmax=3, rate=0.005}
- I VDD input for performance testing
vddTest = CT.signal [(0, 玍 }->2),(1,\mp@subsup{\lambda}{-}{}->1.5),(2,\mp@subsup{\lambda}{-}{}->1)] :: CT.Signal Rational
- I VI input for performance testing of the RC filter model

```

```

    Signal Rational
    - I Switch control signal for performance testing
ctlTest = DE.signal [(0,()),(0.5,()),(1,()),(1.5,()),(2,()),(2.5,())] :: DE.Signal ()

```

DISCLAIMER: we have measured the performance for all experiments using the :set +s directive in a ghci interpreter session, which offers some crude information about the run-time. The measurements are by no means realistic for compiled and optimized programs, but rather give a rough idea about the raw, un-optimized functions run directly in the interpreter. In any case, this offers a common ground for comparing implementations between them to observe trends and the influences of the design decisions, but not to judge ForSyDe-Atom as such.
```

\lambda> :set +s
\lambda> fromRational \$ osc1 vddTest 'CT.at' 2.8
4.9787192469339395e-2
(0.01 secs, 398,792 bytes)
\lambda> fromRational \$ osc2 vddTest ctlTest 'CT.at' 2.8

```
```

4.9787192469339395e-2
(0.01 secs, 385,952 bytes)
\lambda> fromRational \$ rcfilter viTest 'CT.at' 2.8
5.169987618408467e-2
(0.08 secs, 34,044,128 bytes)
\lambda> fromRational \$ osc4 vddTest ctlTest 'CT.at` 2.8
5.169987618408467e-2
(0.07 secs, 34,047,720 bytes)

```
\begin{tabular}{r||c|c||c|c}
\multicolumn{1}{c|}{ Model } & \multicolumn{2}{c|}{\begin{tabular}{c} 
Experiment 1 \\
time (s)
\end{tabular}} & \multicolumn{2}{c}{ Exmory (MB) }
\end{tabular}\(|\) time (s) \begin{tabular}{rl} 
memory (MB) \\
\hline osc1 & 0.01
\end{tabular}

Table 3.1: Experimental results for testing the RC models
The experimental results on an computer with Intel \(®\) Core \({ }^{\mathrm{TM}} \mathrm{i} 7-3770 \mathrm{CPU} @ 3.40 \mathrm{GHz} \times\) 8 threads, and \(31,4 \mathrm{GiB}\) RAM are shown in table 3.1. As expected, osc1 and osc2 perform in almost negligible time for both experiments due to lazy evaluation which performs computation only at the requested evaluation point. However, the lazy evaluation is costly in terms of runtime memory, fact noticed especially for rcfiler and osc4 due to the fact that intermediate structures need to be stored before evaluating. The complexity of rcfilter and osc4 are seen both in the execution time and in the memory consumption. Apart from the cost in performance, model fidelity for unknown inputs came also with a high price in loss of precision due to chained numerical computation, fact seen from the evaluation results of Experiment 1 in the interpreter listing above. Choosing a better solver than euler, e.g. a Runge-Kutta solver, or even better, a symbolic solver, could improve both performance and precision, but that is out of the scope of this report. Another, rather surprising fact is that osc2 performs slightly better than osc1, probably to the lack of tag calculus in the SY domain, i.e. tags are mainly passed untouched between interfaces, whereas calculations are performed in a synchronous reactive manner on values only.


Figure 3.11: Execution time and memory consumption for Experiment 2

To get a feeling of the cost complexity of the models for osc1, osc2 and osc4, we plot the execution time and memory consumption during Experiment 2 for three resolutions: 10 milliseconds, 50 milliseconds and 100 milliseconds. This way we can observe the trends in fig. 3.11, where the most notable one is the exponential growth in execution time for osc4. This can be explained recalling the so-called "time leaks" observed by hudak03 (among others), which occur due to the fact that lazy evaluation forces to recalculate all previous states of the ODE solver in order to evaluate the current sample. In functional reactive programming (FRP), a solution to time leaks is the concept of "continuation", which is roughly what happens when the synchronous reactive FSM receives a new discrete event: it stores the current state to the ODE to be used in
the future. Knowing this, let us see what happens when we double the number of discrete events within the input \(V_{D D}\) signal.
```

- I VDD input for performance testing, having twice more discrete events than 'vddTest'

```

```

    Signal Rational
    ```

\subsection*{3.3 Conclusion}

This project is still under development...

\subsection*{3.4 References}

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\section*{Part II}

\section*{Tools \& Libraries Documentation}


\section*{The ForSyDe-Аtom Standard Library}

This chapter is an extended API documentation of ForSyDe-Atom Standard Library version \(0.1 .1^{1}\) which was organized to also serve the purpose of technical report. It treats both theoretical and practical aspects, justifying the implementation and providing examples of usage for most of the library-exported functions. It has been generated from the inline documentation using a Haddock \({ }^{2}\) generator.

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\subsection*{4.1 Introduction}

The ForSyDe (Formal System Design) methodology has been developed with the objective to move system design to a higher level of abstraction and to bridge the abstraction gap by transformational design refinement. It targets the modelling and characterization of cyber-physical systems inheriting the theory of Models of Computation (MoCs), providing both a correct-by-construction execution model and analyzable entry point for further synthesis flows. For more information about ForSyDe and its associated projects please consult the ForSyDe webpage.

The forsyde-atom library is a shallow-embedded DSL implementing the execution semantics of an atom-based approach to ForSyDe. Its purpose is to provide a modeling framework for cyberphysical systems and to serve as a proof-of-concept for a future (non-strict typed) DSL. Adhering to the formalism set by the PhD Thesis of Ingo Sander, the current framework views systems as networks of processes communicating through signals. The atom-based approach to ForSyDe adds some new important concepts:
- the separation of concerns through semantically-independent layers of computation, behavior, synchronization and structure. Each layer offers a different analyzable view of a given system and is described using the concept of higher-order functions.
- the description of each layer as a network of primitive building blocks called atoms. Each atom embeds an undividable operation, and wraps functions of lower layers with the semantics dictated by a higher layer. They are described using the powerful concept of applicative functors.
- the complete autonomy of atoms in relation to the patterns they build. As such, process networks or constructors are nothing but structural ad-hoc compositions that aid in achieving complex behaviors whereas the actual behavior is dictated by the atoms alone. This is proved (for the time being) for the Synchronization Layer by implementing all MoCs as instances of only one type class.

While the host language limits the possibility of providing general (e.g. proces) constructors due to type-strictness, the forsyde-atom framework is by all means complete. In this sense the user can create her own custom correct-by-design constructors and networks as compositions of the provided atoms and utilities. Also, this haddock-generated page is organized both as an API documentation and as a technical report to facilitate te use and understanding of the formal principles behing the design process.

\subsection*{4.1.1 Naming convention}

All multi-parameter patterns and utilities provided by the library API as higher-order functions are named along the lines of functionMN where \(M\) represents the number of curried inputs (i.e. a1 -> a2 -> ... -> aM), while N represents the number of tupled outputs (i.e. (b1,b2, .., bN)). To avoid repetition we only provide documentation for functions with 2 imputs and 2 outputs (i.e. function22), while the available ones are mentioned as a regex (i.e. function[1-4][1-4]). In case the provided functions do not suffice, feel free to implement your own patterns following the example in section 2.4.

\subsection*{4.2 ForSyDe.Atom}
```

module ForSyDe.Atom (

```
    \(\operatorname{ExB}(\) extend, \((/ . \backslash),(/ * \backslash),(/ \& \backslash),(/!\backslash)), \quad \operatorname{Stream}(N u l l S,(:-))\),
    module ForSyDe.Atom.MoC.Stream, Time, TimeStamp,
    MoC(Fun, Ret, (-.-), (-*-), (-*), (-<-), (-\&-)),
    Skeleton \(((=.=),(=*=),(=\backslash=),(=\ll=)\), first, last),
    module ForSyDe.Atom.Utility, (||<)
    ) where

The formal foundation upon which ForSyDe I. Sander and Jantsch, 2006 defines its semantics is the tagged signal model Lee and Sangiovanni-Vincentelli, 1998. This is a denotational framework introduced by Lee and Sangiovanni-Vincentelli as a common meta model for describing properties of concurrent systems in general terms as sets of possible behaviors. Systems are regarded as compositions of processes acting on signals which are sets of tagged events. Signals are characterized by a tag system which determines causality between events, and could model time, precedence relationships, synchronization points, and other key properties. Based on how tag systems are defined, one can identify several Models of Computations (MoCs) as classes of behaviors dictating the semantics of execution and concurrency in a network of processes.

These concepts are the supporting pillars of ForSyDe's philosophy, and state the purpose of the forsyde-atom library: it is supposed to be a modelling framework used as a proof-of-concept for the atom-based approach to cyber-physical systems Ungureanu and Ingo Sander, 2017. This approach extends the tagged signal model by systematically deconstructing processes to their basic semantics and recreating them using a minimal language of primitive building blocks called atoms. It also tries to expand the scope of this model by exploiting more aspects than just timing, by adding primitives for parallelism, behavior, etc.

The API documentation is structured as follows: this page provides an overview of the general notions and concepts, gently introducing the separate modules and the motivation behind them. Each major module corresponds to a separate layer Ungureanu and Ingo Sander, 2017 implemented as a type class. The documentation pages for each layer and for each of their instances
contains in-depth knowledge and examples, and can be accessed from the contents page or by following the links suggested. For more complex examples and tutorials follow the links in the project web page.

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.

\subsection*{4.2.1 The layered process model}

The forsyde-atom project is led by three main policies:
1. in order to cope with the complexity of cyber-physical systems (CPS) it tries to separate the concerns such as computation, timing, synchronization, parallelism, structure, behavior, etc.
2. in order to have a small, ideally minimal grammar to reason about systems correctness, it aims to provide primitive (indivisible) operators called atoms as building blocks for independently developing complex aspects of a system's execution through means of composition or generalization.
3. in order to express complex behaviors with a minimal grammar, it decouples structure (composition) from meaning (semantics), the only semantics carriers being atoms. Thus complex behaviors can be described in terms of patterns of atoms. Using ad-hoc polymorphism, atoms can be overloaded with different semantics triggered by the data type they input, whereas their composition is always the same.
atom the elementary (primitive, indivisible) constructor which embeds a set of semantics relevant for their respective layer (e.g. timing, behavioural, structural, etc.)
atom patterns meaningful compositions of atoms. They are provided as constructors which need to be properly instantiated in order to be used. We also use the term "pattern" to differentiate atom compositions as constructors from atoms as constructors.

The first policy, i.e. the separation of concerns led to the so-called layered process model which is reflected in the library by providing separate major modules associated with each layer. Layers as such are independent collections of entities for modeling different aspects of CPS. These aspects interact through means of higher-order functions, wrapping each other in as structured fashion in a way which can be visualized as below.

Layers are implemented as type classes which imply:
- atoms as function signatures belonging to the type class;
- patterns which are compositions atoms, provided as constructors;
- data types for all the classes of behaviors concerning the aspect described by the layer in question. These types instantiate the above type class and overload the atoms with semantics in accordance to the behavior described. For example, the MoC layer is currently instantiated by types describing the CT, DE, SY and SDF MoCs.


In order to model interleaving aspects of CPS, layers interact with each other through means of higher order functions. As such, each layer describes some atoms as higher-order functions which take entities belonging to another layer as arguments. Intrinsically, the data types belonging to a layer may be wrapping types of other layers, as depicted in the figure above. For a short comprehensive overview on layers, please refer to Ungureanu and Ingo Sander, 2017.

By convention, the first (innermost) layer is always the function layer which describes arbitrary functions on data and expresses the system's functional aspects. In the following paragraphs we will give an overview of the "outer" layers currently implemented in forsyde-atom, which in comparison, express the extra-functional aspects of a system (timing, behavior, synchronization, and so on).

\subsection*{4.2.2 The Extended Behavior (ExB) Layer}

As seen in layered process model, the extended behavior layer expands the set of possible behaviors implied by a layer (typically the function layer), by defining a set of symbols with known semantics, and adding it to (i.e. wrapping) the pool of possible values or states.

While semantically the ExB layer extends the value pool in order to express special events (e.g. error messages or even the complete absence of events), it practically provides an independent environment to model events with a default/known response, independently of the data path. These responses are particularly captured by atoms, thus enforcing the high-level separation of concerns between e.g. control and data paths.

This layer provides:
- a set of extended behavior atoms defining the interfaces for the resolution and response functions, as part of the ExB type class (see below).
- a library of function wrappers as specific atom patterns (Check the ForSyDe.Atom.ExB module for extensive documentation).
- a set of data types defining classes of behaviors and instantiating the ExB type class (check the links in the instances section for extensive documentation).

\section*{class Functor b => ExB b where}

Class which defines the atoms for the extended behavior layer.

As its name suggests, this layer is extending the behavior of processes (or merely of functions if we ignore timing semantics completely), by expanding the domains of the wrapped layer (e.g. the set of values) with symbols having clearly-defined semantics (e.g. special events with known responses).
The types associated with this layer can simply be describes as:
\[
\mathcal{B}(\alpha)=\alpha \cup b
\]
where \(\alpha\) is a base type and \(b\) is the type extension, i.e. a set of symbols with clearly-defined semantics.
Extended behavior atoms are functions of these types, defined as interfaces in the ExB type class.

\section*{Methods}
```

extend :: a -> b a

```

Extends a value (from a layer below) with a set of symbols with known semantics, as described by a type instantiating this class.
(/.\\) :: (a -> a) -> b a -> b a
Basic functor operator. Lifts a function (from a layer below) into the domain of the extended behavior layer.
\[
\triangle:(\alpha \rightarrow \beta) \rightarrow \mathcal{B}(\alpha) \rightarrow \mathcal{B}(\beta)
\]
(/*\\) :: b (a -> a) -> b a -> b a
Applicative operator. Defines a res between two extended behavior symbols.
\[
\triangle: \mathcal{B}(\alpha \rightarrow \beta) \rightarrow \mathcal{B}(\alpha) \rightarrow \mathcal{B}(\beta)
\]
(/\& \()\) ) : b Bool \(->\mathrm{b}\) a \(->\mathrm{b}\) a
Predicate operator. Generates a defined behavior based on an extended Boolean predicate.
\[
\notin \mathcal{B}(\mathrm{Bool}) \rightarrow \mathcal{B}(\alpha) \rightarrow \mathcal{B}(\alpha)
\]
(/! \\) :: a -> b a -> a
\(\quad\) Degrade operator. Degrades a behavior-extended value into a non-extended one (from a layer below), based on a kernel value. Used also to throw exceptions.
instance ExB AbstExt
Implements the absent semantics of the extended behavior atoms.

\subsection*{4.2.3 The Model of Computation (MoC) Layer}

This layer represents a major part of the forsyde-atom library and is concerned in modeling the timing aspects of CPS. While its foundations have been layered in the classical ForSyDe I. Sander and Jantsch, 2006, it is mainly inspired from Lee and Sangiovanni-Vincentelli, 1998 an it tries to follow the tagged signal model as closely as it is permitted by the host language, and with the adaptations require by the atom approach.

Although a short introduction of the tagged signal model has been written in the introduction of this documentation, we feel obliged to provide a primer in the classical ForSyDe theory in order to understand how everything fits together.

\section*{Signals}

Lee and Sangiovanni-Vincentelli, 1998 defines signals as (ordered) sets of events where each event is composed of a tag \(T\) and a value \(V\). Similarly, in ForSyDe a signal is defined as a (partially
or totally) ordered sequence of events that enables processes to communicate and synchronize. Sequencing might infer an implicit order of events, but more importantly it determines an order of evaluation, which is a key piece of a simulation engine.
\[
\begin{aligned}
& s=\left\{e_{0}, e_{1}, \ldots\right\} \in S \\
& \text { where } e_{j}=\left(t_{j}, v_{j}\right) \\
& \quad v_{j} \in V, t_{j} \in T \\
& \quad t_{j} \leq t_{j+1}, \forall j \in \mathbb{N}
\end{aligned}
\]

In ForSyDe-Atom, sequencing is achieved using the Stream data type, inspired from Reekie, 1995. In ForSyDe-Atom, signals are streams that carry events, where each type of event is identified by a type constructor which defines its tag system. In other words, we can state that through its tag system, a signal is bound to a MoC.
```

data Stream e

```
\(=\) NullS terminates a signal
I e (:-) (Stream e) the default constructor appends an event to the head of the stream
Defines a stream of events, encapsulating them in a structure isomorphic to an infinite list R. S. Bird, 1987, thus all properties of lists may also be applied to Streams. While, in combination with lazy evaluation, it is possible to create and simulate infinite signals, we need to ensure that the first/previous event is always fully evaluated. This can be translated into the following rule:
zero-delay feedbacks are forbidden, due to un-evaluated self-referential calls. In a feedback loop, there always has to be enough events to ensure the data flow.

This rule imposes that the stream of data is uninterrupted in order to have an evaluatable kernel every time a new event is produced (i.e. to avoid deadlocks). Thus we can add the rule:
cleaning of output events is also forbidden. In other words, for each new input at any instant in time, a process must react with at least one output event.
```

instance Functor Stream

```
allows for the mapping of an arbitrary function (a -> b) upon all the events of a (Stream a).
instance Applicative Stream
enables the Stream to behave like a ZipList
instance Foldable Stream
provides folding functions useful for implementing utilities, such as length.
instance Read \(a \quad\) => Read (Stream a)
signal (1 :- 2 :- Nulls) is read using the string " \(\{1,2\}\) ".
instance Show a => Show (Stream a)
signal (1 :- \(2:-\) NullS) is represented as \(\{1,2\}\).
instance Plottable a => Plot (Signal a)
SY signals.
instance Plottable a => Plot (Signal a)
SDF signals.
instance Plottable a => Plot (Signal a)
DE signals.
```

instance Plottable a => Plot (Signal a)
CT signals.

```

For extended documentation and a list of all utilities associated with the Stream type you can consult:
module ForSyDe.Atom.MoC.Stream

\section*{Processes}

As described in Lee and Sangiovanni-Vincentelli, 1998, processes are either "set of possible behaviors" of signals or "relations" between multiple signals. One can describe complex systems by composing processes, which in this case is interpreted as the "intersection of the behaviors of each of the processes being involved".
monotonicity In order to ensure causal order and determinancy, processes need to be monotonic Lee and Sangiovanni-Vincentelli, 1998. A signal's tags (if explicit) must be a partial or total order and all tag alterations must be monotonic.

ForSyDe inherits this definition with respect to a functional view, thus a process \(p\) is a functional mapping over (the history of) signals. A process can only be instantiated using a process constructor \(p c\), which is a higher order function embedding MoC semantics and/or a specific composition, but lacking functionality.
\[
\begin{array}{ll}
p: S^{m} \rightarrow S^{n} \in P & p c_{\mathrm{M}}: V^{n} \rightarrow P \in P C \\
p(s, \ldots)=(s, \ldots) & p c_{\mathrm{M}}(v, \ldots)=p
\end{array}
\]

Since processes are functions, process composition is equivalent to function composition. This means that composing two processes p 1 and p 2 grants the process p 2 . p1
```

p1 :: Signal }\alpha->\mathrm{ Signal }
p2 :: Signal }\beta->\mathrm{ Signal }
p2. p1 :: Signal \alpha -> Signal }

```

This implies that there is a signal Signal \(\beta\) that "streams" the result from p 1 to p 2 , as suggested in the drawing:
\[
\begin{gathered}
p_{1}: S_{\alpha} \rightarrow S_{\beta} \\
p_{2}: S_{\beta} \rightarrow S_{\gamma} \\
p_{2} \circ p_{1}: S_{\alpha} \rightarrow S_{\gamma} \\
\left(p_{2} \circ p_{1}\right)(s) \equiv p_{2}\left(p_{1}(s)\right)
\end{gathered}
\]


Process networks describe ForSyDe systems in terms of compositions of processes and originate from Reekie's process nets Reekie, 1995. A process network is a process itself, i.e. function from signal(s) to signal(s). The composition above p2 . p1 can also be regarded as a process network.

In ForSyDe-Atom atoms can be regarded as process constructors as their instantiations are functions on signals of events. Instantiations of atom patterns are the exact equivalent of process networks, which themselves are also processes, depending on the level of abstraction you are working with (hierarchical blocks vs. flat structures).

To understand the versatility of composition and partial application in building process constructors, consider the example above where composition of two processes infers a signal between them. This mechanism also works when composing constructors (un-instantiated atoms), which yields another constructor. By instantiating (fully applying) the new constructor we obtain a process network equivalent to the composition of the respective primitive processes obtained by instantiating (fully applying) the component atoms, like in the example below:
\[
\begin{aligned}
p_{1}\left(s_{1}\right) & =(+) \odot s_{1} \\
p_{2}\left(s_{1}, s_{2}\right) & =s_{1} \circledast s_{2} \\
p n\left(s_{1}, s_{2}\right) & =\left(p_{2}\left(s_{2}\right) \circ p_{1}\right)\left(s_{1}\right) \\
p c^{\prime}(f)\left(s_{1}, s_{2}\right) & =\left(\circledast s_{2}\right) \circ\left(f \odot s_{1}\right) \\
p^{\prime}\left(s_{1}, s_{2}\right) & =p c^{\prime}(+)\left(s_{1}, s_{2}\right) \\
p n & \equiv p^{\prime}
\end{aligned}
\]

Now if we visualize process networks as graphs, where processes are nodes and signals are edges, a meaningful process composition could be regarded as graph patterns. Therefore it is safe to associate process constructors as patterns in process networks.

\section*{Models of Computation}

As mentioned in the introduction, MoCs are classes of behaviors dictating the semantics of execution and concurrency in a network of processes. Based on the definitions of their tag systems ForSyDe identifies MoCs as:
1. timed where \(T\) is a totally ordered set and \(t\) express the notion of physical time (e.g. continuous time CT, discrete event DE) or precedence (e.g. synchronous SY);
2. untimed, where \(T\) is a partially ordered set and \(t\) is expressed in terms of constraints on the tags in signals (e.g. dataflow, synchronous data flow SDF).

As concerning MoCs, ForSyDe implements the execution semantics through process constructors, abstracting the timing model and inferring a schedule of the process network. In ForSyDe-Atom all atoms embed operating semantics dictated by a certain MoC and are side-effect-free. This ensures the functional correctness of a system even from early design stages.

\section*{Representing Time}

For explicit time representation, ForSyDe-atom provides two distinct data types.

\section*{type Time = Rational}

Type alias for the type to represent metric (continuous) time. Underneath we use Rational that is able to represent any \(t\) between \(t_{1}<t_{2} \in T\).
type TimeStamp = DiffTime
Alias for the type representing discrete time. It is inherently quantizable, the quantum being a picosecond ( \(10^{-12}\) seconds), thus it can be considered order-isomorphic with a set of integers, i.e. between any two timestamps there is a finite number of timestamps. Moreover, a timestamp can be easily translated into a rational number representing fractions of a second, so the conversion between timestamps (discrete time) and rationals (analog/continuous time) is straightforward.

This type is used in the explicit tags of the DE MoC (and subsequently the discrete event evaluation engine for simulating the CT MoC ).

\section*{MoC Layer Overview}

This layer consists of:
- 4 atoms as infix operators, implemented as methods of the type class Moc. Since each MoC is determined by its tag system, we expose this which are instances of this class. Thus an event's type will trigger an atom to behave in accordance to its associated MoC.
- a library of meaningful atom patterns as process constructors. (Check the ForSyDe.Atom.MoC module for extensive documentation).
- a set of data types defining tag systems through the structure of events (i.e. \(T \times V\) ). They are instances of the MoC type class and define the rules of execution that will trigger an atom to behave in accordance to its associated MoC. For each supported MoC, forsyde-atom provides a module which defines the signal (event) type, but also a set of utilities and process constructors as specific instantiations of atom patterns. (Check the links in the instances section for extensive documentation).

\section*{class Applicative e => MoC e where}

This is a type class defining interfaces for the MoC layer atoms. Each model of computation exposes its tag system through a unique event constructor as an instance of this class, which defines \(T \times V\).

To express all possible MoCs which can be described using a tagged signal model we need to capture the most general form of their atoms. Recall that all atoms in the layered framework are represented as higher-order functions on structured types (instances of this class), taking functions of other (lower) layers as arguments. While this principle stands also for this layer, the functions taken as arguments need to be formatted for each MoC in particular in order to capture additional information, which we can call in general terms as the execution context.
One typical example of additional information is the consumption and production rates of for data flow MoCs (e.g. SDF). In this case the passed functions are defined over "partitions" of events, i.e. groupings of events with the same partial order in relation to, for example, a process firing. The formal description of such a "formatted function" taken as argument by a MoC entity is:
\[
\dot{a} \rightarrow \dot{b} \equiv \dot{\alpha}_{1} \times \dot{\alpha}_{2} \times \ldots \rightarrow \dot{\beta}_{1} \times \dot{\beta}_{2} \times \ldots
\]
where \(a\) and \(b\) might be Cartesian products of different types, corresponding to how many signals the constructor is applied to or how many signals it yields, and each type is expressed as:
\[
\begin{aligned}
& \dot{\alpha}=\alpha^{r} \\
& r \in \begin{cases}\{1\}, & \text { for all timed MoCs (e.g. SY, DE, CT) } \\
\mathbb{N}, & \text { for static variants of SDF } \\
\mathbb{N}^{n}, & \text { for CSDF } \\
S \times \mathbb{N} \rightarrow \mathbb{N}, & \text { where } S \text { is a state space, } \\
& \text { in the most general case of untimed data flow }\end{cases}
\end{aligned}
\]

While, as you can see above, the execution context can be extracted from the type information, working with type-level parameters is not a trivial task in Haskell, especially if we want to describe a general and extensible type class. This is why we have chosen a pragmatic approach in implementing the MoC class:
- any (possible) Cartesian product of \(\alpha\) is represented using a recursive type, namely a list \([\alpha]\).
- as the execution context cannot (or can hardly) be extracted from the recursive type, in the most general case we pass both context and argument as a pair (see each instance in particular). To aid in pairing contexts with each argument in a function, the ctxt utilities are provided (see ctxt22).
- this artifice was masked using the generic type families Fun and Res.

\section*{Methods}

Fun \(e \mathrm{a} b\)
This is a type family alias for a context-bound function passed as an argument to a MoC atom. In the most simple case it can be regarded as an enhanced \(->\) type operator. While hiding the explicit definition of arguments, this implementation choice certainly has its advantages in avoiding unnecessary or redundant type constructors (see version 0.1 .1 and prior). Aliases are replaced at compile time, thus not affecting run-time performance.
\[
\dot{\alpha} \rightarrow \beta
\]

Ret e b
As with Fun, this alias hides a context-bound value (e.g. function return). Although the definition seems to be redundant with Fun, this alias is needed for utilities to recreate clean types again (see -*).

\section*{(-.-) : : Fun e a b -> Stream (e a) -> Stream (e b)}

This atom is mapping a function on values (in the presence of a context) to a signal, i.e. stream of tagged events. As ForSyDe deals with determinate, functional processes, this atom defines the (only) behavior of a process in rapport to one input signal Lee and Sangiovanni-Vincentelli, 1998.
\[
\begin{gathered}
\odot:(\dot{\alpha} \rightarrow \beta) \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\beta) \\
(-*-): \text { : Stream (e (Fun e a b)) } \rightarrow \text { Stream (e a) } \rightarrow \text { Stream (e b) }
\end{gathered}
\]

This atom synchronizes two signals, one carrying functions on values (in the presence of a context), and the other containing values, during which it applies the former on the latter. As concerning the process created, this atom defines a relation between two signals Lee and Sangiovanni-Vincentelli, 1998.
\[
\begin{aligned}
* & : \mathcal{S}(\dot{\alpha} \rightarrow \beta) \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\beta) \\
(-*): & \text { Stream (e (Ret e b)) } \gg \operatorname{Stream}(\mathrm{e} \text { b) }
\end{aligned}
\]

Artificial utility which drops the context and/or partitioning yielding a clean signal type.
\[
\mathcal{S}(\dot{\beta}) \rightarrow \mathcal{S}(\beta)
\]
(-<-) : : Stream (e a) -> Stream (e a) -> Stream (e a)
This atom appends a (partition of) events at the beginning of a signal. This atom is necessary to ensure complete partial order of a signal and assures the least upper bound necessary for example in the evaluation of feedback loops Lee and SangiovanniVincentelli, 1998.
\[
\theta: \dot{\alpha} \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\alpha)
\]

Notice the difference between the formal and the implemented type signatures. In the implementation the value/partition is wrapped inside an event type to enable smooth composition. You might also notice the type for the initial event(s) as being wrapped inside a signal constructor. This allows defining an DSL for this layer which is centered around signals exclusively, while also enabling to define atoms as homomorphisms to certain extent R. Bird and Moor, 1997. Certain MoCs might have additional constraints on the first operand to be finite.
(-\&-) :: Stream (e a) -> Stream (e a) -> Stream (e a)
This atom allows the manipulation of tags in a signal in a restrictive way which preserves monotonicity and continuity in a process Lee and Sangiovanni-Vincentelli, 1998, namely by phase-shifting all tags in a signal with the appropriate metric corresponding to each MoC. Thus it preserves the characteristic function intact I. Sander and Jantsch, 2006.
\[
\oplus(T \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\alpha)
\]

As with the -<- atom, we can justify the type signature for smooth composition and the definition of atoms as homomorphisms to certain extent. This in turn allows the interpretation of the -Methods- operator as aligning the phases of two signals: the second operand is aligned based on the first.
```

instance MoC SY

```

Implenents the execution and synchronization semantics for the SY MoC through its atoms.
```

instance MoC SDF

```

Implenents the SDF semantics for the MoC atoms

\section*{instance MoC DE}

Implenents the execution and synchronization semantics for the DE MoC through its atoms.
```

instance MoC CT

```

Implenents the execution and synchronization semantics for the CT MoC through its atoms.

\subsection*{4.2.4 The Skeleton Layer}

The skeleton layer describes recursive and regular composition of processes which expose inherent potential for parallelism. As such, it wraps lower layer entities (i.e. processes, signals), into regular structures called categorical types. Most of the ground work for this layer is based on the categorical type theory R. Bird and Moor, 1997, which enable the description of algorithmic skeletons as high-level constructs encapsulating parallelism and communication with an associated cost complexity.

This layer provides:
- 3 atoms as infix operators which, as demonstrated in R. Bird and Moor, 1997 and Skillicorn, 1994, are enough to describe all algorithmic skeletons.
- a library of generic skeletons as specific atom patterns. (Check the ForSyDe.Atom.Skeleton module for extensive documentation).
- a set of different categorical types which implement these atoms, as instances of the Skeleton type class. These types provide additional skeletons patterns of atoms which takes as arguments their own type constructors. (Check the links in the instances section for extensive documentation).
class Functor c => Skeleton c where
Class containing all the Skeleton layer atoms.
This class is instantiated by a set of categorical types, i.e. types which describe an inherent potential for being evaluated in parallel. Skeletons are patterns from this layer. When skeletons take as arguments entities from the MoC layer (i.e. processes), the results themselves are parallel process networks which describe systems with an inherent potential to be implemented on parallel platforms. All skeletons can be described as composition of the three atoms below (=«= being just a specific instantiation of \(=\backslash=\) ). This possible due to an existing theorem in the categorical type theory, also called the Bird-Merteens formalism R. Bird and Moor, 1997:
factorization A function on a categorical type is an algorithmic skeleton (i.e. catamorphism) iff it can be represented in a factorized form, i.e. as a map composed with a reduce.

Consequently, most of the skeletons for the implemented categorical types are described in their factorized form, taking as arguments either:
- type constructors or functions derived from type constructors
- processes, i.e. MoC layer entities

Most of the ground-work on algorithmic skeletons on which this module is founded has been laid in R. Bird and Moor, 1997, Skillicorn, 1994 and it founds many of the frameworks collected in Fischer, Gorlatch, and Bischof, 2003.

\section*{Methods}

\section*{(=.=) : : (a -> b) -> c a -> c b}

Atom which maps a function on each element of a structure (i.e. categorical type), defined as:
\[
\stackrel{\diamond}{ }(\alpha \rightarrow \beta) \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\beta)
\]
\(=.=\) together with \(=*=\) form the map pattern.
(=*=) :: c (a -> b) -> c a -> c b
Atom which applies the functions contained by as structure (i.e. categorical type), on the elements of another structure, defined as:
\[
\Leftrightarrow: \mathcal{C}(\alpha \rightarrow \beta) \rightarrow \mathcal{C}(\alpha) \rightarrow \mathcal{C}(\beta)
\]
\(=.=\) together with \(=*=\) form the map pattern.
(=\=) :: (a -> a -> a) -> c a -> a
Atom which reduces a structure to an element based on an associative function, defined as:
(=《=)
\[
\otimes:(\alpha \times \alpha \rightarrow \alpha) \rightarrow \mathcal{C}(\alpha) \rightarrow \alpha
\]

Skeleton which pipes an element through all the functions contained by a structure. N.B.: this is not an atom. It has an implicit definition which might be augmented by instances of this class to include edge cases.
\[
\begin{aligned}
& \forall: \mathcal{C}(\alpha \rightarrow \alpha) \times \alpha \rightarrow \alpha \\
& \forall=(\circ) \diamond
\end{aligned}
\]

As the composition operation is not associative, we cannot treat pipe as a true reduction. Alas, it can still be exploited in parallel since it exposes another type of parallelism: time parallelism.
first : : c a -> a
Returns the first element in a structure.N.B.: this is not an atom. It has an implicit definition which might be replaced by instances of this class with a more efficient implementation.
\[
\begin{aligned}
& \text { first }_{S}: \mathcal{C}(\alpha) \rightarrow \alpha \\
& \text { first }_{S}=(\lambda a b \rightarrow a) \diamond
\end{aligned}
\]
last : : c a -> a
Returns the last element in a structure.N.B.: this is not an atom. It has an implicit definition which might be replaced by instances of this class with a more efficient implementation.
\[
\begin{aligned}
& \text { last }_{S}: \mathcal{C}(\alpha) \rightarrow \alpha \\
& \text { last }_{S}=(\lambda a b \rightarrow b) \diamond
\end{aligned}
\]

Ensures that Vector is a structure associated with the Skeleton Layer.

\subsection*{4.2.5 Utilities}

The Atom module exports a set of utility functions, mainly for aiding the designer to avoid working with zipped tuples which might pollute the design. Utilities are function without any semantical value (thus not considered atoms). They operate on and might alter the structure of some datum, but it does not affect its state.

For a list of all the provided utilities, please consult the following module:
```

module ForSyDe.Atom.Utility

```

Among the most useful utilities we mentions the unzip function. Recall that in all our definitions for patterns, they were expressed in the most general form as functions from \(n\)-ary Cartesian products to \(m\)-ary Cartesian products. While partial application provides a versatile mechanism that can translate n-ary inputs into curried arguments (which is very powerful in combination with an applicative style), we cannot do so for return types. For the latter we must rely on tuples. But working with tuples of data wrapped in several layers of structures becomes extremely cumbersome. Take for example the case of a process constructed with \(p c\) in equation (1) below. Using only the provided atoms to implement \(p c\) would give us a process which returns only one signal of a tuple and not, as we would like, a tuple of signals of events.
\[
\begin{aligned}
& p c:\left(\alpha \rightarrow \beta^{n}\right) \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}\left(\beta^{n}\right) \\
& \triangleleft: \mathcal{S}\left(\beta^{n}\right) \rightarrow \mathcal{S}(\beta)^{n} \\
&(p c f \triangleleft): \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\beta)^{n}
\end{aligned}
\]


Therefore, by implementing all data types associated with signals and events as instances of Functor, we were able to provide a (set of) unzip utility functions defined as in equation (2) above, in the ForSyDe.Atom.Utility module. Mind that we call unzip a utility and not an atom, since it has no synchronization nor behavior semantic. It just conveniently "lifts" the wrapped tuples in order to create "collections" of events and signals, and it is imposed by the mechanisms of the type system in the host language.
(||<) :: (Functor f1, Functor f2) \(\Rightarrow\)
f1 (f2 (a1, a2)) -> (f1 (f2 a1), f1 (f2 a2))
This set of utility functions "unzip" nested n-tuples, provided as postfix operators. They are crucial for reconstructing data types from higher-order functions which input functions with multiple outputs. It relies on the nested types being instances of Functor.

The operator convention is ( \(1+<+\) ), where the number of I represent the number of layers the n-tuple is lifted, while the number of \(<+1\) is the order \(n\) of the n-tuple.

ForSyDe.Atom.Utility exports the constructors below. Please follow the examples in the source code if they do not suffice:


\section*{Example:}
```

\lambda> :set -XPostfixOperators
\lambda> ([Just (1,2,3), Nothing, Just (4,5,6)] ||<<)
([Just 1,Nothing, Just 4],[Just 2,Nothing, Just 5], [Just 3,Nothing, Just 6])

```

\subsection*{4.3 ForSyDe.Atom.ExB}
```

module ForSyDe.Atom.ExB (
ExB(extend, (/.\), (/*\), (/\&\), (/!\)), res22, filter, filter',
degrade, ignore22
where

```

This module exports the core entities of the extended behavior layer: interfaces for atoms and common patterns of atoms. It does NOT export any implementation or instantiation of any specific behavior extension type.

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.

\subsection*{4.3.1 Atoms}

\section*{class Functor b => ExB b where}

Class which defines the atoms for the extended behavior layer.
As its name suggests, this layer is extending the behavior of processes (or merely of functions if we ignore timing semantics completely), by expanding the domains of the wrapped layer (e.g. the set of values) with symbols having clearly-defined semantics (e.g. special events with known responses).
The types associated with this layer can simply be describes as:
\[
\mathcal{B}(\alpha)=\alpha \cup b
\]
where \(\alpha\) is a base type and \(b\) is the type extension, i.e. a set of symbols with clearly-defined semantics.
Extended behavior atoms are functions of these types, defined as interfaces in the ExB type class.

\section*{Methods}
extend : : a -> b a Source
Extends a value (from a layer below) with a set of symbols with known semantics, as described by a type instantiating this class.
\((/ . \backslash)::(\mathrm{a}->\mathrm{a})->\mathrm{b}\) a \(->\mathrm{b}\) a
Basic functor operator. Lifts a function (from a layer below) into the domain of the extended behavior layer.
\[
\triangle:(\alpha \rightarrow \beta) \rightarrow \mathcal{B}(\alpha) \rightarrow \mathcal{B}(\beta)
\]
\((/ * \backslash):: \mathrm{b}(\mathrm{a}->\mathrm{a})->\mathrm{b} \mathrm{a}->\mathrm{b} \mathrm{a}\)
Applicative operator. Defines a res between two extended behavior symbols.
\[
\triangle: \mathcal{B}(\alpha \rightarrow \beta) \rightarrow \mathcal{B}(\alpha) \rightarrow \mathcal{B}(\beta)
\]
(/\&\\) :: b Bool -> b a -> b a Source
Predicate operator. Generates a defined behavior based on an extended Boolean predicate.
\[
\text { A }: \mathcal{B}(\text { Bool }) \rightarrow \mathcal{B}(\alpha) \rightarrow \mathcal{B}(\alpha)
\]
(/! \\) : : a -> b a -> a Source
Degrade operator. Degrades a behavior-extended value into a non-extended one (from a layer below), based on a kernel value. Used also to throw exceptions.
\[
\triangle: \alpha \rightarrow \mathcal{B}(\alpha) \rightarrow \alpha
\]
instance ExB AbstExt
Implements the absent semantics of the extended behavior atoms.

\subsection*{4.3.2 Patterns}
```

res22

```
    : : ExB b
    \(\Rightarrow \quad\left(\mathrm{a} 1->\mathrm{a} 2->\left(\mathrm{a} 1^{\prime}, \mathrm{a} 2^{\prime}\right)\right)\) function on values
    \(->\) b a1 first input
    -> b a2 second input
    -> (b a1', b a2') tupled output
\[
\begin{aligned}
& \mathbb{A}:\left(V^{n} \rightarrow V^{m}\right) \rightarrow \mathcal{B}^{n} \rightarrow \mathcal{B}^{m} \\
& f \Perp\left(b_{1}, b_{2}, \ldots b_{n}\right)=f \triangleleft b_{1} \triangle b_{2} \triangle \ldots \star b_{n} \triangleleft
\end{aligned}
\]

The res behavior pattern lifts a function on values to the extended behavior domain, and applies a resolution between two extended behavior symbols.
Constructors: res [1-8] [1-4].
```

filter :: ExB b => b Bool -> b a -> b a

```

Prefix name for the prefix operator \(/ \& \backslash\).
```

filter' :: ExB b => Bool -> a -> b a

```

Same as filter but takes base (non-extended) values as input arguments.
```

degrade :: ExB b => a -> b a -> a

```

Prefix name for the degrade operator \(/!\backslash\).

\[
\begin{aligned}
& \text { ignore }_{B}:\left(V^{n} \times V^{m} \rightarrow V^{n}\right) \rightarrow V^{n} \times \mathcal{B}^{m} \rightarrow V^{n} \\
& \text { ignore }_{B}(f)\left(a^{n}, b^{m}\right)=a^{n} ₫\left(f\left(a^{n}\right) \Perp b^{m}\right)
\end{aligned}
\]

The ignoreXY pattern takes a function of \(\mathrm{Y}+\mathrm{X}\) arguments, Y basic inputs followed by X behavior-extended inputs. The X behavior-extended arguments are subjugated to a res, and the result is then degraded using the first \(Y\) arguments as fallback. The effect is similar to "ignoring" a the result of a res function if \(\in b\).
The main application of this pattern is as extended behavior wrapper for state machine functions which do not "understand" extended behavior semantics, i.e. it simply propagates the current state \((\in \alpha)\) if the inputs (their res) belongs to the set of extended values \((\in b)\).
Constructors: ignore[1-4] [1-4].

\subsection*{4.4 ForSyDe.Atom.ExB.Absent}
```

module ForSyDe.Atom.ExB.Absent (
AbstExt(Abst, Prst)
) where

```

This module implements the constructors and assocuated utilities of a type which extends the behavior of a function to express "absent events" (see Halbwachs et al., 1991)

The AbstExt type can be used directly with the atom patterns defined in ForSyDe.Atom.ExB, and no helpers or utilities are needed. Example usage:
```

\lambda> res21 (+) (Prst 1) (Prst 2)
3
\lambda> res21 (+) Abst Abst
\lambda> filter Abst (Prst 1)
\perp
\lambda> filter (Prst False) (Prst 1)
\perp
\lambda> filter (Prst True) (Prst 1)
1
\lambda> filter' False 1 :: AbstExt Int
\lambda> filter' True 1 :: AbstExt Int
1
\lambda> degrade 0 (Prst 1)
\lambda> degrade 0 Abst
O
\lambda> ignore11 (+) 1 (Prst 1)
2
\lambda> ignore11 (+) 1 Abst
1
\lambda> res21 (+) (Prst 1) Abst
*** Exception: [ExB.Absent] Illegal occurrence of an absent and present event

```

\section*{data AbstExt a}
\(=\) Abst \(\perp\) denotes the absence of a value
1 Prst a \(T\) a present event with a value
The AbstExt type extends the base type with the ' \(\perp\) ' symbol denoting the absence of a value/event (see Halbwachs et al., 1991).

\section*{instance Functor AbstExt}

Functor instance. Bypasses the special values and maps a function to the wrapped value. Provides the (<\$>) operator.
```

instance Applicative AbstExt

```

Applicative instance. Check source code for the lifting rules. Provides the (<*>) operator

\section*{instance ExB AbstExt}

Implements the absent semantics of the extended behavior atoms.
```

instance Eq a => Eq (AbstExt a)

```
instance Read a \(=>\) Read (AbstExt a)

Reads the '_' character to an Abst and a normal value to Prst-wrapped one.
```

instance Show a => Show (AbstExt a)

```

Shows Abst as ' \(\perp\) ', while a present event is represented with its value.
```

instance (Show a, Plottable a) => Plottable (AbstExt a)

```

Absent-extended plottable types

\subsection*{4.5 ForSyDe.Atom.MoC}


This module exports the core entities of the MoC layer: interfaces for atoms and process constructors as patterns of atoms. It does NOT export any implementation or instantiation of any specific MoC.

Current MoC implementations can be used by importing their respective modules:
- ForSyDe.Atom.MoC.CT
- ForSyDe.Atom.MoC.DE
- ForSyDe.Atom.MoC.SY
- ForSyDe.Atom.MoC.SDF

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.

\subsection*{4.5.1 Atoms}

\section*{class Applicative e => MoC e where}

This is a type class defining interfaces for the MoC layer atoms. Each model of computation exposes its tag system through a unique event constructor as an instance of this class, which defines \(T \times V\).
To express all possible MoCs which can be described using a tagged signal model we need to capture the most general form of their atoms. Recall that all atoms in the layered framework are represented as higher-order functions on structured types (instances of this class), taking functions of other (lower) layers as arguments. While this principle stands also for this layer, the functions taken as arguments need to be formatted for each MoC in particular in order to capture additional information, which we can call in general terms as the execution context.
One typical example of additional information is the consumption and production rates of for data flow MoCs (e.g. SDF). In this case the passed functions are defined over "partitions" of events, i.e. groupings of events with the same partial order in relation to, for example, a
process firing. The formal description of such a "formatted function" taken as argument by a MoC entity is:
\[
\dot{a} \rightarrow \dot{b} \equiv \dot{\alpha}_{1} \times \dot{\alpha}_{2} \times \ldots \rightarrow \dot{\beta}_{1} \times \dot{\beta}_{2} \times \ldots
\]
where \(a\) and \(b\) might be Cartesian products of different types, corresponding to how many signals the constructor is applied to or how many signals it yields, and each type is expressed as:
\[
\begin{aligned}
& \dot{\alpha}=\alpha^{r} \\
& r \in \begin{cases}\{1\}, & \text { for all timed MoCs (e.g. SY, DE, CT) } \\
\mathbb{N}, & \text { for static variants of SDF } \\
\mathbb{N}^{n}, & \text { for CSDF } \\
S \times \mathbb{N} \rightarrow \mathbb{N}, & \text { where } S \text { is a state space, } \\
& \text { in the most general case of untimed data flow }\end{cases}
\end{aligned}
\]

While, as you can see above, the execution context can be extracted from the type information, working with type-level parameters is not a trivial task in Haskell, especially if we want to describe a general and extensible type class. This is why we have chosen a pragmatic approach in implementing the MoC class:
- any (possible) Cartesian product of \(\alpha\) is represented using a recursive type, namely a list \([\alpha]\).
- as the execution context cannot (or can hardly) be extracted from the recursive type, in the most general case we pass both context and argument as a pair (see each instance in particular). To aid in pairing contexts with each argument in a function, the ctxt utilities are provided (see ctxt22).
- this artifice was masked using the generic type families Fun and Res.

\section*{Methods}

Fun e a b Source
This is a type family alias for a context-bound function passed as an argument to a MoC atom. In the most simple case it can be regarded as an enhanced \(->\) type operator. While hiding the explicit definition of arguments, this implementation choice certainly has its advantages in avoiding unnecessary or redundant type constructors (see version 0.1.1 and prior). Aliases are replaced at compile time, thus not affecting run-time performance.

Ret e b Source
\[
\dot{\alpha} \rightarrow \beta
\]

As with Fun, this alias hides a context-bound value (e.g. function return). Although the definition seems to be redundant with Fun, this alias is needed for utilities to recreate clean types again (see -*).
\[
\dot{\beta}
\]
(-.-) :: Fun e a b -> Stream (e a) -> Stream (e b) Source
This atom is mapping a function on values (in the presence of a context) to a signal, i.e. stream of tagged events. As ForSyDe deals with determinate, functional processes, this atom defines the (only) behavior of a process in rapport to one input signal Lee and Sangiovanni-Vincentelli, 1998.
\[
\odot:(\dot{\alpha} \rightarrow \beta) \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\beta)
\]
(-*-) : : Stream (e (Fun e a b)) -> Stream (e a) -> Stream (e b) Source
This atom synchronizes two signals, one carrying functions on values (in the presence of a context), and the other containing values, during which it applies the former on the latter. As concerning the process created, this atom defines a relation between two signals Lee and Sangiovanni-Vincentelli, 1998.
\[
\circledast: \mathcal{S}(\dot{\alpha} \rightarrow \beta) \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\beta)
\]
(-*) : : Stream (e (Ret e b)) -> Stream (e b) Source
Artificial utility which drops the context and/or partitioning yielding a clean signal type.
\[
\mathcal{S}(\dot{\beta}) \rightarrow \mathcal{S}(\beta)
\]
(-<-) : : Stream (e a) -> Stream (e a) -> Stream (e a) Source
This atom appends a (partition of) events at the beginning of a signal. This atom is necessary to ensure complete partial order of a signal and assures the least upper bound necessary for example in the evaluation of feedback loops Lee and SangiovanniVincentelli, 1998.
\[
\theta: \dot{\alpha} \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\alpha)
\]

Notice the difference between the formal and the implemented type signatures. In the implementation the value/partition is wrapped inside an event type to enable smooth composition. You might also notice the type for the initial event(s) as being wrapped inside a signal constructor. This allows defining an DSL for this layer which is centered around signals exclusively, while also enabling to define atoms as homomorphisms to certain extent R. Bird and Moor, 1997. Certain MoCs might have additional constraints on the first operand to be finite.
(-\&-) : : Stream (e a) -> Stream (e a) -> Stream (e a) Source
This atom allows the manipulation of tags in a signal in a restrictive way which preserves monotonicity and continuity in a process Lee and Sangiovanni-Vincentelli, 1998, namely by phase-shifting all tags in a signal with the appropriate metric corresponding to each MoC. Thus it preserves the characteristic function intact I. Sander and Jantsch, 2006.
\[
\oplus: T \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\alpha)
\]

As with the -<- atom, we can justify the type signature for smooth composition and the definition of atoms as homomorphisms to certain extent. This in turn allows the interpretation of the -Methods- operator as aligning the phases of two signals: the second operand is aligned based on the first.
```

instance MoC SY

```

Implenents the execution and synchronization semantics for the SY MoC through its atoms.
```

instance MoC SDF

```

Implenents the SDF semantics for the MoC atoms

\section*{instance MoC DE}

Implenents the execution and synchronization semantics for the DE MoC through its atoms.
```

instance MoC CT

```

Implenents the execution and synchronization semantics for the CT MoC through its atoms.

\subsection*{4.5.2 Process constructors}

Process constructors are defined as patterns of MoC atoms. Check the naming convention of the API in the page description.
delay : : MoC e => Stream (e a) -> Stream (e a) -> Stream (e a)
\[
\begin{aligned}
& \theta: T \times \dot{\alpha} \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\alpha)
\end{aligned}
\]

The delay process provides both initial token(s) and shifts the phase of the signal. In other words, it "delays" a signal with one or several events.

There is also an infix variant -\&>- (infixl 3 ). To justify the first argument, see the documentation of the -<- atom.
comb22
```

: MoCe
$\Rightarrow$ Fun e a1 (Fun e a2 (Ret e b1, Ret e b2)) combinational function (*)
-> Stream (e a1)
-> Stream (e a2)
-> (Stream (e b1), Stream (e b2))
first input signal
second input signal
two output signals

```
\(\left.{ }^{*}\right)\) to be read a1 -> a2 -> (b1, b2) where each argument and result might be individually wrapped with a context and might also express a partition.
\[
\begin{aligned}
& \oplus:\left(\dot{V}^{n} \rightarrow \dot{V}^{m}\right) \rightarrow \mathcal{S}^{n} \rightarrow \mathcal{S}^{m} \\
& f \oplus\left(s_{1}, s_{2}, \ldots s_{n}\right)=f \odot s_{1} \circledast s_{2} \circledast \ldots \circledast s_{n} \triangleleft
\end{aligned}
\]


The comb processes takes care of synchronization between signals and maps combinatorial functions on their event values.

This library exports constructors of type comb [1-8] [1-4].

\section*{reconfig22}
:: \(\quad \mathrm{MoCe}\)
\(\Rightarrow\) Stream (e (Fun e a1 (Fun e a2 (Ret e b1, Ret e signal carrying functions (*) b2))))
-> Stream (e a1) first input signal
-> Stream (e a2) second input signal
-> (Stream (e b1), Stream (e b2)) two output signals
\({ }^{*}\) ) to be read a1 -> a2 -> (b1, b2) where each argument and result might be individually wrapped with a context and might also express a partition.
\[
\begin{aligned}
& \operatorname{reconfig}_{M}: \mathcal{S}\left(\dot{V}^{n} \rightarrow \dot{V}^{m}\right) \rightarrow \mathcal{S}^{n} \rightarrow \mathcal{S}^{m} \\
& \operatorname{reconfig}_{M}\left(s f, s_{1}, s_{2}, \ldots s_{n}\right)=f \odot s_{1} \circledast s_{2} \circledast \ldots \circledast s_{n} \triangleleft
\end{aligned}
\]


The reconfig processes constructs adaptive processes, where the first signal carries functions, and it is synchronized with all the other signals.

This library exports constructors of type reconfig[1-8] [1-4].
```

:: MoC e
=> Fun e st1 (Fun e st2 (Fun e a1 (Fun e a2 (Ret e next state function (*)
st1, Ret e st2))))
-> (Stream (e st1), Stream (e st2))
-> Stream (e a1)
-> Stream (e a2)
-> (Stream (e st1), Stream (e st2))

```
initial state(s) (**)
first input signal
second input signal
output signals mirroring the
next state(s).
(*) meaning st1 -> st2 -> a1 -> a2 -> (st1,st2) where each argument and result might be individually wrapped with a context and might also express a partition.
\(\left.{ }^{* *}\right)\) see the documentation for -<- for justification of the type
\[
\begin{aligned}
& \operatorname{state}_{M}:\left(\dot{V}_{s t}^{m} \times \dot{V}^{n} \rightarrow \dot{V}_{s t}^{m}\right) \times \dot{E}_{s t}^{m} \rightarrow \mathcal{S}^{n} \rightarrow \mathcal{S}_{s t}^{m} \\
& \operatorname{state}_{M}\left(n s, i^{m}\right)\left(s^{n}\right)=n s \oplus\left(s t^{m}, s^{n}\right) \\
& \quad \text { where } s t^{m}=i^{m} \otimes n s \oplus\left(s t^{m}, s^{n}\right)
\end{aligned}
\]


The state processes generate process networks corresponding to a simple state machine like in the graph above.

This library exports constructors of type state [1-4] [1-4].
```

stated22

```
\[
:: \quad \mathrm{MoCe}
\]
\[
\Rightarrow \text { Fun e st1 (Fun e st2 (Fun e a1 (Fun e a2 (Ret e next state function (*) }
\]
st1, Ret e st2))))
\[
->\quad\left(\text { Stream } \left(\text { e st1), Stream (e st2)) } \quad \text { initial state(s) }\left(^{* *}\right)\right.\right.
\]
-> Stream (e a1)
first input signal
-> Stream (e a2)
second input signal
-> (Stream (e st1), Stream (e st2))
output signals mirroring the next state(s).
\(\left.{ }^{*}\right)\) meaning st1 -> st2 \(\rightarrow>\) a1 -> a2 -> (st1,st2) where each argument and result might be individually wrapped with a context and might also express a partition.
(**) see the documentation for -<- for justification of the type

The state processes generate process networks corresponding to a simple state machine like in the graph above. The difference between state22 and stated22 is that the latter outputs the current state rather than the next one. There exists a variant with 0 input signals, in which case the process is a signal generator.

This library exports constructors of type stated [0-4] [1-4].
\[
\begin{aligned}
& \operatorname{stated}_{M}:\left(\dot{V}_{s t}^{m} \times \dot{V}^{n} \rightarrow \dot{V}_{s t}^{m}\right) \times \dot{E}_{s t}^{m} \rightarrow \mathcal{S}^{n} \rightarrow \mathcal{S}_{s t}^{m} \\
& \operatorname{stated}_{M}\left(n s, i^{m}\right)\left(s^{n}\right)=s t^{m} \\
& \text { where } s t^{m}=i^{m} \otimes n s \oplus\left(s t^{m}, s^{n}\right)
\end{aligned}
\]
moore22
:: MoCe
\begin{tabular}{lll}
\(\Rightarrow\) & Fun e st (Fun e a1 (Fun e a2 (Ret e st))) & next state function \(\left(^{*}\right)\) \\
\(->\) & Fun e st (Ret e b1, Ret e b2) & output decoder (**) \\
-> & Stream (e st) & initial state \(\left(^{* * *}\right)\) \\
-> & Stream (e a1) & first input signal \\
\(->\) & Stream (e a2) & second input signal \\
\(->\) & (Stream (e b1), Stream (e b2)) & output signals
\end{tabular}
(*) meaning st -> a1 -> a2 -> st where each argument and result might be individually wrapped with a context and might also express a partition.
\({ }^{(* *)}\) meaning st -> (b1, b2) where each argument and result might be individually wrapped with a context and might also express a partition.
\(\left(^{* * *}\right)\) see the documentation for \(-<-\) for justification of the type
\[
\begin{aligned}
& \operatorname{moore}_{M}:\left(\dot{V}_{s t} \times \dot{V}^{n} \rightarrow \dot{V}_{s t}\right) \times\left(\dot{V}_{s t} \rightarrow \dot{V}^{m}\right) \times \dot{E}_{s t} \rightarrow \mathcal{S}^{n} \rightarrow \mathcal{S}^{m} \\
& \operatorname{moore}_{M}\left(n s, o d, i^{m}\right)\left(s^{n}\right)=o d \oplus s t^{m} \\
& \text { where } s t^{m}=i^{m} \otimes n s \oplus\left(s t^{m}, s^{n}\right)
\end{aligned}
\]

The moore processes model Moore state machines.
This library exports constructors of type moore [1-4] [1-4].

\section*{mealy22}
:: MoCe
\(\Rightarrow\) Fun e st (Fun e a1 (Fun e a2 (Ret e st))) next state function (*)
\(\rightarrow\) Fun e st (Fun e a1 (Fun e a2 (Ret e b1, Ret e b2))) output decoder (**)
-> Stream (e st)
-> Stream (e a1)
-> Stream (e a2)
(Stream (e b1), Stream (e b2)) initial state ( \({ }^{* * *}\) ) first input signal second input signal output signals
(*) meaning st -> a1 -> a2 -> st where each argument and result might be individually wrapped with a context and might also express a partition.
(**) meaning st -> a1 -> a2 -> (b1, b2) where each argument and result might be individually wrapped with a context and might also express a partition.
\(\left(^{* * *}\right)\) see the documentation for \(-<-\) for justification of the type
\[
\begin{aligned}
& \text { mealy }_{M}:\left(\dot{V}_{s t} \times \dot{V}^{n} \rightarrow \dot{V}_{s t}\right) \times\left(\dot{V}_{s t} \times \dot{V}^{n} \rightarrow \dot{V}^{m}\right) \times \dot{E}_{s t} \rightarrow \mathcal{S}^{n} \rightarrow \mathcal{S}^{m} \\
& \text { mealy }{ }_{M}\left(n s, o d, i^{m}\right)\left(s^{n}\right)=\operatorname{od} \oplus\left(s t^{m}, s^{n}\right) \\
& \quad \text { where } s t^{m}=i^{m} \otimes n s \oplus\left(s t^{m}, s^{n}\right)
\end{aligned}
\]


The mealy processes model Mealy state machines.
This library exports constructors of type mealy [1-4] [1-4].

\subsection*{4.5.3 Utilities}
```

ctxt22
-> (ctx, ctx)
-> (a1 -> a2 -> (b1, b2))
-> (ctx, a1 -> (ctx, a2 -> ((ctx, b1),
(ctx, b2))))

```
    \(:\) (ctx, ctx) argument contexts (e.g. consumption
        rates in SDF)
        result contexts (e.g. production rates in
        SDF)
        function on values/partitions of values
        context-wrapped form of the previous
        function
        ctxt \(: \Gamma \times\left([\alpha]^{n} \rightarrow[\beta]^{m}\right) \rightarrow\left(\dot{\alpha}^{n} \rightarrow \dot{\beta}^{m}\right)\)

Wraps a function with the context needed by some MoCs for their constructors (e.g. rates in SDF).
This library exports wrappers of type ctxt[1-8] [1-4].
warg :: c -> (a -> b) -> (c, a -> b)
Attaches a context parameter to a function argument (e.g consumption rates in SDF). Used as kernel function in defining e.g. ctxt22.
```

wres :: p -> b -> (p, b)

```

Attaches a context parameter to a function's result (e.g production rates in SDF). Used as kernel function in defining e.g. ctxt22.
```

(-*<) :: MoC e =>
Stream (e (Ret e b1, Ret e b2)) -> (Stream (e b1), Stream (e b2))

```

Utilities for extending the -* atom for dealing with tupled outputs. This library exports operators of form \(-*<\{1,8\}\).

\subsection*{4.6 ForSyDe.Atom.MoC.Stream} module ForSyDe.Atom.MoC.Stream (

Stream(NullS, (:-)), , , , stream, fromStream, headS, tailS,
lastS, repeatS, takeS, dropS, takeWhileS, (+-+)
where

This module defines the shallow-embedded Stream datatype and utility functions operating on it. In ForSyDe a signal is represented as a (partially or totally) ordered sequence of events that enables processes to communicate and synchronize. The Stream type is but an ordered structure to encapsulate events as infinite streams.

\section*{data Stream e}
\[
\begin{array}{ll}
=\text { Nulls } & \text { terminates a signal } \\
\text { I e (:-) (Stream e) } & \text { the default constructor appends an } \\
& \text { event to the head of the stream }
\end{array}
\]

Defines a stream of events, encapsulating them in a structure isomorphic to an infinite list R. S. Bird, 1987, thus all properties of lists may also be applied to Streams. While, in combination with lazy evaluation, it is possible to create and simulate infinite signals, we need to ensure that the first/previous event is always fully evaluated. This can be translated into the following rule:
zero-delay feedbacks are forbidden, due to un-evaluated self-referential calls. In a feedback loop, there always has to be enough events to ensure the data flow.

This rule imposes that the stream of data is uninterrupted in order to have an evaluatable kernel every time a new event is produced (i.e. to avoid deadlocks). Thus we can add the rule:
cleaning of output events is also forbidden. In other words, for each new input at any instant in time, a process must react with at least one output event.

\section*{instance Functor Stream}
allows for the mapping of an arbitrary function ( \(\mathrm{a}-\mathrm{>}\) b) upon all the events of a (Stream a).
```

instance Applicative Stream

```
enables the Stream to behave like a ZipList
instance Foldable Stream
provides folding functions useful for implementing utilities, such as length.
instance Read a => Read (Stream a)
signal (1 :- 2 :- NullS) is read using the string "\{1,2\}".
instance Show a => Show (Stream a)
signal (1 :- \(2:-\mathrm{NullS}\) ) is represented as \(\{1,2\}\).
```

instance Plottable a => Plot (Signal a)
SY signals.
instance Plottable a => Plot (Signal a)
SDF signals.
instance Plottable a => Plot (Signal a)
DE signals.
instance Plottable a => Plot (Signal a)
CT signals.

```
stream :: [a] -> Stream a

The function signal converts a list into a signal.
```

fromStream :: Stream a -> [a]

```

The function fromStream converts a signal into a list.
headS : : Stream a -> a

Returns the head of a signal.
```

tailS :: Stream e -> Stream e

```

Returns the tail of a signal
```

lastS :: Stream p -> p

```

Returns the last event in a signal.
```

repeatS :: a -> Stream a

```

Returns an infinite list containing the same repeated event.
```

takeS :: (Ord t, Num t) => t -> Stream e -> Stream e

```

Returns the first n events in a signal.
dropS :: (Ord t, Num t) => t -> Stream e -> Stream e
Drops the first n events in a signal.
takeWhileS :: (a -> Bool) -> Stream a -> Stream a
Returns the first events of a signal which comply to a condition.
(+-+) :: Stream e -> Stream e -> Stream e
Concatenates two signals.

\subsection*{4.7 ForSyDe.Atom.MoC.SY}
```

module ForSyDe.Atom.MoC.SY (
SY(SY, val), Signal, unit2, signal, readSignal, delay, comb22,
reconfig22, constant2, generate2, stated22, state22, moore22,
mealy22, when, when', is, whenPresent, filter, filter', fill, hold,
reactAbst2, toDE2, toSDF2, zipx, unzipx, unzipx'
) where

```

The sy library implements the atoms holding the sematics for the synchronous computation model. It also provides a set of helpers for properly instantiating process network patterns as process constructors.

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.

\subsection*{4.7.1 Synchronous (sy) event}

According to Lee and Sangiovanni-Vincentelli, 1998, "two events are synchronous if they have the same tag, and two signals are synchronous if all events in one signal are synchronous to an event in the second signal and vice-versa. A system is synchronous if every signals in the system is synchronous to all the other signals in the system."

The synchronous (Sy) MoC defines no notion of physical time, its tag system suggesting in fact the precedence among events. To further demystify its mechanisms, we can formulate the following proposition:

The SY MoC is abstracting the execution semantics of a system where computation is assumed to perform instantaneously (with zero delay), at certain synchronization points, when data is assumed to be available.

Below is a possible behavior in time of the input and the output signals of a SY process, to illustrate these semantics:


Implementing the SY tag system is straightforward if we consider the synchronous Signal as an infinite list. In this case the tags are implicitly defined by the position of events in a signal: \(t_{e}\) would correspond with the event at the head of a signal \(t_{1}\) with the following event, etc... The only explicit parameter passed to a SY event constructor is the value it carries \(\in V_{e}\). As such, we can state the following particularities:
1. tags are implicit from the position in the Stream, thus they are completely ignored in the type constructor.
2. the type constructor wraps only a value
3. being a timed \(M o C\), the order between events is total Lee and Sangiovanni-Vincentelli, 1998.
4. from the previous statement we can conclude that there is no need for an execution context (see section 4.5) and we can ignore the formatting of functions in ForSyDe.Atom.MoC, thus we can safely assume:
\[
\dot{a}=a
\]
newtype SY a
\(=\) SY
> val :: a
The SY event. It identifies a synchronous signal.
instance Functor SY
Allows for mapping of functions on a SY event.
instance Applicative SY
Allows for lifting functions on a pair of SY events.
instance MoC SY
Implenents the execution and synchronization semantics for the SY MoC through its atoms.
```

instance Read a => Read (SY a)

```

Reads the value wrapped
instance Show a => Show (SY a)
Shows the value wrapped
instance Plottable \(a=>\) Plot (Signal a)
SY signals.
instance type Ret \(S Y\) b \(=\mathrm{b}\)
instance type Fun SY a b \(=\mathrm{a}->\mathrm{b}\)

\subsection*{4.7.2 Aliases \& utilities}

A set of type synonyms and utilities are provided for convenience. The API type signatures will feature these aliases to hide the cumbersome construction of atoms and patters as seen in ForSyDe.Atom.MoC.
```

type Signal a = Stream (SY a)

```

Type synonym for a SY signal, i.e. "an ordered stream of SY events"
unit2 :: (a1, a2) -> (Signal a1, Signal a2)
Wraps a (tuple of) value(s) into the equivalent unit signal(s).
Helpers: unit, unit2, unit3, unit4.
signal :: [a] -> Signal a
Transforms a list of values into a SY signal.
readSignal :: Read a => String -> Signal a
Reads a signal from a string. Like with the read function from Prelude, you must specify the tipe of the signal.
```

|> readSignal "{1,2,3,4,5}" :: Signal Int

```
\(\{1,2,3,4,5\}\)

\subsection*{4.7.3 sy process constuctors}

These SY process constructors are basically specific instantiations of the patterns of atoms defined in ForSyDe.Atom.MoC. Some are also wrapping functions in an extended behavioural model.

\section*{Simple}
delay
\[
\begin{array}{lll}
:: & \text { a } & \text { initial value } \\
\text {-> } & \text { Signal a } & \text { input signal } \\
\text {-> } & \text { Signal a } & \text { output signal }
\end{array}
\]

The delay process "delays" a signal with one event. Instantiates the delay pattern.
```

|> let s = signal [1,2,3,4,5]
\lambda> delay 0 s
{0,1,2,3,4,5}

```
comb22
    \(:: \quad(\mathrm{a} 1->\mathrm{a} 2->(\mathrm{b} 1, \mathrm{~b} 2))\)
    -> Signal a1 first input signal
        alue
    -> Signal a2 second input signal
    (Signal b1, Signal b2) two output signals
    comb processes map combinatorial functions on signals and take care of synchronization
    between input signals. It instantiates the comb pattern (see comb22).
    Constructors: comb [1-4] [1-4].
    \(\lambda>\) let s1 = signal [1..]
    \(\lambda>\) let s2 \(=\) signal \([1,1,1,1,1]\)
\(\lambda>\) comb11 (+1) s2
\(\{2,2,2,2,2\}\)
\(\lambda>\) comb22 ( \(\backslash \mathrm{a}\) b-> (a+b, a-b)) s1 s2
( \(\{2,3,4,5,6\},\{0,1,2,3,4\}\) )

reconfig22
    :: Signal (a1 -> a2 -> (b1, b2)) signal carrying functions
    -> Signal a1
    first input signal carrying arguments
    -> Signal a2
    second input signal carrying arguments
    -> (Signal b1, Signal b2)
                                two output signals
reconfig creates an synchronous adaptive process where the first signal carries functions and the other carry the arguments. It instantiates the reconfig atom pattern (see reconfig22).
Constructors: reconfig[1-4] [1-4].
```

\lambda> let sf = signal [(+1),(*2),(+1),(*2),(+1),(*2),(+1)]
\lambda> let s1 = signal [1..]
\lambda> reconfig11 sf s1
{2,4,4,8,6,12,8}

```

constant2
\[
\begin{array}{lll}
:: & (b 1, b 2) & \text { values to be repeated } \\
-> & (\text { Signal b1, Signal b2) } & \text { generated signals }
\end{array}
\]

A signal generator which keeps a value constant. It is actually an instantiation of the statedoX constructor (check stated22).
Constructors: constant [1-4].
```

\lambda> let (s1, s2) = constant2 (1,2)
\lambda> takeS 3 s1
{1,1,1}
\lambda> takeS 5 s2
{2,2,2,2,2}

```

\section*{generate2}
:: (b1 -> b2 -> (b1, b2)) function to generate next value
-> (b1, b2) kernel values
-> (Signal b1, Signal b2) generated signals
A signal generator based on a function and a kernel value. It is actually an instantiation of the statedoX constructor (check stated22).

Constructors: generate [1-4].
\(\lambda>\) let \((s 1, s 2)=\) generate2 \((\backslash a b->(a+1, b+2))(1,2)\)
\(\lambda>\) takeS 5 s1
\(\{1,2,3,4,5\}\)
入> takeS 7 s2
\(\{2,4,6,8,10,12,14\}\)
stated22
\begin{tabular}{lll}
\(::\) & \((b 1->b 2->~ a 1 ~\) & a2 \(->(b 1, b 2))\) \\
-> & next state function \\
-> b2) & Signal a1 & \\
initial state values
\end{tabular}
stated is a state machine without an output decoder. It is an instantiation of the state MoC constructor (see stated22).
Constructors: stated [1-4] [1-4].
```

\lambda> let s1 = signal [1,2,3,4,5]
\lambda> stated11 (+) 1 s1
{1,2,4,7,11,16}

```

state22

state is a state machine without an output decoder. It is an instantiation of the stated MoC constructor (see state22).

Constructors: state[1-4] [1-4].
\(\lambda>\) let \(s 1=\) signal \([1,2,3,4,5]\)
\(\lambda>\) state11 (+) 1 s1
\(\{2,4,7,11,16\}\)
moore22
:: (st -> a1 -> a2 -> st) next state function
-> (st -> (b1, b2)) output decoder
-> st initial state
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)
moore processes model Moore state machines. It is an instantiation of the moore MoC constructor (see moore22).

Constructors: moore [1-4] [1-4].

mealy processes model Mealy state machines. It is an instantiation of the mealy MoC constructor (see mealy22).
Constructors: mealy [1-4] [1-4].
```

\lambda> let s1 = signal [1,2,3,4,5]
\lambda> mealy11 (+) (-) 1 s1
{0,0,1,3,6}

```


\section*{Predicate behavior}

These processes manipulate the behavior of a signal based on predicates on their status.
```

when
:: Signal (AbstExt Bool) Signal of predicates
-> Signal (AbstExt a) Input signal
-> Signal (AbstExt a) Output signal

```

This process predicates the existence of values in a signal based on a signal of boolean values (conditions). It is similar to the when construct in the synchronous language Lustre Halbwachs et al., 1991, based on which clock calculus can be performed.

OBS: this process assumes that all signals carry absent-extended values, which is appropriate in describing multi-clock systems. For a version which inputs signals of non-extended values, check when'
```

\lambda> let s1 = (signal . map Prst) [1,2,3,4,5]
\lambda> let pred = (signal . map Prst) [False,False,False,True,True]
\lambda> when pred s1
{\perp,\perp,\perp,4,5}

```
when'
\begin{tabular}{lll}
\(::\) & Signal Bool & Signal of predicates \\
\(->\) & Signal a & Input signal \\
\(->\) & Signal (AbstExt a) & Output signal
\end{tabular}

Same as when but inputs signals of non-extended values.
```

\lambda> let s1 = signal [1,2,3,4,5]
\lambda> let pred = signal [False,False,False,True,True]
\lambda> when' pred s1
{\perp,\perp,\perp,4,5}

```
is : : Signal (AbstExt a) -> (a -> Bool) -> Signal (AbstExt Bool)

Simple wrapper for applying a predicate function on a signal of absent-extended events.
\(\lambda>\) let s1 = signal \$ map Prst [1, 2, 3, 4,5]
\(\lambda>\) s1 'is' (>3)
\{False, False, False, True, True\}
whenPresent :: Signal (AbstExt a1)
```

-> Signal (AbstExt a2) -> Signal (AbstExt a2)

```

Same as when but triggering the output events merely based on the presence of the first input rather than a predicate function.
```

\lambda> let s1 = signal \$ map Prst [1,2,3,4,5]
\lambda> let sp = signal [Prst 1, Prst 1, Abst, Prst 1, Abst]
\lambda> whenPresent sp s1
{1,2,\perp,4,\perp}

```
filter
\begin{tabular}{lll} 
:: & (a -> Bool) & Predicate function \\
-> & Signal (AbstExt a) & Input signal \\
-> & Signal (AbstExt a) & Output signal
\end{tabular}

Filters out values to Abst if they do not fulfill a predicate function.
OBS: this process assumes that all signals carry absent-extended values, which is appropriate in describing multi-clock systems. For a version which inputs signals of non-extended values, check filter'.
```

\lambda> let s1 = (signal . map Prst) [1,2,3,4,5]
\lambda> filter (>3) s1
{\perp,\perp,\perp,4,5}

```

filter'
:: (a -> Bool) Predicate function
-> Signal a Input signal
-> Signal (AbstExt a) Output signal

Same as filter but inputs signals of non-extended values.
```

\lambda> let s1 = signal [1,2,3,4,5]
\lambda> filter' (>3) s1
{\perp,\perp,\perp,4,5}

| $::$ | a | Value to fill with |
| :--- | :--- | :--- |
| $->$ | Signal (AbstExt a) | Input |
| -> | Signal a | Output |

```
fill

Fills absent events with a pre-defined value.
```

\lambda> let s1 = signal [Abst, Abst, Prst 1, Prst 2, Abst, Prst 3]
\lambda> fill 0 s1
{0,0,1,2,0,3}

```
hold
\(: \quad \mathrm{a}\)

Value to fill with in case there was no previous value
-> Signal (AbstExt a) Input
\(\rightarrow\) Signal a Output
Similar to fill, but holds the last non-absent value if there was one. It implements a state pattern (see state22).
\(\lambda>\) let s1 = signal [Abst, Abst, Prst 1, Prst 2, Abst, Prst 3]
\(\lambda>\) hold 0 s1
\(\{0,0,1,2,2,3\}\)

reactAbst2
\begin{tabular}{lll}
\(::\) & \begin{tabular}{l} 
(Signal (AbstExt a1) -> Signal
\end{tabular} & \begin{tabular}{l} 
process which degrades the absent extension, \\
(AbstExt a2) -> Signal b)
\end{tabular} \\
\(->\) & Signal (AbstExt a1) & \\
\(->\) Signal (AbstExt a2) & \\
-> Signal (AbstExt b) & \begin{tabular}{l} 
absent-extended signal, properly reacting to \\
the inputs
\end{tabular}
\end{tabular}

Creates a wrapper enabling a reactive behvior to absent-extended signals for processes which would otherwise degrade the absent-extension (e.g. state machines with ignore22 behavior).
Constructors: reactAbst[1-4].
```

\lambda> let s1 = readSignal "{1,1,1,_,1,_,1}" :: Signal (AbstExt Int)
\lambda> let proc = stated11 (B.ignore11 (+)) 0
\lambda> proc s1
{0,1,2,3,3,4,4,5}
\lambda> reactAbst1 proc s1
{0,1,2,\perp,3,\perp,4}

```


\subsection*{4.7.4 Interfaces}
toDE2
\begin{tabular}{lll} 
:: & Signal TimeStamp & SY signal carrying DE timestamps \\
-> & Signal a & first input SY signal \\
-> & Signal b & second input SY signal \\
-> & (Signal a, Signal b) & two output DE signals
\end{tabular}

Wraps explicit timestamps to a (set of) SY signal(s), rendering the equivalent synchronized DE signal(s).

Constructors: toDE, toDE2, toDE3, toDE4.
```

\lambda> let s1 = SY.signal [0,3,4,6,9]
\lambda> let s2 = SY.signal [1,2,3,4,5]
\lambda> toDE s1 s2
{ 1 @0s, 2 @3s, 3 @4s, 4 @6s, 5 @9s}

```
toSDF2 :: Signal a -> Signal b -> (Signal a, Signal b)
Transforms a (set of) SY signal(s) into the equivalent SDF signal(s). The only change is the event consructor. The total order of SY is interpreted as partial order by the next SDF process downstream.
Constructors: toSDF, toSDF2, toSDF3, toSDF4.
\(\lambda>\) let \(s=\) SY.signal \([1,2,3,4,5]\)
\(\lambda>\) toSDF s
\(\{1,2,3,4,5\}\)

zipx :: Vector (Signal a) -> Signal (Vector a)
Synchronizes all the signals contained by a vector and zips them into one signal of vectors. It instantiates the zipx skeleton.
```

\lambda> let s1 = SY.signal [1,2,3,4,5]
\lambda> let s2 = SY.signal [11,12,13,14,15]
\lambda> let v1 = V.vector [s1,s1,s2,s2]
\lambda> v1
<{1,2,3,4,5},{1,2,3,4,5},{11,12,13,14,15},{11,12,13,14,15}>
\lambda> zipx v1
{<1,1,11,11>,<2,2,12,12>,<3,3,13,13>,<4,4,14,14>,<5,5,15,15>}

```

```

unzipx : : Integer -> Signal (Vector a) -> Vector (Signal a)

```

Unzips the vectors carried by a signal into a vector of signals. It instantiates the unzipx skeleton. To avoid infinite recurrence, the user needs to provide the length of the output vector.
\(\lambda>\) let \(\mathrm{v} 1=\mathrm{V}\).vector \([1,2,3,4]\)
\(\lambda>\) let s1 = SY.signal [v1,v1,v1,v1,v1]
\(\lambda>\mathrm{s} 1\)
\(\{\langle 1,2,3,4\rangle,\langle 1,2,3,4\rangle,\langle 1,2,3,4\rangle,\langle 1,2,3,4\rangle,\langle 1,2,3,4\rangle\}\)
\(\lambda>\) unzipx 4 s1
\(<\{1,1,1,1,1\},\{2,2,2,2,2\},\{3,3,3,3,3\},\{4,4,4,4,4\}>\)

unzipx' : : Signal (Vector a) -> Vector (Signal a)
Same as unzipx, but "sniffs" the first event to determine the length of the output vector. Might have unsafe behavior!
```

\lambda> let v1 = V.vector [1,2,3,4]
\lambda> let s1 = SY.signal [v1,v1,v1,v1,v1]
\lambda> s1
{<1,2,3,4>,<1,2,3,4>,<1,2,3,4>,<1,2,3,4>,<1,2,3,4>}
\lambda> unzipx' s1
<{1,1,1,1,1},{2,2,2,2,2},{3,3,3,3,3},{4,4,4,4,4}>

```

\subsection*{4.8 ForSyDe.Atom.MoC.DE}
module ForSyDe.Atom.MoC.DE (
TimeStamp, DE(DE, tag, val), Signal, unit2, infinite, until, signal, checkSignal, readSignal, delay, delay', comb22, reconfig22, sync2, constant2, generate2, stated22, state22, moore22, mealy22, toSY2, toCT2, zipx, unzipx, unzipx', embedSY22
) where

The DE library implements the atoms holding the sematics for the discrete event computation model. It also provides a set of helpers for properly instantiating process network patterns as process constructors.

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.

\subsection*{4.8.1 Discrete event (DE)}

According to Lee and Sangiovanni-Vincentelli, 1998, "a discrete-event system is a timed system \(Q\) where for all \(s \in Q\), the tag sytem is order-isomorphic to a subset of the integers. Order-isomorphic means simply that there exists an order-preserving bijection between the events in \(T\) and a subset of the integers (or the entire set of integers)."

The discrete event (DE) MoC does suggest the notion of physical time through its tags, also called timestamps. As the definition above implies, an important property of the DE tag system is that between any two timestamps \(t_{u}\) and \(t_{v}\) there is a finite number of possible timestamps. Based on this we can formulate the folowing specialized definition:

The DE MoC is abstracting the execution semantics of a system where synchronization is discretized and time-aware, and it is performed whenever a new event occurs.

There are many variants of discrete event simulators, each of them implementing slight variations of the semantics stated in Lee and Sangiovanni-Vincentelli, 1998. The execution model covered by the DE implementation of ForSyDe-atom may be described as a simplified "cycle simulator" with no delta-delay nor superdense time. The signals behave as "latched channels" (similar to an HDL simulator), and processes react instantaneously to any new event. While the simplicity of the execution engine is a desired one, more complex behaviors such as zero-time feedback, non-zero reaction time and communication protocols (e.g. lossy buffers) may be achieved by composing patterns from the ForSyDe.Atom.MoC and/or ForSyDe.Atom.ExB layers. Nevertheless, the DE behaviors possible within ForSyDe-Atom are included in the class of conservative simulators as presented in Fujimoto00 due to the dataflow nature of the evaluation mechanisms. Below you can see an example of a simple DE process, without behavior extensions:


Below are stated a few particularities of our DE MoC implementation:
1. According to Lee and Sangiovanni-Vincentelli, 1998, our DE MoC is a one-sided system, i.e. time starts from an absolute 0 . While negative time cannot be represented, signals can be phase-aligned with the help of the -\&- atom. All signals need to start from timestamp 0 , and events need to be positioned with their tags in strict ascending order. The checkSignal utility enforces these rules.
2. tags are explicit and a DE event will construct a type around both a tag and a value. Tags represent the start time of the event, the end time being implicit from the start time of the next event. By doing so, we ensure that the time domain is non-disjoint, i.e. a sub-case of continous time.
3. according to the previous point, events are assumed to persist from their time of arrival until the next event arrives or, if there is no incoming event, until infinity. This default behavior can be be interpret signals as being either persistent channels (e.g. latched wires), or non-blocking buffers of size 1 .
4. as a consequence to the previous is that feedback loops will generate an infinite number of events (strictly preceding each other), since a loop updates the value after a certain delay, and any input is assumed to go to infinity. Thus we can now fully jutify the definition of the delay pattern as consisting in a prepend (i.e. generating the new value) and a phase shift (i.e. advancing time with a positive integer). This is done in order to both preserve causality and avoid deadlock.
5. due to the reactive and dataflow natures of the execution system, DE processes are forbidden to clean up events. Doing so might lead to deadlock wherever any feedback is involved. This means that a new event is created every time a new event arrives, regardless of what value it carries. This means that all values are propagated, justifying our system's conservative approach Fujimoto00 Atoms themselves do not clean signals, but using interfaces that do should be treated with extreme special care, as it is considered unsafe and deadlock-prone.
6. due to the conservative approach, and to the fact that MoC atoms are independent synchronization entities, ForSyDe simulators are completely parallelizable, since processes are self-sufficient and do not depend on a global event queue (as compared to other cycle simulators).
7. any signal from outside needs to be sane ( \(T\) must be a total order) before being injected into a ForSyDe process network. Helper functions are equipped with sanity checkers. Inside a ForSyDe process network, transformations are rate-monotonic, thus output signals are guaranteed to be sane.
8. since \(T\) is a total order, there is no need for an execution context (see section 4.5) and we can ignore the formatting of functions in ForSyDe.Atom.MoC, thus we can safely assume:
\[
\dot{a}=a
\]
type TimeStamp = DiffTime
Alias for the type representing discrete time. It is inherently quantizable, the quantum being a picosecond ( \(10^{-12}\) seconds), thus it can be considered order-isomorphic with a set of integers, i.e. between any two timestamps there is a finite number of timestamps. Moreover, a timestamp can be easily translated into a rational number representing fractions of a second, so the conversion between timestamps (discrete time) and rationals (analog/continuous time) is straightforward.

This type is used in the explicit tags of the DE MoC (and subsequently the discrete event evaluation engine for simulating the CT MoC ).
```

data DE a

```
    \(=\mathrm{DE}\)
    \(>\) tag :: TimeStamp timestamp
    > val :: a the value

The DE event. It identifies a discrete event signal.
```

instance Functor DE

```

Allows for mapping of functions on a DE event.
instance Applicative DE
Allows for lifting functions on a pair of DE events.
instance MoC DE
Implenents the execution and synchronization semantics for the DE MoC through its atoms.
```

instance Eq a => Eq (DE a)
instance Read a => Read (DE a)

```

Reads the string of type v@t as an event DE t v.
```

instance Show a => Show (DE a)

```

Shows the event with tag \(t\) and value \(v\) as \(v\) @t.
```

instance Plottable a => Plot (Signal a)

```
    DE signals.
instance type Ret DE b \(=\mathrm{b}\)
instance type Fun DE a b \(=\mathrm{a}->\mathrm{b}\)

\subsection*{4.8.2 Aliases \& utilities}

A set of type synonyms and utilities are provided for convenience. The API type signatures will feature these aliases to hide the cumbersome construction of atoms and patters as seen in ForSyDe.Atom.MoC.
```

type Signal a = Stream (DE a)
Type synonym for a SY signal, i.e. "a signal of SY events"
unit2 :: ((TimeStamp, a1), (TimeStamp, a2))
-> (Signal a1, Signal a2)

```
    Wraps a (tuple of) pair(s) (tag, value) into the equivalent unit signal(s), in this case a
    signal with one event with the period tag carrying value.
    Helpers: unit and unit[2-4].
infinite :: a -> Signal a
    Creates an infinite signal.
until :: TimeStamp -> Signal a -> Signal a
    Takes the first part of the signal util a given timestamp. The last event of the resulting
    signal is at the given timestamp and carries the previous value. This utility is useful when
    plotting a signal, to specify the interval of plotting.
signal :: [(TimeStamp, a)] -> Signal a
    Transforms a list of tuples (tag, value) into a DE signal. Checks if it is well-formed.
checkSignal :: Stream (DE a) -> Stream (DE a)
    Checks if a signal is well-formed or not, according to the DE MoC interpretation in ForSyDe-Atom.
readSignal : : Read a => String -> Signal a
    Reads a signal from a string and checks if it is well-formed. Like with the read function from
    Prelude, you must specify the type of the signal.
    \(\lambda>\) readSignal "\{ 1@0, 2@2, 3@5, 4@7, 5@10 \}" :: Signal Int
    \{ 1 @Os, 2 @2s, 3 @5s, 4 @7s, 5 @10s\}
    \(\lambda>\) readSignal "\{ 1@0, 2@2, 3@5, 4@10, 5@7 \}" :: Signal Int
    \{ 1 @Os, 2 @2s, 3 @5s*** Exception: [MoC.DE] malformed signal
    \(\lambda>\) readSignal "\{ 1@1, 2@2, 3@5, 4@7, 5@10 \}" :: Signal Int
    *** Exception: [MoC.DE] signal does not start from global 0

\subsection*{4.8.3 DE process constuctors}

The DE process constructors are basically specific instantiations of patterns defined in ForSyDe .Atom.MoC. Some might also be wrapping functions in an extended behavioural model.

\section*{Simple}

These are mainly direct instantiations of patterns defined in ForSyDe.Atom.MoC, using DE-specific utilities.
```

delay
:: TimeStamp time delay
-> a initial value
-> Signal a input signal
-> Signal a output signal

```

The delay process "delays" a signal with one event. Instantiates the delay pattern.
```

\lambda> let s = readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: Signal Int
\lambda> delay 3 0 s
{0 @0s, 1 @3s, 2 @5s, 3 @9s, 4 @11s, 5 @12s}

```

delay,
:: Signal a signal "borrowing" the initial event
-> Signal a input signal
-> Signal a output signal
Similar to the previous, but this is the raw instantiation of the delay pattern. It "borrows" the first event from one signal and appends it at the head of another signal.
```

\lambda> let s1 = readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: Signal Int
\lambda> let s2 = readSignal "{3@0, 4@4, 5@5, 6@8, 7@9}" :: Signal Int
\lambda> delay' s1 s2
{1 @0s, 3 @2s, 4 @6s, 5 @7s, 6 @10s, 7 @11s}

```


\section*{comb22}
\[
\begin{array}{lll}
:: & (\mathrm{a} 1->\mathrm{a} 2->(\mathrm{b} 1, \mathrm{~b} 2)) & \text { function on values } \\
-> & \text { Signal a1 } & \text { first input signal } \\
-> & \text { Signal a2 } & \text { second input signal } \\
-> & \text { (Signal b1, Signal b2) } & \text { two output signals }
\end{array}
\]
comb processes map combinatorial functions on signals and take care of synchronization between input signals. It instantiates the comb pattern (see comb22).
Constructors: comb [1-4] [1-4].
```

\lambda> let s1 = infinite 1
\lambda> let s2 = readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: Signal Int
\lambda> comb11 (+1) s2
{ 2 @Os, 3 @2s, 4 @6s, 5 @8s, 6 @9s}
\lambda> comb22 (\a b-> (a+b,a-b)) s1 s2
({ 2 @0s, 3 @2s, 4 @6s, 5 @8s, 6 @9s},{ 0 @0s, -1 @2s, -2 @6s, -3 @8s, -4 @9s})

```

reconfig22
:: Signal (a1 -> a2 -> (b1, b2)) signal carrying functions
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)
first input signal carrying arguments second input signal carrying arguments two output signals
reconfig creates a DE adaptive process where the first signal carries functions and the other carry the arguments. It instantiates the reconfig atom pattern (see reconfig22).
Constructors: reconfig[1-4] [1-4].
```

\lambda> let sf = signal [(0,(+1)),(2,(*2)),(5,(+1)),(7,(*2))]
\lambda> let s1 = signal [(0,1),(3,2),(5,3),(9,4)]
\lambda> reconfig11 sf s1
{ 2 @0s, 2 @2s, 4 @3s, 4 @5s, 6 @7s, 8 @9s}

```

sync2
\begin{tabular}{lll} 
:: & Signal a1 & first input signal \\
-> & Signal a2 & second input signal \\
-> & (Signal a1, Signal a2) & two output signals
\end{tabular}
sync synchronizes multiple signals, so that they have the same set of tags, and consequently, the same number of events. It instantiates the comb atom pattern (see comb22).
Constructors: sync [1-4]
```

\lambda> let s1 = readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: Signal Int
\lambda> let s2 = readSignal "{1@0, 2@5, 3@6, 4@10, 5@12}" :: Signal Int
\lambda> sync2 s1 s2
({ 1 @0s, 2 @2s, 2 @5s, 3 @6s, 4 @8s, 5 @9s, 5 @10s, 5 @12s},{ 1 @0s, 1 @2s, 2 @5s, 3 @6s, 3
@8s, 3 @9s, 4 @10s, 5 @12s})

```

constant2
\[
:: \quad(\mathrm{b} 1, \mathrm{~b} 2) \quad \text { values to be repeated }
\]
-> (Signal b1, Signal b2) generated signals
A signal generator which keeps a value constant. As compared with the Sy, it just constructs an infinite signal with constant value (i.e. a signal with one event starting from time 0 ).
Constructors: constant [1-4].
```

\lambda> constant1 2
{ 2 @Os}

```

\section*{generate2}

\section*{\(:: \quad(\mathrm{b} 1->\mathrm{b} 2\)-> \((\mathrm{b} 1, \mathrm{~b} 2))\)}
((TimeStamp, b1), (TimeStamp, b2) )
-> (Signal b1, Signal b2)
function to generate next value
kernel values tupled with their generation rate.
generated signals

A signal generator based on a function and a kernel value. It is actually an instantiation of the stated0X constructor (check stated22).
Constructors: generate [1-4].
```

\lambda> let (s1,s2) = generate2 (\a b -> (a+1,b+2)) ((3,1),(1,2))
\lambda> takeS 5 s1
{ 1 @Os, 2 @3s, 2 @4s, 2 @5s, 3 @6s}
\lambda> takeS 7 s2
{ 2 @0s, 4 @1s, 6 @2s, 8 @3s, 10 @4s, 12 @5s, 14 @6s}

```
stated22
    :: (b1 -> b2 -> a1 -> a2 -> (b1, next state function
    b2))
    -> ((TimeStamp, b1), (TimeStamp, initial state values tupled with their initial delay
        b2))
-> Signal a1 first input signal
-> Signal a2 second input signal
-> (Signal b1, Signal b2) output signals

stated is a state machine without an output decoder. It is an instantiation of the state MoC constructor (see stated22).

Constructors: stated[1-4] [1-4].
```

\lambda> let s = readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: Signal Int
\lambda> takeS 7 \$ stated11 (+) (6,1) s
{1 @0s, 2 @6s, 3 @8s, 5 @12s, 7 @14s, 8 @15s, 10 @18s}

```

```

state22

```
    :: (b1 -> b2 -> a1 -> a2 -> (b1, next state function
        b2))
    -> ((TimeStamp, b1), (TimeStamp,
        b2) )
    -> Signal a1
    -> Signal a2
    -> (Signal b1, Signal b2) output signals
        initial state values tupled with their initial delay
    first input signal
    second input signal
    state is a state machine without an output decoder, and the state non-transparent. It is an
    instantiation of the state MoC constructor (see state22).
    Constructors: state[1-4] [1-4].
\(\mid \lambda>\) let \(s=\) readSignal "\{1@0, 2@2, 3@6, 4@8, 5@9\}" :: Signal Int
\(\lambda>\) takeS 7 \$ state11 (+) \((6,1)\) s
\{ 2 @ \(\mathrm{s}, 3\) @ \(3 \mathrm{~s}, 5\) @ \(5 \mathrm{~s}, 7\) @ \(8 \mathrm{~s}, 8\) @9s, 10 @12s, 12 @14s\}
moore22
    \(\therefore \quad(\) st \(\rightarrow\) a1 -> a2 -> st) next state function
    -> (st -> (b1, b2)) output decoder
    -> (TimeStamp, st) initial state: tag and value
    -> Signal a1
    -> Signal a2
        (Signal b1, Signal b2)
moore processes model Moore state machines. It is an instantiation of the moore MoC constructor (see moore22).
Constructors: moore [1-4] [1-4]
```

\lambda> let s = readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: Signal Int
\lambda> takeS 7 \$ moore11 (+) (+1) (6,1) s
{ 2 @0s, 3 @6s, 4 @8s, 6 @12s, 8 @14s, 9 @15s, 11 @18s}

```
mealy22
    :: (st -> a1 -> a2 -> st) next state function
    -> (st -> a1 -> a2 -> (b1, b2)) outpt decoder
    -> (TimeStamp, st) initial state: tag and value
    -> Signal a1
    -> Signal a2
    -> (Signal b1, Signal b2)

mealy processes model Mealy state machines. It is an instantiation of the mealy MoC constructor (see mealy22).
Constructors: mealy [1-4] [1-4]
```

\lambda> let s = readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: Signal Int
\lambda> takeS 7 \$ mealy11 (+) (-) (6,1) s
{ 0 @0s, -1 @2s, -1 @6s, -1 @8s, -2 @9s, 0 @12s, 2 @14s}

```


\section*{Interface processes}
toSY2
```

:: Signal a first input DE signal
-> Signal b second input DE signal
-> (Signal TimeStamp, Signal a, signal carrying timestamps tupled with the two out-
Signal b)
put SY signals

```

Synchronizes a (set of) DE signal(s) an strips off their explicit tags, outputting the equivalent SY signal(s), tupled with an SY signal carrying the timestamps for the synchronization points.
Constructors : toSy [1-4].
```

\lambda> let s1 = DE.infinite 1
\lambda> let s2 = DE.readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: DE.Signal Int
\lambda> toSY2 s1 s2
({0s,2s,6s,8s,9s},{1,1,1,1,1},{1,2,3,4,5})

```

```

:: Signal (Time -> a) first input DE signal
-> Signal (Time -> b) second input DE signal
-> (Signal a, Signal b) two output CT signals

```
toCT2

Semantic preserving transformation between a (set of) DE signal(s) and the equivalent CT signals. The DE events must carry a function of Time which will be lifted by providing it with CT implicit time semantics.
Constructors: toct [1-4].
zipx :: Vector (Signal a) -> Signal (Vector a)
Synchronizes all the signals contained by a vector and zips them into one signal of vectors. It instantiates the zipx skeleton.
```

\lambda> let s1 = DE.readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: DE.Signal Int
\lambda> let s2 = DE.readSignal "{1@0, 2@2, 3@4, 4@8, 5@9}" :: DE.Signal Int
\lambda> let v1 = V.vector [s1,s1,s2,s2]
\lambda> v1
<{ 1 @0s, 2 @2s, 3 @6s, 4 @8s, 5 @9s},{ 1 @0s, 2 @2s, 3 @6s, 4 @8s, 5 @9s},{ 1 @0s, 2 @2s, 3
@4s, 4 @8s, 5 @9s},{ 1 @Os, 2 @2s, 3 @4s, 4 @8s, 5 @9s}>
\lambda> zipx v1
{<1,1,1,1> @0s, <2,2,2,2> @2s, <2,2,3,3>@4s, <3,3,3,3> @6s, <4,4,4,4> @8s, <5,5,5,5> @9s}

```

unzipx : : Integer -> Signal (Vector a) -> Vector (Signal a)
Unzips the vectors carried by a signal into a vector of signals. It instantiates the unzipx skeleton. To avoid infinite recurrence, the user needs to provide the length of the output vector.
```

\lambda> let v1 = V.vector [1,2,3,4]
\lambda> let s1 = DE.signal [(0,v1),(2,v1),(6,v1),(8,v1),(9,v1)]
\lambda> s1
{<1,2,3,4> @0s, <1,2,3,4> @2s, <1,2,3,4> @6s, <1,2,3,4> @8s, <1,2,3,4> @9s}
\lambda> unzipx 4 s1
<{ 1 @Os, 1 @2s, 1 @6s, 1 @8s, 1 @9s},{ 2 @Os, 2 @2s, 2 @6s, 2 @8s, 2 @9s},{ 3 @0s, 3 @2s, 3
@6s, 3 @8s, 3 @9s},{ 4 @0s, 4 @2s, 4 @6s, 4 @8s, 4 @9s}>

```

unzipx' :: Signal (Vector a) -> Vector (Signal a)
Same as unzipx, but "sniffs" the first event to determine the length of the output vector. Might have unsafe behavior!
```

\lambda> let v1 = V.vector [1,2,3,4]
\lambda> let s1 = DE.signal [(0,v1),(2,v1),(6,v1),(8,v1),(9,v1)]
\lambda> s1
{<1,2,3,4> @0s, <1,2,3,4> @2s, <1,2,3,4> @6s, <1,2,3,4> @8s, <1,2,3,4> @9s}
\lambda> unzipx' s1
<{ 1 @Os, 1 @2s, 1 @6s, 1 @8s, 1 @9s},{ 2 @0s, 2 @2s, 2 @6s, 2 @8s, 2 @9s},{ 3 @0s, 3 @2s, 3
@6s, 3 @8s, 3 @9s},{ 4 @0s, 4 @2s, 4 @6s, 4 @8s, 4 @9s}>

```

\section*{Hybrid processes}

\section*{embedSY22}
:: (Signal a1 -> Signal a2 -> (Signal b1, Signal b2)) SY process
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)
first input DE signal second input DE signal two output DE signals

Embeds a SY process inside a DE environment. Internally, it synchronizes the input signals, translates them to SY, feeds them to a SY process and translates the result back to DE using the same input tags. Seen from outside, this process behaves like a DE process with "instantaneous response", even for feedback loops.
Constructors: embedSY [1-4] [1-4].
For the following example, see the difference between its output and the one of stated22
```

\lambda> let s = readSignal "{1@0, 2@2, 3@6, 4@8, 5@9}" :: DE.Signal Int
\lambda> embedSY11 (SY.stated11 (+) 1) s
{ 1 @0s, 2 @2s, 4 @6s, }7\mathrm{ @8s, 11 @9s}

```


\subsection*{4.9 ForSyDe.Atom.MoC.CT}
```

module ForSyDe.Atom.MoC.CT (
TimeStamp, Time, CT(CT, tag, phase, func), Signal, unit2, infinite,
signal, checkSignal, delay, delay', comb22, reconfig22, constant2,
infinite2, generate2, stated22, state22, moore22, mealy22, toDE1,
sampDE2, zipx, unzipx, unzipx'
) where

```

The CT library implements the atoms holding the sematics for the continuous time computation model. It also provides a set of helpers for properly instantiating process network patterns as process constructors.

For working with time or timestamps please check the utilities provided by the ForSyDe.Atom.MoC.Time and ForSyDe.Atom.MoC.TimeStamp modules.

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.

\subsection*{4.9.1 Continuous time (ст) event}

According to Lee and Sangiovanni-Vincentelli, 1998, "[regarding metric time] at a minimum, \(T\) is an Abelian group, in addition to being totally ordered. A consequence is that \(t_{2}-t_{1}\) is itself a tag \(\forall t_{1}, t_{2} \in T\). In a slightly more elaborate model of computation, \(T\) has a metric. (...) A continuous-time system is a metric timed system \(Q\) where \(T\) is a continuum (a closed connected set)."

The continuous time (ст) MoC defines the closest behavior to what we could call "physical time", where signals cover the full span of a simulation as "functions of time" rather than "values". As such, we can state:

The CT MoC is abstracting the execution semantics and describes a system where computation is performed continuously over a (possibly infinite) span of time.

Below is an illustration of the behavior in time of the input and the output signals of a CT process:


Our CT MoC is implemented as an enhanced version of DE with respect to the CT MoC definition, in the sense that:
1. tags \(t\) are also represented with TimeStamps, thus we can say that changes in a CT signal happen at discrete times (see below).
2. values are represented as functions over a continuous span of time \(f(t)\) rather than just a value or a series of values. The time domain is represented with rational numbers which, as compared to floating point numbers, do not suffer from inherent quantisation, being able to model true continuity, i.e. between any two arbitrary points in time there is an infinite amount of intermediate moments.
3. The event constructor has also a phase component \(\phi\), which is taken into consideration only when evaluating the event function, i.e. \(f(t+\phi)\). This enables the modeling of "phase dispacements" of delay lines without altering the function itself (and thus increasing the complexity of the un-evaluated function graph). The phase needs to be reset during event synchronization.

These seemingly minor changes have deep implications in the expressiveness of a FoSyDe CT system and how we interpret it. Capturing the particularities of this MoC, we can formulate the following properties:

1. CT signals, due to their formation as streams of tagged events, represent discrete changes in a continuous function over time (e.g. analog signal). While the functions carried by events are infinite (have always happened and will always happen), being carried by events in a tag system suggests that changes occur at discrete times. A CT signal can be represented by the analog circuit above, where the inputs are continuous signals, but the switch is discrete. Like in the DE MoC, the absolute time 0 represent the time when the system started to be observed.
2. the previous property is also proven by the fact that the evaluation engine of ForSyDe-Atom is inherently discrete, i.e. evaluation is performed whenever a new event occurs, in a dataflow manner. Allowing infinitely small distances between tags would hinder the advancement of simulation time.
3. events carry functions and not values. In a lazy evaluation system like Haskell's, functions are kept symbolic until evaluation. This means that in a CT system computations are propagated as function graphs until a result is needed, e.g. a signal is plotted for arbitrary positions in time. This way intermediate quantization errors are eliminated, and the cost of higher plot resolution is the cost of evaluating the final results only.
4. needless to say, for each \(t \in T\), a signal is able to return (e.g. plot) the exact value \(v\) for that particular \(t\).
5. since itself the CT MoC is just an enhanced DE system, all atom evaluation properties are inherited from it: feedback loops need to advance time, atoms are forbidden to clean signals, and the conservative approach makes it ideal for parallel/distributed simulation.
6. since \(T\) is a total order, there is no need for an execution context (see section 4.5) and we can ignore the formatting of functions in ForSyDe.Atom.MoC, thus we can safely assume:
\[
\dot{a}=a
\]
type TimeStamp = DiffTime
Alias for the type representing discrete time. It is inherently quantizable, the quantum being a picosecond ( \(10^{-12}\) seconds), thus it can be considered order-isomorphic with a set of integers, i.e. between any two timestamps there is a finite number of timestamps. Moreover, a timestamp can be easily translated into a rational number representing fractions of a second, so the conversion between timestamps (discrete time) and rationals (analog/continuous time) is straightforward.

This type is used in the explicit tags of the DE MoC (and subsequently the discrete event evaluation engine for simulating the CT MoC ).
```

type Time = Rational

```

Type alias for the type to represent metric (continuous) time. Underneath we use Rational that is able to represent any \(t\) between \(t_{1}<t_{2} \in T\).
data CT a
\(=\quad \mathrm{CT}\)
> tag : : TimeStamp start time of event
> phase :: Time phase. Models function delays
\(>\) func :: Time \(->\) a function of time
The CT type, identifying a continuous time event and implementing an instance of the MoC class.
```

instance Functor CT

```

Allows for mapping of functions on a CT event.
```

instance Applicative CT

```

Allows for lifting functions on a pair of CT events.
```

instance MoC CT

```

Implenents the execution and synchronization semantics for the CT MoC through its atoms.
```

instance Show a => Show (CT a)

```

A non-ideal instance meant for debug purpose only, For each event it evaluates the function at the tag time only!
```

instance Plottable a => Plot (Signal a)

```

CT signals.
```

instance type Ret CT b = b

```
instance type Fun CT a b \(=\mathrm{a}->\mathrm{b}\)

\subsection*{4.9.2 Aliases \& utilities}

A set of type synonyms and utilities are provided for convenience. The API type signatures will feature these aliases to hide the cumbersome construction of atoms and patters as seen in ForSyDe.Atom.MoC.
```

type Signal a = Stream (CT a)
Type synonym for a CT signal, i.e. "a signal of CT events"
unit2 :: ((TimeStamp, Time -> a1), (TimeStamp, Time -> a2))
-> (Signal a1, Signal a2)

```

Wraps a (tuple of) pair(s) (time, function) into the equivalent unit signal(s), i.e. signal(s) with one event with the period time carrying function.
Helpers: unit and unit[2-4].
```

infinite :: (Time -> a) -> Signal a

```

Creates an infinite signal.
```

signal :: [(TimeStamp, Time -> a)] -> Signal a

```

Transforms a list of tuples such as the ones taken by event into a CT signal
```

checkSignal :: Stream (CT a) -> Stream (CT a)

```

Checks if a signal is well-formed or not, according to the CT MoC interpretation in ForSyDe-Atom.

\subsection*{4.9.3 Ст process constuctors}

The CT process constructors are basically specific instantiations of patterns defined in ForSyDe.Atom. MoC. Some might also be wrapping functions in an extended behavioural model.

In the examples below we have imported and instantiated the functions such as e' pi ', sin' and cos' from the collection of utilities in ForSyDe.Atom.MoC.Time and ForSyDe.Atom.MoC.TimeStamp. Also, for the sake of documentation the interactive examples are only dumping the CT signals in data files using the dumpDat utility defined in ForSyDe.Atom.Utility.Plot, according to the custom cfg structure. These files can be further plotted by any tool of choice, or using the plotting utilities provided in the ForSyDe.Atom.Utility. Plot module.
```

import ForSyDe.Atom.Utility.Plot
import ForSyDe.Atom.MoC.Time as Time
import ForSyDe.Atom.MoC.TimeStamp as TimeStamp
let pi' = TimeStamp.pi
let exp' = Time.exp
let sin' = Time.sin
let cos' = Time.cos
let cfg = defaultCfg {xmax=10, rate=0.1}

```

\section*{Simple}

These are mainly direct instantiations of patterns defined in ForSyDe.Atom.MoC, using DE-specific utilities.
\begin{tabular}{|c|}
\hline \multirow[b]{3}{*}{\begin{tabular}{l}
-> Signal a input signal \\
-> Signal a output signal \\
The delay process "delays" a signal with one e CT MoC, this process can be interpreted as an
```

\lambda> let s = infinite (sin')
\lambda> let s' = delay 2 (\_ -> 0) s
\lambda> dumpDat \$ prepare cfg {labels=["delay"]} \$ s,
Dumped delay in ./fig
["./fig/delay.dat"]

```
\end{tabular}} \\
\hline \\
\hline \\
\hline
\end{tabular}

delay'
:: Signal a signal "borrowing" the initial event
-> Signal a input signal
-> Signal a output signal

Similar to the previous, but this is the raw instantiation of the delay pattern. It "borrows" the first event from one signal and appends it at the head of another signal.
```

\lambda> let s = infinite (sin')
\lambda> let s' = signal [(0, \_ -> 0), (2, \_ -> 1)]
\lambda> dumpDat \$ prepare cfg {labels=["delayp"]} \$ delay' s's
Dumped delayp in ./fig
["./fig/delayp.dat"]

```

comb22
:: (a1 -> a2 -> (b1, b2)) function on values
-> Signal a1 first input signal
-> Signal a2 second input signal
-> (Signal b1, Signal b2) two output signals
comb processes map combinatorial functions on signals and take care of synchronization between input signals. It instantiates the comb pattern (see comb22).
Constructors: comb[1-4] [1-4].
```

\lambda> let s1 = infinite (sin')
\lambda> let s2 = signal [(0,\_->0),(pi',\_->1),(2*pi',\_->0),(3*pi',\_->1)]
\lambda> let o1 = comb11 (+1) s2
\lambda> let (o2_1, o2_2) = comb22 (\a b-> (a+b,a*b)) s1 s2
\lambda> dumpDat \$ prepare cfg {labels=["comb11"]} o1
Dumped comb11 in ./fig
["./fig/comb11.dat"]
\lambda> dumpDat \$ prepareL cfg {labels=["comb22-1","comb22-2"]} [o2_1, o2_2]
Dumped comb22-1, comb22-2 in ./fig
["./fig/comb22-1.dat","./fig/comb22-2.dat"]

```


\section*{reconfig22}
:: Signal (a1 -> a2 -> (b1, b2)) signal carrying functions
-> Signal a1 first input signal carrying arguments
-> Signal a2
-> (Signal b1, Signal b2) second input signal carrying arguments two output signals
reconfig creates a CT adaptive process where the first signal carries functions and the other carry the arguments. It instantiates the reconfig atom pattern (see reconfig22).
Constructors: reconfig[1-4] [1-4].
```

\lambda> let s1 = infinite (sin')
\lambda> let sf = signal [(0,\_->(*0)),(pi',\_-> (+1)),(2*pi',\_->(*0)),(3*pi',\_-> (+1))]
\lambda> dumpDat \$ prepare cfg {labels=["reconfig"]} \$ reconfig11 sf s1
Dumped reconfig in ./fig
["./fig/reconfig.dat"]

```
constant2
\(::(b 1, b 2) \quad\) values to be repeated
-> (Signal b1, Signal b2) generated signals

A generator for a constant signal. As compared with the constant2, it just constructs an infinite signal with constant value (i.e. a signal with one event starting from time 0 ).
Constructors: constant[1-4].
```

|> dumpDat \$ prepare cfg {labels=["constant"]} \$ constant1 2
Dumped constant in ./fig
["./fig/constant.dat"]

```

```

infinite2

```
:: (Time -> b1, Time -> b2) values to be repeated
-> (Signal b1, Signal b2) generated signals

A generator for an infinite signal. Similar to constant2.
Constructors: infinite[1-4].
\(\mid \lambda>\) let (inf1, inf2) \(=\) infinite2 (sin', cos')
\(\lambda>\) dumpDat \(\$\) prepareL cfg \{labels=["infinite2-1", "infinite2-2"]\} [inf1, inf2]
Dumped infinite2-1, infinite2-2 in ./fig
["./fig/infinite2-1.dat","./fig/infinite2-2.dat"]


\section*{generate2}
```

:: (b1 -> b2 -> (b1, b2))
-> ((TimeStamp, Time -> b1), (TimeStamp,
Time -> b2))
-> (Signal b1, Signal b2)

```
function to generate next value kernel values tupled with their generation rate. generated signals

A signal generator based on a function and a kernel value. It is actually an instantiation of the statedoX constructor (check stated22).
Constructors: generate [1-4].
```

\lambda> let { osc 0=1 ; osc 1 = 0 }
\lambda> dumpDat \$ prepare cfg {labels=["generate"]} \$ generate1 osc (pi', \_->0)
Dumped generate in ./fig
["./fig/generate.dat"]

```


Another example simulating an RC oscillator:
\[
\lambda>\text { let vs }=2 \quad \text {-- Vs : supply voltage }
\]
\(\begin{array}{ll}\lambda>\text { let } r=100 & --R: \text { resistance } \\ \lambda>\text { let } c=0.0005 & --C: \text { capacitance }\end{array}\)
\(\lambda>\) let c \(=0.0005\)
\(\lambda>\) let vc \(\mathrm{t}=\mathrm{vs} *(1-\exp ,(-\mathrm{t} /(\mathrm{r} * \mathrm{c})))\)-- Vc(t): voltage charging through capacitor
\(\lambda>\) let \(\mathrm{ns} \mathrm{v}=\mathrm{vs}+(-1 * \mathrm{v}) \quad--\) next state : charging / discharging
\(\lambda>\) let rc0sc \(=\) generate1 ns (milisec 500, vc)
\(\lambda>\) dumpDat \(\$\) prepare cfg \{labels=["rcosc"]\} \$ rcOsc
Dumped rcosc in ./fig
["./fig/rcosc.dat"]
```

stated22

```
```

:: (b1 -> b2 -> a1 -> a2 -> (b1, b2))
-> ((TimeStamp, Time -> b1), (TimeStamp,
Time -> b2))
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)

```
next state function
initial state values tupled with their initial delay
first input signal second input signal
output signals

stated is a state machine without an output decoder. It is an instantiation of the state MoC constructor (see stated22).

Constructors: stated[1-4] [1-4].
```

\lambda> let { osc 0 a = a; osc _ a = 0 }
\lambda> let s1 = signal [(0,\_->1), (6,\_->0)]
\lambda> dumpDat \$ prepare cfg {labels=["stated"]} \$ stated11 osc (1,\_->0) s1
Dumped stated in ./fig
["./fig/stated.dat"]

```


\section*{state22}
```

:: (b1 -> b2 -> a1 -> a2 -> (b1, b2))
-> ((TimeStamp, Time -> b1), (TimeStamp,
Time -> b2))
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)

```
next state function
initial state values tupled with their initial delay
first input signal
second input signal
output signals
state is a state machine without an output decoder, and the state non-transparent. It is an instantiation of the state MoC constructor (see state22).

Constructors: state [1-4] [1-4].
```

\lambda> let { osc 0 a = a; osc _ a = 0 }
\lambda> let s1 = signal [(0,\_->1), (6,\_->0)]
\lambda> dumpDat \$ prepare cfg {labels=["state"]} \$ state11 osc (1,\_->0) s1
Dumped state in ./fig
["./fig/state.dat"]

```
moore22
```

:: (st -> a1 -> a2 -> st) next state function
-> (st -> (b1, b2)) output decoder
-> (TimeStamp, Time -> st) initial state: tag and value
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)

```
moore processes model Moore state machines. It is an instantiation of the moore MoC constructor (see moore22).

Constructors: moore[1-4] [1-4].
```

\lambda> let { osc 0 a = a; osc _ a = 0 }
\lambda> let s1 = signal [(0,\_->1), (6,\_->0)]
\lambda> dumpDat \$ prepare cfg {labels=["moore"]} \$ moore11 osc (*3) (1,\_->0) s1
Dumped moore in ./fig
["./fig/moore.dat"]

```

mealy22
```

:: (st -> a1 -> a2 -> st) next state function
-> (st -> a1 -> a2 -> (b1, b2)) outpt decoder
-> (TimeStamp, Time -> st) initial state: tag and value
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)

```
    mealy processes model Mealy state machines. It is an instantiation of the mealy MoC con-
    structor (see mealy22).
    Constructors: mealy [1-4] [1-4].
```

\lambda> let { osc (-1) _ = 1; osc 1 _ = (-1) }
\lambda> let s1 = infinite sin'
\lambda> dumpDat \$ prepare cfg {labels=["mealy"]} \$ mealy11 osc (*) (pi',\_->1) s1
Dumped mealy in ./fig
["./fig/mealy.dat"]

```

\section*{Interfaces}
toDE1 : : Signal a -> Signal (Time -> a)
Translates a (set of) CT signal(s) into DE semantics without loss of information. In DE, the abstract function of time inferred by the CT event loses its abstraction and it is "dropped" to explicit form, under a lower layer. In other words the implicit time semantics are lost, the carried value simply becoming an ordinary function.
Constructors: toDE[1-4].


Synchronizes a (set of) CT signal(s) with a DE carrier which holds the timestamps at which the CT signal must be sampled, and outputs the respective (set of) DE signal(s).
Constructors: sampDE [1-4].
```

\lambda> let s = CT.infinite (fromRational . sin')
\lambda> let c = DE.generate1 id (pi'/2, 1)
\lambda> takeS 6 \$ sampDE1 c s
{ 0.0 @Os, 1.0 @1.570796326794s, 1.793238520564752e-12 @3.141592653588s, -1.0 @4.712388980382
s, 0.0 @6.283185307176s, 1.0 @7.85398163397s}

```
zipx :: Vector (Signal a) -> Signal (Vector a)

Synchronizes all the signals contained by a vector and zips them into one signal of vectors.
It instantiates the zipx skeleton.

```

\lambda> let s1 = CT.signal [(0,const 1), (2,const 2), (6,const 3)]
\lambda> let s2 = CT.signal [(0,const 1), (2,const 2), (4,const 3)]
\lambda> let v1 = V.vector [s1,s1,s2,s2]
\lambda> zipx v1
{ <1,1,1,1> @0s, <2,2,2,2> @2s, <2,2,3,3> @4s, <3,3,3,3> @6s}

```

See zipx from the ForSyDe.Atom.MoC.DE library for a comprehensive visual example.
```

unzipx :: Integer -> Signal (Vector a) -> Vector (Signal a)

```

Unzips the vectors carried by a signal into a vector of signals. It instantiates the unzipx skeleton. To avoid infinite recurrence, the user needs to provict the length of the output vector.
```

\lambda> let v1 = V.vector [1,2,3,4]
\lambda> let s1 = CT.signal [(0,const v1),(2,const v1),(6,const v1)]
\lambda> unzipx 4 s1
<{4 @0s, 4 @2s, 4 @6s},{ 3 @0s, 3 @2s, 3 @6s},{ 2 @0s, 2 @2s, 2 @6s},{ 1 @0s, 1 @2s, 1 @6s}>

```

See unzipx from the ForSyDe.Atom.MoC.DE library for a comprehensive visual example.
```

unzipx' :: Signal (Vector a) -> Vector (Signal a)

```

Same as unzipx, but "sniffs" the first event to determine the length of the output vector. Might have unsafe behavior!
```

\lambda> let v1 = V.vector [1,2,3,4]
\lambda> let s1 = CT.signal [(0,const v1),(2,const v1),(6,const v1)]
\lambda> unzipx' s1
<{ 4 @0s, 4 @2s, 4 @6s},{ 3 @0s, 3 @2s, 3 @6s},{ 2 @0s, 2 @2s, 2 @6s},{ 1 @0s, 1 @2s, 1 @6s}>

```

\subsection*{4.10 ForSyDe.Atom.MoC.SDF}
module ForSyDe.Atom.MoC.SDF (
SDF (SDF, val), Signal, Prod, Cons, signal, readSignal, delay,
delay', comb22, reconfig22, constant2, generate2, stated22, state22, moore22, mealy22, toSY2, zipx, unzipx
) where

The SDF library implements the atoms holding the sematics for the synchronous data flow computation model. It also provides a set of helpers for properly instantiating process network patterns as process constructors.

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.

\subsection*{4.10.1 Synchronous data flow (SDF) event}

The synchronous data flow (SDF) MoC is the first untimed MoC implemented by the forsyde-atom framework. On untimed MoCs, Lee and Sangiovanni-Vincentelli, 1998 states that: "when tags are partially ordered rather than totally ordered, we say that the system is untimed. Untimed systems cannot have the same notion of causality as timed systems [see Sy]. (...) Processes defined in terms of constraints on the tags in the signals (...) have a consistent cut rather than simultaneity." Regarding SDF, it states that "is a special case of Kahn process networks Kahn and MacQueen, 1976. A dataflow process is a Kahn process that is also sequential, where the events on the selfloop signal denote the firings of the dataflow actor. The firing rules of a dataflow actor are partial ordering constraints between these events and events on the inputs. (...) Produced/consumed events are defined in terms of relations with the events in the firing signal. It results that for the same firing \(i, \mathrm{e}_{\mathrm{i}}<\mathrm{e}_{\mathrm{o}}\), as an intuitive sort of causality constraint."

Based on the above insights, we can formulate a simplified definition of the forsyde-atom interpretation of SDF:

The SDF MoC is abstracting the execution semantics of a system where computation is performed according to firing rules where the production and the consumption rates are fixed.

Below is a possible behavior in time of the input and the output signals of a SDF process. Events sharing the same partial ordering in relation to one firing are overlined:


Implementing the SDF tag system implied a series of engineering decisions which lead to the following particularities:
1. signals represent FIFO channels, and tags are implicit from their position in the Stream structure. Internally, SDF signals have exactly the same structure as SY signals, whereas the partial ordering is imposed by the processes alone.
2. the SDF event constructor wraps only a value.
3. being an untimed \(M o C\), the order between events is partial to the firings of processes. An SDF atom will fire only when there are enough events to trigger its inputs. Once a firing occurs, it will take care of partitioning the input or output signals.
4. SDF atoms \(d o\) require a context: the consumption \(c\) and production \(p\) rates. Also, the functions passed as arguments reflect the fact that multiple events are handled during a firing.
5. the previous statement can be synthesized into the following execution context (see section 4.5), which also justifies the SDF implementation of Fun and for Ret:
\[
\dot{a}=a^{n}, \text { where } n \in \mathbb{N}
\]
```

newtype SDF a

```
\(=\quad \mathrm{SDF}\)
> val : : a
The CT type, identifying a discrete time event and implementing an instance of the MoC class. A discrete event explicitates its tag which is represented as an integer.
instance Functor SDF
Allows for mapping of functions on a SDF event.

\section*{instance Applicative SDF}

Allows for lifting functions on a pair of SDF events.
```

instance Foldable SDF
instance Traversable SDF
instance MoC SDF
Implenents the SDF semantics for the MoC atoms

```
```

instance Read a => Read (SDF a)

```

Reads the value wrapped
```

instance Show a => Show (SDF a)

```

Shows the value wrapped
```

instance Plottable a => Plot (Signal a)

```

SDF signals.
```

instance type Ret SDF a = (Prod, [a])
instance type Fun SDF a b = (Cons, [a] -> b)

```

\subsection*{4.10.2 Aliases \& utilities}

A set of type synonyms and utilities are provided for convenience. The API type signatures will feature these aliases to hide the cumbersome construction of atoms and patters as seen in ForSyDe.Atom.MoC.
```

type Signal a = Stream (SDF a)

```
    Type synonym for a SY signal, i.e. "a signal of SY events"
type Prod = Int

Type synonym for consumption rate
```

type Cons = Int

```

Type synonym for production rate
signal :: [a] -> Signal a
Transforms a list of values into a SDF signal with only one partition, i.e. all events share the same (initial) tag.
readSignal : : Read a \(=>\) String \(->\) Signal a
Reads a signal from a string. Like with the read function from Prelude, you must specify the tipe of the signal.
```

\lambda> readSignal "{1,2,3,4,5}" :: Signal Int
{1,2,3,4,5}

```

These SY process constructors are basically specific instantiations of the patterns of atoms defined in ForSyDe.Atom.MoC. Some are also wrapping functions in an extended behavioural model.

\section*{Simple}
delay
:: [a] list of initial values
-> Signal a input signal
-> Signal a output signal
The delay process "delays" a signal with initial events built from a list. It is an instantiation of the delay constructor.
```

\lambda> let s = signal [1,2,3,4,5]
\lambda> delay [0,0,0] s
{0,0,0,1,2,3,4,5}

```

delay,
:: Signal a signal containing the initial tokens
-> Signal a input signal
-> Signal a output signal
Similar to the previous, but this is the raw instantiation of the delay pattern. It appends the contents of one signal at the head of another signal.
```

\lambda> let s1 = signal [0,0,0]
\lambda> let s2 = signal [1,2,3,4,5]
\lambda> delay' s1 s2
{0,0,0,1,2,3,4,5}

```
comb22
:: ((Cons, Cons), (Prod, Prod), [a1] -> function on lists of values, tupled with con[a2] -> ([b1], [b2])) sumption / production rates
-> Signal a1 first input signal Signal a2 second input signal
-> (Signal b1, Signal b2) two output signals
comb processes map combinatorial functions on signals and take care of synchronization between input signals. It instantiates the comb atom pattern (see comb22).
Constructors: comb [1-4] [1-4].
```

\lambda> let s1 = signal [1..]
\lambda> let s2 = signal [1,1,1,1,1,1,1]
\lambda> let f [a,b,c] [d,e] = [a+d, c+e]
\lambda> comb21 ((3,2),2,f) s1 s2
{2,4,5,7,8,10}

```

Incorrect usage (not covered by doctest):
\[
\lambda>\operatorname{comb21}((3,2), 3, f) \mathrm{s} 1 \mathrm{~s} 2
\]
*** Exception: [MoC.SDF] Wrong production
```

reconfig22
:: ((Cons, Cons), (Prod,
Prod))
([b1], [b2]))
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)

```
    -> Signal ([a1] -> [a2] -> function on lists of values, tupled with consumption /
    production rates
    first input signal
    second input signal
    two output signals
reconfig creates an SDF adaptive process where the first signal carries functions and the other carry the arguments. It instantiates the reconfig atom pattern (see reconfig22). According to our SDF definition, the production and consumption rates need to be fixed, so they are passed as parameters to the constructor, whereas the first signal carries adaptive functions only. For the adaptive signal it only makes sense that the consumption rate is always 1 .
Constructors: reconfig[1-4] [1-4].
```

\lambda> let f1 a = [sum a]
\lambda> let f2 a = [maximum a]
\lambda> let sf = signal [f1,f2,f1,f2,f1,f2,f1]
\lambda> let s1 = signal [1..]
\lambda> reconfig11 (4,1) sf s1
{10,8,42,16,74,24,106}

```
constant2

:: ([b1], [b2])
-> (Signal b1, Signal b2) generated signals
A signal generator which repeats the initial tokens indefinitely. It is actually an instantiation of the stated0X constructor (check stated22).
Constructors: constant [1-4].
```

\lambda> let (s1, s2) = constant2 ([1,2,3], [2,1])

```
\(\lambda>\) takeS 7 s 1
\(\{1,2,3,1,2,3,1\}\)
\(\lambda>\) takeS 5 s2
\(\{2,1,2,1,2\}\)
generate2
: : ((Cons, Cons), (Prod, Prod), [b1]
\(->[b 2]->([b 1],[b 2]))\)
-> ([b1], [b2])
-> (Signal b1, Signal b2)
function to generate next value, tupled with consumption / production rates values of initial tokens generated signals

A signal generator based on a function and a kernel value. It is actually an instantiation of the statedoX constructor (check stated22).
Constructors: generate[1-4].
```

\lambda> let f a b = ([sum a, sum a],[sum b, sum b, sum b])
\lambda> let (s1,s2) = generate2 ((2,3),(2,3),f) ([1,1],[2,2,2])
\lambda> takeS 7 s1
{1,1,2,2,4,4,8}
\lambda> takeS 8 s2
{2,2,2,6,6,6,18,18}

```

```

stated22
:: ((Cons, Cons, Cons, Cons), (Prod, Prod), [b1]
-> [b2] -> [a1] -> [a2] -> ([b1], [b2]))
-> ([b1], [b2])
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)

```
next state function, tupled with consumption / production rates initial state partitions of values first input signal second input signal output signals
\(\qquad\)
\(\qquad\)

stated is a state machine without an output decoder. It is an instantiation of the state MoC constructor (see stated22).
Constructors: stated[1-4] [1-4].
:: ((Cons, Cons, Cons, Cons), (Prod, Prod), [b1]
        -> [b2] -> [a1] -> [a2] -> ([b1], [b2]))
    -> ([b1], [b2])
    -> Signal a1
    -> Signal a2
    -> (Signal b1, Signal b2)
state is a state machine without an output decoder. It is an instantiation of the stated MoC constructor (see state22).
Constructors: state [1-4] [1-4].
\(\lambda>\) let \(f[a][b, c]=[a+b+c]\)
\(\lambda>\) let \(s=\) signal \([1,2,3,4,5,6,7]\)
\(\lambda>\) state11 \(((1,2), 1, f)\) [1] \(s\)
\(\{4,11,22\}\)

moore22
:: ((Cons, Cons, Cons), Prod, [st] -> [a1] -> [a2] -> [st])
-> (Cons, (Prod, Prod), [st] -> ([b1], [b2]))
-> [st]
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)
moore processes model Moore state machines. It is an instantiation of the moore MoC constructor (see moore22).
Constructors: moore [1-4] [1-4].
next state function, tupled with consumption / production rates initial partitions of values
first input signal
second input signal output signals
\(\qquad\) It is an in
```

\lambda> let f [a] [b,c] = [a+b+c]

```
\lambda> let f [a] [b,c] = [a+b+c]
\lambda> let s = signal [1,2,3,4,5,6,7]
\lambda> let s = signal [1,2,3,4,5,6,7]
\lambda> stated11 ((1,2),1,f) [1] s
\lambda> stated11 ((1,2),1,f) [1] s
{1,4,11,22}
```

{1,4,11,22}

```

```

\lambda> let ns [a] [b,c] = [a+b+c]
\lambda> let od [a] = [a+1,a*2]
\lambda> let s = signal [1,2,3,4,5,6,7]
\lambda> moore11 ((1,2),1,ns) (1,2,od) [1] s
{2,2,5,8,12,22,23,44}
$\xrightarrow{1,2,3,4,5,6,7} \xrightarrow{\lambda(a)(b, c) \rightarrow(a+b+c)} \overline{\lambda(a) \rightarrow(a+1, a \times 2)} \underline{(1)} \quad \underline{2,2,5,8}, \underline{12,22}, \underline{23,44}$

```
mealy22
:: ((Cons, Cons, Cons), Prod, [st] -> [a1] -> next state function, tupled with con-
        [a2] -> [st]) sumption / production rates
-> ((Cons, Cons, Cons), (Prod, Prod), [st] -> outpt decoder, tupled with consump-
        [a1] -> [a2] -> ([b1], [b2]))
-> [st]
-> Signal a1
-> Signal a2
-> (Signal b1, Signal b2)
mealy processes model Mealy state machines. It is an instantiation of the mealy MoC constructor (see mealy22).

Constructors: mealy [1-4] [1-4].
```

\lambda> let ns [a] [b,c] = [a+b+c]
\lambda> let od [a] [b] = [a+b,a*b]
\lambda> let s = signal [1,2,3,4,5,6,7]
\lambda> mealy11 ((1,2),1,ns) ((1,1),2,od) [1] s
{2,1,6,8,14,33,26,88}

```

\section*{Interfaces}
toSY2 : : Signal a \(\rightarrow\) Signal b -> (Signal a, Signal b)
Transforms a (set of) SDF signal(s) into the equivalent SY signal(s). The only change is the event consructor. The partial order of DE is interpreted as SY's total order, based on the positioning of events in the signals (e.g. FIFO buffers) at that moment.

Constructors: toSY[1-4].
```

\lambda> let s = SDF.signal [1, 2, 3,4,5]
\lambda> toSY s
{1,2,3,4,5}

```
zipx
\[
\begin{array}{lll}
:: & \text { Vector Cons } & \text { consumption rates } \\
-> & \text { Vector (Signal a) } & \text { vector of signals } \\
-> & \text { Signal (Vector a) } & \text { signal of vectors }
\end{array}
\]

Consumes tokens from a vector of signals and merges them into a signal of vectors, with a production rate of 1 . It instantiates the zipx skeleton.
```

\lambda> let s1 = SDF.signal [1,2,3,4,5]
\lambda> let s2 = SDF.signal [11,12,13,14,15]
\lambda> let v1 = V.vector [s1,s1,s2,s2]
\lambda> let r = V.vector [2,1,2,1]
\lambda> v1
<{1,2,3,4,5},{1,2,3,4,5},{11,12,13,14,15},{11,12,13,14,15}>
\lambda> zipx r v1
{<1, 2,1, 11, 12,11>,<3,4,2,13,14, 12>}

```

unzipx
\begin{tabular}{lll}
\(::\) & Vector Prod & production rates (in reverse order) \\
-> & Signal (Vector a) & signal of vectors \\
-> & Vector (Signal a) & vector of signals
\end{tabular}

Consumes the vectors carried by a signal with a rate of 1 , and unzips them into a vector of signals based on the user provided rates. It instantiates the unzipx skeleton.

OBS: due to the recur pattern contained by unzipx, the vector of production rates needs to be provided in reverse order (see ForSyDe.Atom.Skeleton.Vector).
```

\lambda> let s1 = SDF.signal [1,2,3,4,5]
\lambda> let s2 = SDF.signal [11,12,13,14,15]
\lambda> let v1 = V.vector [s1,s1,s2,s2]
\lambda> let r = V.vector [2,1,2,1]
\lambda> let sz = zipx r v1
\lambda> v1
<{1,2,3,4,5},{1,2,3,4,5},{11,12,13,14,15},{11,12,13,14,15}>
\lambda> sz
{<1,2,1,11,12,11>,<3,4,2,13,14,12>}
\lambda> unzipx (V.reverse r) sz
<{1,2,3,4},{1,2},{11,12,13,14},{11,12}>

```

\subsection*{4.11 ForSyDe.Atom.MoC.Time}
```

module ForSyDe.Atom.MoC.Time (
Time, time, const, e, (*^*), pi, sin, cos, tan, atan, asin,
acos, sqrt, exp, cosh, sinh, tanh, atanh, asinh, acosh, log
) where

```

Collection of utility functions for working with Time. While the Ст MoC describes time as being a non-disjoint continuum (represented in ForSyDe-Atom with Rational numbers), most of the
functions here are non-ideal approximations or conversions from floating point equivalents. The trigonometric functions are imported from the numbers package, with a fixed eps parameter.

These utilities are meant to get started with using the CT MoC, and should be used with caution if result fidelity is a requirement. In this case the user should find a native Rational implementation for a particular function.
```

type Time = Rational

```
    Type alias for the type to represent metric (continuous) time. Underneath we use Rational
    that is able to represent any \(t\) between \(t_{1}<t_{2} \in T\).
time :: TimeStamp -> Time
    Converts TimeStamp into Time representation.
const : : a -> Time -> a
    Returns a constant function.
e : : Time
    Euler's number in Time format. Converted from the Prelude equivalent, which is Floating.
(*~*) :: Time -> Time -> Time
    "Power of" function taking Times as arguments. Converts back and forth to Floating, as it
    uses the \(* *\) operator, so it is prone to conversion errors.
pi :: Time
    Time representation of the number \(\pi\). Rational representation with a precision of 0.000001 .
sin : : Time -> Time
    Sine of Time. Rational representation with a precision of 0.000001 .
cos : : Time -> Time
    Cosine of Time. Rational representation with a precision of 0.000001 .
tan :: Time -> Time
    Tangent of Time. Rational representation with a precision of 0.000001 .
atan : : Time -> Time

Arctangent of Time. Rational representation with a precision of 0.000001 .
asin :: Time \(->\) Time

Arcsine of Time. Rational representation with a precision of 0.000001 .
acos :: Time -> Time
Arccosine of Time. Rational representation with a precision of 0.000001 .
sqrt : : Time -> Time
Square root of Time. Rational representation with a precision of 0.000001 .
exp :: Time -> Time
Exponent of Time. Rational representation with a precision of 0.000001 .
cosh : : Time -> Time
Hyperbolic cosine of Time. Rational representation with a precision of 0.000001 .
sinh :: Time -> Time
Hyperbolic sine of Time. Rational representation with a precision of 0.000001 .
tanh :: Time -> Time
Hyperbolic tangent of Time. Rational representation with a precision of 0.000001 .
```

atanh :: Time -> Time

```

Hyperbolic arctangent of Time. Rational representation with a precision of 0.000001 .
```

asinh :: Time -> Time

```

Hyperbolic arcsine of Time. Rational representation with a precision of 0.000001 .
```

acosh :: Time -> Time

```

Hyperbolic arccosine of Time. Rational representation with a precision of 0.000001 .

\section*{log :: Time -> Time}

Natural logarithm of Time. Rational representation with a precision of 0.000001 .

\subsection*{4.12 ForSyDe.Atom.MoC.TimeStamp}
```

module ForSyDe.Atom.MoC.TimeStamp (
TimeStamp, picosec, nanosec, microsec, milisec, sec, minutes, hours,
toTime, pi,
) where

```

This module implements a timestamp data type, based on Data.Time.Clock.
```

type TimeStamp = DiffTime
Alias for the type representing discrete time. It is inherently quantizable, the quantum being a picosecond ( $10^{-12}$ seconds), thus it can be considered order-isomorphic with a set of integers, i.e. between any two timestamps there is a finite number of timestamps. Moreover, a timestamp can be easily translated into a rational number representing fractions of a second, so the conversion between timestamps (discrete time) and rationals (analog/continuous time) is straightforward.
This type is used in the explicit tags of the DE MoC (and subsequently the discrete event evaluation engine for simulating the CT MoC).

```
```

picosec :: Integer -> TimeStamp

```
picosec :: Integer -> TimeStamp
Specifies a timestamp in terms of picoseconds.
```

```
nanosec :: Integer -> TimeStamp
```

nanosec :: Integer -> TimeStamp
Specifies a timestamp in terms of nanoseconds.

```
```

microsec :: Integer -> TimeStamp

```
microsec :: Integer -> TimeStamp
Specifies a timestamp in terms of microseconds.
```

```
milisec :: Integer -> TimeStamp
```

milisec :: Integer -> TimeStamp
Specifies a timestamp in terms of miliseconds.

```
```

sec :: Integer -> TimeStamp

```
sec :: Integer -> TimeStamp
Specifies a timestamp in terms of seconds.
```

```
minutes :: Integer -> TimeStamp
```

minutes :: Integer -> TimeStamp
Specifies a timestamp in terms of minutes.

```
```

hours :: Integer -> TimeStamp

```
hours :: Integer -> TimeStamp
Specifies a timestamp in terms of hours.
```

```
toTime :: TimeStamp -> Rational
```

toTime :: TimeStamp -> Rational
Converts a timestamp to a rational number, used for describing continuous time.
pi :: TimeStamp
TimeStamp representation of the number $\pi$. Converted from the Prelude equivalent, which is Floating.

```

\subsection*{4.13 ForSyDe.Atom.Skeleton}
```

module ForSyDe.Atom.Skeleton (
Skeleton((=.=), (=*=), (=\=), (=<<=), first, last), farm22, reduce,
reducei, pipe, pipe2
) where

```

This module exports a type class with the interfaces for the Skeleton layer atoms. It does NOT export any implementation of atoms not any constructor as composition of atoms.

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.

\subsection*{4.13.1 Atoms}
class Functor c => Skeleton c where
Class containing all the Skeleton layer atoms.
This class is instantiated by a set of categorical types, i.e. types which describe an inherent potential for being evaluated in parallel. Skeletons are patterns from this layer. When skeletons take as arguments entities from the MoC layer (i.e. processes), the results themselves are parallel process networks which describe systems with an inherent potential to be implemented on parallel platforms. All skeletons can be described as composition of the three atoms below (=«= being just a specific instantiation of \(=\backslash=\) ). This possible due to an existing theorem in the categorical type theory, also called the Bird-Merteens formalism R. Bird and Moor, 1997:
factorization A function on a categorical type is an algorithmic skeleton (i.e. catamorphism) iff it can be represented in a factorized form, i.e. as a map composed with a reduce.

Consequently, most of the skeletons for the implemented categorical types are described in their factorized form, taking as arguments either:
- type constructors or functions derived from type constructors
- processes, i.e. MoC layer entities

Most of the ground-work on algorithmic skeletons on which this module is founded has been laid in R. Bird and Moor, 1997, Skillicorn, 1994 and it founds many of the frameworks collected in Fischer, Gorlatch, and Bischof, 2003.

\section*{Methods}
\[
(=.=)::(\mathrm{a}->\mathrm{b})->\mathrm{c} a->\mathrm{c} \text { b }
\]

Atom which maps a function on each element of a structure (i.e. categorical type), defined as:
\[
\stackrel{\diamond}{ }(\alpha \rightarrow \beta) \rightarrow \mathcal{S}(\alpha) \rightarrow \mathcal{S}(\beta)
\]
\(=.=\) together with \(=*=\) form the map pattern.
\((=*=):: c(\mathrm{a}->\mathrm{b})->\mathrm{c} \mathrm{a}->\mathrm{c} \mathrm{b}\)
Atom which applies the functions contained by as structure (i.e. categorical type), on the elements of another structure, defined as:
\[
\Leftrightarrow: \mathcal{C}(\alpha \rightarrow \beta) \rightarrow \mathcal{C}(\alpha) \rightarrow \mathcal{C}(\beta)
\]
\(=.=\) together with \(=*=\) form the map pattern.
\((=\backslash=)::(\mathrm{a}->\mathrm{a}->\mathrm{a})\)-> c a \(->\mathrm{a}\)
Atom which reduces a structure to an element based on an associative function, defined as:
\[
\otimes:(\alpha \times \alpha \rightarrow \alpha) \rightarrow \mathcal{C}(\alpha) \rightarrow \alpha
\]
(=«=)
Skeleton which pipes an element through all the functions contained by a structure. N.B.: this is not an atom. It has an implicit definition which might be augmented by instances of this class to include edge cases.
\[
\begin{aligned}
& \forall: \mathcal{C}(\alpha \rightarrow \alpha) \times \alpha \rightarrow \alpha \\
& \diamond=(\circ) \diamond
\end{aligned}
\]

As the composition operation is not associative, we cannot treat pipe as a true reduction. Alas, it can still be exploited in parallel since it exposes another type of parallelism: time parallelism.
first : : c a -> a
Returns the first element in a structure.N.B.: this is not an atom. It has an implicit definition which might be replaced by instances of this class with a more efficient implementation.
last :: c a -> a
Returns the last element in a structure.N.B.: this is not an atom. It has an implicit definition which might be replaced by instances of this class with a more efficient implementation.
\[
\begin{aligned}
& \text { first }_{S}: \mathcal{C}(\alpha) \rightarrow \alpha \\
& \text { first }_{S}=(\lambda a b \rightarrow a) \otimes
\end{aligned}
\]
\[
\begin{aligned}
& \operatorname{last}_{S}: \mathcal{C}(\alpha) \rightarrow \alpha \\
& \text { last }_{S}=(\lambda a b \rightarrow b) \diamond
\end{aligned}
\]
instance Skeleton Vector
Ensures that Vector is a structure associated with the Skeleton Layer.

\subsection*{4.13.2 Skeleton constructors}

Patterns of in the skeleton layer are provided, like all other patterns in ForSyDe-Atom, as constructors. If the layer below this one is the MoC layer, i.e. the functions taken as arguments are processes, then these skeletons can be regarded as process network constructors, as the structures created are process networks with inherent potential for parallel implementation.
farm22 :: Skeleton c =>
( \(\mathrm{a} 1->\mathrm{a} 2->(\mathrm{b} 1, \mathrm{~b} 2)\) ) \(->\mathrm{c}\) a1 -> c a2 -> (c b1, c b2)
farm maps a function on a vector. It is the embodiment of the map homomorphism, and its naming is inspired from the pattern predominant in HPC. Indeed, if we consider the layer below as being the MoC layer (i.e. the passed functions are processes), the resulting structure could be regarded as a "farm of data-parallel processes".
Constructors: farm[1-8] [1-4].
\[
\begin{aligned}
& \diamond:\left(V^{n} \rightarrow V^{m}\right) \rightarrow \mathcal{C}^{n} \rightarrow \mathcal{C}^{m} \\
& f \diamond\left(c_{1}, c_{2}, \ldots c_{n}\right)=f \diamond c_{1} \diamond c_{2} \diamond \ldots \diamond c_{n} \triangleleft
\end{aligned}
\]


\section*{reduce}
:: Skeleton c
=> (a -> a -> a) associative function (*)
-> c a structure
-> a reduced element
Infix name for the \(=\backslash=\) atom operator.
\(\left(^{*}\right)\) if the operation is not associative then the network can be treated like a pipeline.

\section*{reducei}
:: Skeleton c
=> (a -> a -> a) associative function (*)
-> a initial element of structure
-> c a structure
-> a reduced element
reducei is special case of reduce where an initial element is specified outside the reduced vector. It is implemented as a pipe with switched arguments, and the reduction function is constrained to be associative. It is semantically equivalent to the pattern depicted below.
\(\left(^{*}\right)\) if the operation is not associative then the network is semantically equivalent to pipe1 (see pipe2).
\[
\text { reducei }_{S}:(\alpha \times \alpha \rightarrow \alpha) \rightarrow \alpha \rightarrow \mathcal{C}(\alpha) \rightarrow \alpha
\]
\[
\text { reducei }_{S} f e\langle c\rangle=f \diamond\langle c\rangle \diamond e
\]
pipe
:: Skeleton c
\(\Rightarrow \quad c(a->a)\) vector of functions
\(->\) a kernel element
-> a result
Infix name for the \(=«=\) skeleton operator.
pipe2 :: Skeleton c =>
(a1 -> a2 -> a -> a) -> c a1 -> c a2 -> a -> a
The pipe constructors are a more generic form of the \(=«=\) (pipe) skeleton apt for successive partial application and create more robust parameterizable pipeline networks.
Constructors: comb [1-8].
\(\operatorname{pipe}_{S}:\left(\beta^{n} \rightarrow \alpha \rightarrow \alpha\right) \rightarrow \mathcal{C}\left(\beta^{n}\right) \rightarrow \alpha \rightarrow \alpha\)
\(\operatorname{pipe}_{S} f\langle c\rangle^{n} e=f \diamond\langle c\rangle^{n} \forall e\)


\subsection*{4.14 ForSyDe.Atom.Skeleton.Vector}
```

module ForSyDe.Atom.Skeleton.Vector (
Vector(Null, (:>)), null, unit, (<++>), vector, fromVector, indexes,
isNull, (<:), farm22, reduce, prefix, suffix, pipe, (=/=), recur,
cascade2, mesh2, length, index, fanout, fanoutn, generate, iterate,
first, last, inits, tails, init, tail, concat, reverse, group,
shiftr, shiftl, rotr, rotl, take, drop, takeWhile, filterIdx, odds
evens, stride, get, (<@), (<@!), gather1, (<@>), replace, scatter,
bitrev, duals, unduals, zipx, unzipx
) where

```

This module defines the data type Vector as a categorical type, and implements the atoms for the Skeleton class. Algorithmic skeletons for vector are mostly described in their factorized form, which ensures that they are catamorphisms (see the factorization theorem). Where efficiency or practicality is a concern, some skeletons are implemented as recurrences. One can still prove that they are catamorphisms through alternative theorems (see Skillicorn, 1994).

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.

\subsection*{4.14.1 Vector data type}

\section*{data Vector a}
\(=\) Null Null element. Terminates a vector.
I a (:>) (Vector a) appends an element at the head of a vector.
The Vector, or at least its interpretation, is the exact equivalent of an infinite list, as defined in R. Bird and Moor, 1997. Its name though is borrowed from Reekie, 1995, since it is more suggestive in the context of process networks.
According to R. Bird and Moor, 1997, Vector should be implemented as following:
data Vector \(\mathrm{a}=\mathrm{Null} \quad\) - null element
| Unit a - singleton vector
| Vector a <++> Vector a concatenate two vectors
This construction suggests the possibility of splitting a Vector into multiple parts and evaluating it in parallel. Due to reasons of efficiency, and to ensure that the structure is flat and homogeneous, vector is implemented using the same constructors as an infinite list like in R. S. Bird, 1987 (see below). When defining skeletons of vectors we will not use the real constructors though, but the theoretical ones defined above and provided as functions . This way we align ForSyDe-Atom's Vector type with the categorical type theory and its theorems.
Another particularity of vector is that it instantiates the reduction atom \(=\backslash=\) as a right fold, as it is the most efficient implementation in the context of lazy evaluation. As a consequence reduction is performed from right to left. This is noticeable especially in the case of pipeline-based skeletons (where pipe itself is a reduction with the right-associative composition operator .) is performed from right to left, which comes in natural when considering the order of function composition. Thus for reduce-based skeletons (e.g. prefix, suffix, recur, cascade, mesh) the result vectors shall be read from end to beginning.
```

instance Functor Vector

```

Provides an implementation for \(=.=\).
```

instance Applicative Vector

```

Provides an implementation for \(=*=\).

\section*{instance Foldable Vector}

Provides an implementation for \(=\backslash=\).

\section*{instance Skeleton Vector}

Ensures that Vector is a structure associated with the Skeleton Layer.
```

instance Eq a => Eq (Vector a)

```
instance Read a => Read (Vector a)

The vector \(1:>2:>\) Null is read using the string " \(<1,2>\) ".
```

instance Show a => Show (Vector a)

```

The vector \(1:>2:>\) Null is represented as \(<1,2\rangle\).
```

instance Plottable a => Plottable (Vector a)

```

Vectors of plottable types
```

instance Plottable a => Plot (Vector a)

```
vectors of coordinates

\subsection*{4.14.2 "Constructors"}

Theoretical constructors for the vector type, used in the definition of skeletons as catamorphisms.
null : : Vector a
Constructs a null vector.
\(\mid \lambda>\) null
<>
```

unit :: a -> Vector a

```

Constructs a singleton vector.
```

$\lambda>$ unit 1

```
<1>
(<++>) : : Vector a -> Vector a -> Vector a
Constructs a vector by appending two existing vectors.
\(\mid \lambda>\) unit \(1<++>\) unit 2
<1, 2>

\subsection*{4.14.3 Utilities}
```

vector :: [a] -> Vector a

```

Converts a list to a vector.
fromVector : : Vector a -> [a]
Converts a vector to a list.
indexes : : Vector Integer
Creates the infinite vector:
\(|<1,2,3,4, \ldots\rangle\)
Used mainly for operation on indexes.
```

isNull :: Vector a -> Bool

```

Returns True if the argument is a null vector.
```

(<:) :: Vector a -> a -> Vector a

```

Appends an element at the end of a vector.

\subsection*{4.14.4 Skeletons}

Algorithmic skeletons on vectors are mainly presented in terms of compositions of the atoms associated with the Skeleton Layer. When defining them, we use the following operators:
\[
\begin{align*}
\rangle & : \alpha \rightarrow \mathcal{V}(\alpha)  \tag{1}\\
+ & : \mathcal{V}(\alpha) \rightarrow \mathcal{V}(\alpha) \rightarrow \mathcal{V}(\alpha)  \tag{2}\\
v[n] & : \mathcal{V}(\alpha) \times \operatorname{Int} \rightarrow \alpha  \tag{3}\\
v[2 . .] & : \mathcal{V}(\alpha) \times \mathcal{V}(\text { Int }) \rightarrow \mathcal{V}(\alpha)  \tag{4}\\
\infty & : \alpha \rightarrow \mathcal{V}(\alpha) \tag{5}
\end{align*}
\]
where:
- (1) is the unit constructor, constructing a singleton vector.
- (2) is the <++> constructor, concatenating two vectors.
- (3) is the <@!> selector. The subscript notation is used to denote element at position \(n\) in a vector.
- (4) suggests an arbitrary selector which returns a vector with another one's elements, based on some indices. The shown example is an alternative notation for the tail skeleton.

\section*{Functional networks}

This sub-category denotes skeletons (patterns) which are take functions as arguments. If the functions are MoC layer entities, i.e. processes, then these patterns are capable of constructing parallel process networks. Using the applicative mechanism, the designer has a high degree of freedom when customizing process networks through systematic partial application, rendering numerous possible usages for the same pattern. To avoid over-encumbering the figures, they depict small test cases, which might not expose the full potential of the constructors.
see the naming convention rules on how to interpret, use and develop your own constructors.
```

farm22

```
\begin{tabular}{lll}
\(::\) & (a1 -> a2 -> (b1, b2)) & function (e.g. process) \\
-> & Vector a1 & first input vector \\
-> & Vector a2 & second input vector \\
-> & (Vector b1, Vector b2) & two output vectors
\end{tabular}
farm is simply the Vector instance of the skeletom farm pattern (see farm22). If the function taken as argument is a process, then it creates a farm network of data parallel processes.
Constructors: farm[1-4] [1-4].
```

\lambda> let v1 = vector [1,2,3,4,5]
\lambda> S.farm21 (+) v1 v1
<2,4,6,8,10>
\lambda> let s1 = SY.signal [1,2,3,4,5]
\lambda> let v2 = vector [s1,s1,s1]
\lambda> S.farm11 (comb11 (+1)) v2
<{2,3,4,5,6},{2,3,4,5,6},{2,3,4,5,6}>
\lambda> S.farm21 (\x -> comb11 (+x)) v1 v2
<{2,3,4,5,6},{3,4,5,6,7},{4, 5,6,7,8}>

```
    \(\operatorname{farm}_{S}:\left(\alpha^{m} \rightarrow \beta^{n}\right) \rightarrow\langle\alpha\rangle^{m} \rightarrow\langle\beta\rangle^{n}\)
    \(\operatorname{farm}_{S}: \ominus\)

reduce : : ( \(\mathrm{a}->\mathrm{a}->\mathrm{a}\) ) \(->\) Vector \(\mathrm{a}->\mathrm{a}\)
As the name suggests, it reduces a vector to an element based on an associative function. If the function is not associative, it can be treated like a pipeline.

Vector instantiates the skeletons for both reduce and reducei.
```

\lambda> let v1 = vector [1,2,3,4,5]
\lambda> S.reduce (+) v1
15
\lambda> let s1 = SY.signal [1,2,3,4,5]
\lambda> let s2 = SY.signal [10,10,10,10,10]
\lambda> let v2 = vector [s1,s1,s1]
\lambda> S.reduce (comb21 (+)) v2
{3,6,9,12,15}
\lambda> S.reducei (comb21 (+)) s2 v2
{13,16,19,22,25}

```

prefix :: (b -> b \(->\) b) -> Vector b \(->\) Vector b
prefix peforms the parallel prefix operation on a vector. Equivalent process networks are constructed if processes are passed as arguments.
Similar to reduce and reducei, two versions prefix and prefixi are provided.
```

\lambda> let v1 = vector [1,2,3,4,5]
\lambda> prefix (+) v1
<15,14,12,9,5>
\lambda> let s1 = SY.signal [1,2,3,4,5]
\lambda> let s2 = SY.signal [10,10,10,10,10]
\lambda> let v2 = vector [s1,s1,s1]
\lambda> prefix (comb21 (+)) v2
<{3,6,9,12,15},{2,4,6,8,10},{1,2,3,4,5}>
\lambda> prefixi (comb21 (+)) s2 v2
<{13,16,19,22,25},{12,14,16,18,20},{11,12,13,14,15}>

```
prefix \(_{S}:(\alpha \times \alpha \rightarrow \alpha) \rightarrow\langle\alpha\rangle \rightarrow\langle\alpha\rangle\)
prefix \(_{S} f\langle a\rangle=(f \diamond) \diamond\left(\right.\) tails \(\left._{S}\langle a\rangle\right)\)


suffix : : (b -> b -> b) -> Vector b -> Vector b
suffix peforms the parallel suffix operation on a vector. Equivalent process networks are constructed if processes are passed as arguments.
Similar to reduce and reducei, two versions suffix and suffixi are provided.
```

\lambda> let v1 = vector [1,2,3,4,5]
\lambda> suffix (+) v1
<1,3,6,10,15>
\lambda> let s1 = SY.signal [1,2,3,4,5]
\lambda> let s2 = SY.signal [10,10,10,10,10]
\lambda> let v2 = vector [s1,s1,s1]
\lambda> suffix (comb21 (+)) v2
<{1,2,3,4,5},{2,4,6,8,10},{3,6,9,12,15}>
\lambda> suffixi (comb21 (+)) s2 v2
<{11,12,13,14,15},{12,14,16,18,20},{13,16,19,22,25}>

```
                    suffix \(_{S}:(\alpha \times \alpha \rightarrow \alpha) \rightarrow\langle\alpha\rangle \rightarrow\langle\alpha\rangle\)
                    \(\operatorname{suffix}_{S} f\langle a\rangle=(f \otimes) \mapsto\left(\right.\) inits \(\left._{S}\langle a\rangle\right)\)

pipe
Vector (a -> a) vector of functions
input
output
-> a output
pipe creates a pipeline of functions from a vector. pipe simply instantiates the \(=«=\) atom whereas pipeX instantiate their omologi from the ForSyDe.Atom. Skeleton module (see pipe2).
OBS: the pipelining is done in the order dictated by the function composition operator: from right to left.
Constructors: pipe [1-4].
```

\lambda> let v1 = vector [(+1),(+1),(+1)]
\lambda> S.pipe v1 1
4
\lambda> let s1 = SY.signal [1,2,3,4]
\lambda> let v2 = vector [1,2,3,4]
\lambda> S.pipe1 (\x -> comb11 (+x)) v2 s1
{11,12,13,14}

```

```

(=/=) :: Vector (a -> a) -> a -> Vector a

```

Infix operator for recur.
recur
\begin{tabular}{lll}
\(::\) & Vector (a -> a) & vector of functions \\
-> & a & input \\
-> & Vector a & output
\end{tabular}
recur creates a systolic array from a vector of functions. Just like pipe and pipeX, there exists a raw recur version with an infix operator \(=/=\), and the enhanced recurX which is meant for systematic partial application of a function on an arbitrary number of vectors until the desired vector of functions is obtained.

Constructors: (=/=), recur, recuri, recur [1-4] [1-4].
```

\lambda> let v1 = vector [(+1),(+1),(+1)]
\lambda> recur v1 1
<4,3,2>
\lambda> recuri v1 1
<4,3,2,1>
\lambda> let s1 = SY.signal [1,2,3,4]
\lambda> let v2 = vector [1,2,3,4]
\lambda> recur1 (\x -> comb11 (+x)) v2 s1
<{11,12,13,14},{10,11,12,13},{8,9,10,11},{5,6,7,8}>

```
cascade2
                                    \(\diamond:\langle\alpha \rightarrow \alpha\rangle \rightarrow \alpha \rightarrow\langle\alpha\rangle\)
                                    \(\langle f\rangle \diamond i=\left(\diamond_{i}\right) \otimes\left(\operatorname{tail}_{\mathrm{s}_{S}}\langle f\rangle\right)\)

: : (a2 \(->\) a1 \(->\) a \(->\) a \(->\) a) function41 which needs to be applied to function21
-> Vector (Vector a2) fills in the first argument in the function above fills in the second argument in the function above
-> Vector a first input vector (e.g. of signals)
-> Vector a second input vector (e.g. of signals)
-> Vector a output
cascade creates a "cascading mesh" as a result of piping a vector into a vector of recur arrays.
Constructors: cascade, cascade [1-4].
```

\lambda> let v1 = vector [1,2,3,4]
> cascade (+) v1 v1
<238,119,49,14>
\lambda> let s1 = SY.signal [1,2,3,4]
\lambda> let vs = vector [s1, s1, s1]
\lambda> cascade (comb21 (+)) vs vs
<{20,40,60, 80},{10,20,30,40},{4,8,12,16}>
\lambda> let vv = vector [vector [1,-1,1], vector [-1,1,-1], vector [1,-1,1] ]
\lambda> cascade1 (\x -> comb21 (\y z-> x*(y+z))) vv vs vs
<{16,32,48,64},{8,16,24,32},{-2,-4,-6,-8}>

```
```

cascade }S:(\alpha->\alpha->\alpha)->\langle\alpha\rangle->\langle\alpha\rangle->\langle\alpha

```
cascade }S:(\alpha->\alpha->\alpha)->\langle\alpha\rangle->\langle\alpha\rangle->\langle\alpha
cascadeS p \langles\mp@subsup{\rangle}{1}{}\langles\mp@subsup{\rangle}{2}{}=\mp@subsup{\operatorname{scanf}}{S}{}\otimes\langles\mp@subsup{\rangle}{2}{}}\forall\langles\mp@subsup{\rangle}{1}{
cascadeS p \langles\mp@subsup{\rangle}{1}{}\langles\mp@subsup{\rangle}{2}{}=\mp@subsup{\operatorname{scanf}}{S}{}\otimes\langles\mp@subsup{\rangle}{2}{}}\forall\langles\mp@subsup{\rangle}{1}{
where scanf}\mp@subsup{\mp@code{S}}{2}{}\mp@subsup{s}{2}{}\langles\mp@subsup{\rangle}{1}{}=p\forall\langles\mp@subsup{\rangle}{1}{}\diamond\mp@subsup{s}{2}{
```

where scanf}\mp@subsup{\mp@code{S}}{2}{}\mp@subsup{s}{2}{}\langles\mp@subsup{\rangle}{1}{}=p\forall\langles\mp@subsup{\rangle}{1}{}\diamond\mp@subsup{s}{2}{

```

mesh2
```

:: (a2 -> a1 -> a -> a -> a)
-> Vector (Vector a2)
-> Vector (Vector a1)
-> Vector a
-> Vector a
-> Vector (Vector a)

```
function41 which needs to be applied to function21 fills in the first argument in the function above fills in the second argument in the function above first input vector (e.g. of signals)
second input vector (e.g. of signals)
output, a 2D vector
mesh creates a 2D systolic array as a result of piping a vector into a vector of 1 D systolic arrays.

Constructors: mesh, mesh [1-4].
```

\lambda> let v1 = vector [1,2,3,4]
\lambda> mesh (+) v1 v1
<<238,119,49,14> ,<119,70,35,13>,<49,35,22,11>,<14,13,11,8>>
\lambda> let s1 = SY.signal [1,2,3,4]
\lambda> let vs = vector [s1, s1, s1]
\lambda> mesh (comb21 (+)) vs vs
<<{20,40,60,80},{10,20,30,40},{4,8,12,16}>,<{10,20,30,40},{6,12, 18,24},{3,6,9,12}>,<{4,8,12,16},
{3,6,9,12},{2,4,6,8}>>
\lambda> let vv = vector [vector [1,-1,1], vector [-1,1,-1], vector [1,-1,1]]
\lambda> mesh1 (\x -> comb21 (\y z-> x*(y+z))) vv vs vs
<<{16,32,48,64},{8,16,24,32},{-2,-4,-6,-8}>,<{8,16,24,32},{-6, -12,-18,-24},{-3,-6, -9, -12}>,
<{-2,-4,-6,-8},{-3,-6,-9,-12},{2,4,6,8}>>

```
                                    mesh \(_{S}:(\alpha \rightarrow \alpha \rightarrow \alpha) \rightarrow\langle\alpha\rangle \rightarrow\langle\alpha\rangle \rightarrow\langle\langle\alpha\rangle\rangle\)
                                    \(\operatorname{mesh}_{S} p\langle s\rangle_{1}\langle s\rangle_{2}=\operatorname{scanf}_{S} \forall\langle s\rangle_{2} \diamond\langle s\rangle_{1}\)
                        where \(\operatorname{scanf}_{S} s_{2}\langle s\rangle_{1}=p \mapsto\langle s\rangle_{1} \diamond s_{2}\)


\section*{Queries}

Queries return various information about a vector. They are also built as skeletons.
length : : Num p => Vector a \(->\mathrm{p}\) returns the number of elements in a value.
\(\lambda>\) length \(\$\) vector \([1,2,3,4,5]\)
5
```

length $_{S}:\langle\alpha\rangle \rightarrow$ Int
length $_{S}=(+\diamond) \circ(1 \diamond)$

```
index :: Vector a2 -> Vector Integer
returns a vector with the indexes from another vector.
\(\mid \lambda>\) index \(\$\) vector \([1,1,1,1,1,1,1]\)
<1,2,3,4,5,6,7>

\section*{Generators}

Generators are specific applications of the prefix or suffix skeletons.
fanout : : t -> Vector t
fanout repeats an element. As a process network it distributes the same value or signal to all the connected processes down the line. Depending on the target platform and the refinement decisions involved, it may be interpreted in the following implementations:
- global or shared memory in case of a massively parallel platform (e.g. GPU)
- a (static) memory or cache location in memory-driven architectures (e.g. CPU)
- a fanout in case of a HDL system
- a broadcast in case of a distributed system
fanoutn : : (Num t, Ord t) \(\Rightarrow\) t \(\rightarrow\) a \(->\) Vector a
fanoutn is the same as fanout, but the length of the result is also provided.
\(\lambda>\) fanoutn 51
<1,1,1,1,1>
generate : : (Num t, Ord t) \(\Rightarrow \mathrm{t} \rightarrow\) ( \(\mathrm{a}->\mathrm{a}\) ) \(->\mathrm{a}->\) Vector a
generate creates a vector based on a kernel function. It is just a restricted version of recur.
\(\mid \lambda>\) generate \(5(+1) 1\)
<6,5,4,3,2>
\[
\begin{aligned}
& \text { generate }_{S}: \text { Int } \rightarrow(\alpha \rightarrow \alpha) \rightarrow \alpha \rightarrow\langle\alpha\rangle \\
& \text { generate }_{S} n f i=\langle f\rangle_{\times n} \diamond_{i}
\end{aligned}
\]
iterate : : (Num t, Ord t) => t -> (a -> a) -> a -> Vector a
iterate is a version of generate which keeps the initial element as well. It is a restricted version of recuri.
```

    \lambda> iterate 5 (+1) 1
    <5,4,3,2,1>
    ```

\section*{Permutators}

Permutators perform operations on the very structure of vectors, and make heavy use of the vector constructors.
first : : Vector a -> a
Instance of first
\(\mid \lambda>\) S.first \(\$\) vector \([1,2,3,4,5]\)
1
last : : Vector a -> a
Instance of last
| \(\lambda>\) S.last \(\$\) vector \([1,2,3,4,5]\)
5
inits : : Vector a -> Vector (Vector a)
creates a vector of all the initial segments in a vector.
\(\mid \lambda>\) inits \(\$\) vector \([1,2,3,4,5]\)
\(\mid\langle<1\rangle,\langle 1,2\rangle,\langle 1,2,3\rangle,\langle 1,2,3,4\rangle,\langle 1,2,3,4,5\rangle>\)
\[
\begin{aligned}
& \operatorname{inits}_{S}:\langle\alpha\rangle \rightarrow\langle\langle\alpha\rangle\rangle \\
& \text { inits }_{S} n=(\operatorname{sel} \diamond) \circ(\langle\langle \rangle\rangle \otimes)
\end{aligned}
\]
\[
\text { where sel } x y=x+((x[L]+) \oplus y)
\]

tails : : Vector a -> Vector (Vector a)
creates a vector of all the final segments in a vector.
\(\mid \lambda>\) tails \(\$\) vector \([1,2,3,4,5]\)
\(|\ll 1,2,3,4,5\rangle,\langle 2,3,4,5\rangle,\langle 3,4,5\rangle,\langle 4,5\rangle,\langle 5 \gg\)
\[
\begin{aligned}
& \operatorname{tails}_{S}:\langle\alpha\rangle \rightarrow\langle\langle\alpha\rangle\rangle \\
& \operatorname{tails}_{S} n=(\operatorname{sel} \diamond) \circ(\langle\langle \rangle\rangle \diamond)
\end{aligned}
\]
where sel \(x y=((+y[1]) \diamond x)+y\)

init : : Vector a -> Vector a
Returns the initial segment of a vector.
\(\mid \lambda>\) init \(\$\) vector \([1,2,3,4,5]\)
<1,2,3,4>
\[
\begin{aligned}
& \operatorname{init}_{S}:\langle\alpha\rangle \rightarrow\langle\alpha\rangle \\
& \operatorname{init}_{S}=\operatorname{inits}_{S}[L-1]
\end{aligned}
\]
tail :: Vector a -> Vector a
Returns the tail of a vector.
| \(\lambda>\) tail \(\$\) vector \([1,2,3,4,5]\)
<2,3,4,5>
\[
\begin{aligned}
& \operatorname{tail}_{S}:\langle\alpha\rangle \rightarrow\langle\alpha\rangle \\
& \operatorname{tail}_{S}=\operatorname{tails}_{S}[2]
\end{aligned}
\]
concat : : Vector (Vector a) -> Vector a concatenates a vector of vectors.
\(\mid \lambda>\) concat \(\$\) vector [vector \([1,2,3,4]\), vector \([5,6,7]\) ]
<1,2,3,4,5,6,7>
\[
\begin{aligned}
& \operatorname{concat}_{S}:\langle\langle\alpha\rangle\rangle \rightarrow\langle\alpha\rangle \\
& \operatorname{concat}_{S}=\# \otimes
\end{aligned}
\]
reverse : : Vector a -> Vector a
reverses the elements in a vector.
\(\mid \lambda>\) reverse \(\$\) vector \([1,2,3,4,5]\)
\(\langle 5,4,3,2,1>\)

where rev \(x y=y+x\)

group :: Integer -> Vector a -> Vector (Vector a) groups a vector into sub-vectors of \(n\) elements.
\(\left\lvert\, \begin{aligned} & \lambda\rangle \text { group } 3 \$ \text { vector }[1,2,3,4,5,6,7,8] \\ & \langle\langle 1,2,3\rangle,\langle 4,5,6\rangle\langle 7,8\rangle\rangle\end{aligned}\right.\)
<<1,2,3>,<4,5,6>,<7,8>>
\[
\begin{aligned}
& \operatorname{group}_{S}: \text { Int } \rightarrow\langle\alpha\rangle \rightarrow\langle\langle\alpha\rangle\rangle \\
& \operatorname{group}_{S} n=\left(\operatorname{take}_{S} n\right) \diamond \circ\left(\operatorname{drop}_{S} n\right) \diamond
\end{aligned}
\]

shiftr : : Vector a -> a -> Vector a
right-shifts a vector with an element.
```

\lambda> vector [1,2,3,4] 'shiftr' 8
<8,1,2,3>

```


shiftl : : Vector a -> a -> Vector a
left-shifts a vector with an element.
```

|> vector [1,2,3,4] 'shiftl' 8

```
<2,3,4,8>

rotr : : Vector a -> Vector a
rotates a vector to the right.
\(\mid \lambda>\) rotr \(\$\) vector \([1,2,3,4]\)
<4,1,2,3>

rotl : : Vector a -> Vector a rotates a vector to the left.

\(\mid \lambda>\) rotl \(\$\) vector \([1,2,3,4]\)
<2,3,4,1>

take : : Integer \(->\) Vector a -> Vector a takes the first \(n\) elements of a vector.
\(\mid \lambda>\) take \(5 \$\) vector \([1,2,3,4,5,6,7,8,9]\)
<1,2,3,4,5>
\[
\operatorname{take}_{S}:\langle\alpha\rangle \rightarrow\langle\alpha\rangle
\]
\(\operatorname{take}_{S} n=(\operatorname{sel} \otimes) \circ(\langle \rangle \otimes)\)
where sel \(x y= \begin{cases}x+y & \text { if } \operatorname{index}(x)<n \\ x & \text { otherwise }\end{cases}\)
drop :: Integer -> Vector a -> Vector a
drops the first \(n\) elements of a vector.
\(\mid \lambda>\) drop \(5 \$\) vector \([1,2,3,4,5,6,7,8,9]\)
<6,7,8,9>
\[
\begin{aligned}
& \operatorname{drop}_{S}:\langle\alpha\rangle \rightarrow\langle\alpha\rangle \\
& \operatorname{drop}_{S} n=(\operatorname{sel} \diamond) \circ(\langle \rangle \diamond)
\end{aligned}
\]
\[
\text { where sel } x y= \begin{cases}x+y & \text { if } \operatorname{index}(x)>n \\ y & \text { otherwise }\end{cases}
\]
takeWhile : : (a -> Bool) -> Vector a -> Vector a
takes the first elements in a vector until the first element that does not fulfill a predicate.
\(\mid \lambda>\) takeWhile (<5) \$ vector \([1,2,3,4,5,6,7,8,9]\)
<1,2,3,4>
```

takeWhile ${ }_{S}:\langle\alpha\rangle \rightarrow\langle\alpha\rangle$
takeWhile $_{S} f=\operatorname{concat}_{S} \circ(\operatorname{sel} \diamond) \circ(\langle\langle \rangle\rangle \ominus)$

```
where sel \(x y= \begin{cases}x+y & \text { if } f(y[1])=\text { true } \wedge \exists x[L] \\ x & \text { otherwise }\end{cases}\)
filterIdx :: (Integer -> Bool) -> Vector a -> Vector a returns a vector containing only the elements of another vector whose index satisfies a predicate.
\(\mid \lambda>\) filterIdx ( \(\backslash x->x\) 'mod' \(3==0\) ) \(\$\) vector \([0,1,2,3,4,5,6,7,8,9]\)
<2,5,8>
filterIdx \(_{S}:\langle\alpha\rangle \rightarrow\langle\alpha\rangle\)
filterIdx \(_{S} f=(\operatorname{sel} \otimes) \circ(\langle \rangle \otimes)\)
where sel \(x y= \begin{cases}x+y & \text { if } f(\operatorname{index}(x))=\text { true } \\ y & \text { otherwise }\end{cases}\)

odds :: Vector a -> Vector a
\[
\text { odds }_{S}=\text { filterIdx }_{S}(\text { odd })
\]
evens :: Vector a -> Vector a

\section*{stride}
:: Integer first index
-> Integer stride length
-> Vector a
Vector a
does a stride-selection on a vector.
\(\lambda>\) stride \(13 \$\) vector \([1,2,3,4,5,6,7,8,9]\)
<1,4,7>
stride \(_{S}:\langle\alpha\rangle \rightarrow\langle\alpha\rangle\)
stride \(_{S}\) first stride \(=(\operatorname{sel} \diamond) \circ(\langle \rangle \oplus)\)
where sel \(x y=\left\{\begin{array}{ll}x+y & \text { if }(\operatorname{index}(x)-\text { first }) \\ y & \text { otherwise }\end{array}\right.\) mod stride \(=0\)

get :: Integer -> Vector a -> Maybe a
returns the \(n\)-th element in a vector, or Nothing if \(n>l\).
\(\mid \lambda>\) get \(3 \$\) vector \([1,2,3,4,5]\)
Just 3
\[
\begin{aligned}
& \operatorname{get}_{S}:\langle\alpha\rangle \rightarrow \alpha \\
& \operatorname{get}_{S} n=(\operatorname{sel} \diamond)
\end{aligned}
\]
where sel \(x y= \begin{cases}x & \text { if } \operatorname{index}(x)=n \\ y & \text { otherwise }\end{cases}\)
(<@) :: Vector a -> Integer -> Maybe a
the same as get but with flipped arguments.
(<@!) :: Vector p -> Integer -> p
unsafe version of <@>. Throws an exception if \(n>l\).

\section*{gather1}
```

:: Vector Integer vector of indexes
-> Vector a input vector
-> Vector (Maybe a)

```
selects the elements in a vector at the incexes contained by another vector.
The following versions of this skeleton are available, the number suggesting how many nested vectors it is operating upon: gather [1-5]
```

\lambda> let ix = vector [vector [1,3,4], vector [3,5,1], vector [5,8,9]]
\lambda> let v = vector [11,12,13,14,15]
\lambda> gather2 ix v
<<Just 11,Just 13,Just 14>,<Just 13,Just 15,Just 11>,<Just 15,Nothing,Nothing>>

```
gather \(_{S}:\langle\) Int \(\rangle\langle\alpha\rangle \rightarrow\langle\alpha\rangle\)
gather \(_{S}\langle i x\rangle\langle v\rangle=\operatorname{sel} \mapsto\langle i x\rangle\)
                        where sel \(x=\langle v\rangle[x]\)

(<@>)

\section*{\(:\) Vector a input vector}
-> Vector Integer vector of indexes
-> Vector (Maybe a)
the same as gather 1 but with flipped arguments
The following versions of this skeleton are available, the number suggesting how many nested vectors it is operating upon.

replace :: Integer -> p -> Vector p -> Vector p
replaces the \(n\)-th element in a vector with another.
| \(\lambda>\) replace \(515 \$\) vector \([1,2,3,4,5,6,7,8,9]\)
\(\langle 1,2,3,4,15,6,7,8,9>\)
\[
\begin{aligned}
& \text { replace }_{S}:\langle\alpha\rangle \rightarrow\langle\alpha\rangle \\
& \text { replace }_{S} n r=(\operatorname{sel} \otimes) \circ(\langle \rangle \otimes) \\
& \qquad \text { where sel } x y= \begin{cases}\langle r\rangle+y & \text { if } \text { index }(x)=n \\
x+y & \text { otherwise }\end{cases}
\end{aligned}
\]

scatter : : Vector Integer \(->\) Vector p \(->\) Vector p -> Vector p scatters the elements in a vector based on the indexes contained by another vector.
\(\lambda>\) scatter (vector \([2,4,5]\) ) (vector \([0,0,0,0,0,0,0,0]\) ) (vector \([1,1,1])\)
\(\langle 0,1,0,1,1,0,0,0\rangle\)
\[
\begin{aligned}
& \text { scatter }_{S}:\langle\text { Int }\rangle\langle\alpha\rangle \rightarrow\langle\alpha\rangle \rightarrow\langle\alpha\rangle \\
& \text { scatter }_{S}\langle i x\rangle\langle h o s t\rangle=(\text { sel } \diamond) \circ(\langle \rangle \otimes) \\
& \quad \text { where sel } x=\text { replace }_{S} \text { index }(x) x[1]\langle h o s t\rangle
\end{aligned}
\]

bitrev : : Vector a -> Vector a performs a bit-reverse permutation.

| \(\lambda>\) bitrev \(\$\) vector ["000", "001", "010", "011", "100", "101", "110", "111"]
<"111", "011", "101", "001", "110", "010", "100", "000">
duals : : Vector b2 \(->\) (Vector b2, Vector b2)
splits a vector in two equal parts.
```

    \lambda> duals $ vector [1, 2,3,4,5,6,7]
    (<1, 2, 3>, <4, 5,6>)
    ```
unduals :: Vector a -> Vector a -> Vector a
concatenates a previously split vector. See also duals

\section*{Interfaces}
zipx
:: \(\quad \mathrm{MoCe}\)
\(\Rightarrow\) Vector ((Vector a -> Vector a -> Vector a) -> Fun e (Vector a) (Fun e (Vector a) (Ret e (Vector a))) )
-> Vector (Stream (e a))
-> Stream (e (Vector a))
vector of MoC-specific context wrappers for the function <++>
input vector of signals output signal of vectors
zipx is a template skeleton for "zipping" a vector of signals. It synchronizes all signals (of the same MoC ) in a vector and outputs one signal with vectors of the synced values. For each signal in the input vector it requires a function which translates a partition of events (see ForSyDe.Atom.MoC) into sub-vectors.
There exist helper intances of the zipx skeleton interface for all supported MoCs.
\[
\begin{aligned}
& \operatorname{zipx}_{S}:\langle\langle\dot{\alpha}\rangle \times\langle\dot{\alpha}\rangle \rightarrow\langle\dot{\alpha}\rangle\rangle \rightarrow\langle\mathcal{S}(\alpha)\rangle \rightarrow \mathcal{S}(\langle\alpha\rangle) \\
& \operatorname{zipx}_{S}\langle\text { part }\rangle=((+\oplus) \diamond\langle\text { part }\rangle) \circ\left(\text { unit }_{S} \diamond\right)
\end{aligned}
\]

```

unzipx :: MoC e =>
(Vector a -> Vector (Ret e a))
-> Integer -> Stream (e (Vector a)) -> Vector (Stream (e a))

```
unzipx is a template skeleton to unzip a signal carrying vectors into a vector of multiple signals. It required a function that splits a vector of values into a vector of event partitions belonging to output signals. Unlike zipx, it also requires the number of output signals. The reason for this is that it is impossible to determine the length of the output vector without "sniffing" the content of the input events, which is out of the scope of skeletons and may lead to unsafe behavior. The length of the output vector is needed in order to avoid infinite recurrence.

There exist helper intances of the unzipx skeleton interface for all supported MoCs.
\[
\begin{aligned}
& \text { unzipx }_{S}:(\langle\alpha\rangle \rightarrow\langle\dot{\alpha}\rangle) \rightarrow \text { Int } \rightarrow \mathcal{S}(\langle\alpha\rangle) \rightarrow\langle\mathcal{S}(\alpha)\rangle \\
& \text { unzipx }_{S} \text { part } n=\left(\left(\text { first }_{S} \oplus\right) \oplus\right) \circ\left(\left\langle\operatorname{tail}_{S}\right\rangle_{\times n} \diamond\right) \circ(\operatorname{part} \oplus)
\end{aligned}
\]
fig/skel-vector-comm-unzipx.pdf

\subsection*{4.15 ForSyDe.Atom.Utility.Plot}
module ForSyDe.Atom.Utility.Plot
Config(Cfg, verbose, path, title, rate, xmax, labels, fire, other), defaultCfg, silentCfg, noJunkCfg, prepare, prepareL, prepareV, showDat, dumpDat, plotGnu, heatmapGnu, showLatex, dumpLatex, plotLatex, Plottable(toCoord), Plot(sample, sample', takeUntil, getInfo), PInfo(Info, typeid, command, measure, style, stacking, sparse), Samples, PlotData
) where

This module imports plotting and data dumping functions working with "plottable" data types, i.e. instances of the Plot and Plottable type classes.

\subsection*{4.15.1 User API}

The following commands are frequently used as part of the normal modeling routine.

\section*{Configuration settings}
```

data Config
= Cfg
> verbose :: Bool verbose printouts on terminal
> path :: String directory where all dumped files will be found
> title :: String base name for dumped files
> rate :: Float sample rate if relevant. Useful for explicit-tagged signals, ignored
otherwise.
> xmax :: Float Maximum X coordinate. Mandatory for infinite structures, optional
otherwise.
> labels:: list of labels with the names of the structures plotted
[String]
> fire :: Bool if relevant, fires a plotting or compiling program.
> other :: Bool if relevant, dumps additional scripts and plots.

```

Record structure containing configuration settings for the plotting commands.
```

instance Show Config

```
defaultCfg :: Config

Default configuration: verbose, dump everything possible, fire whatever program needed. Check source for settings.

Example usage:
```

\lambda> defaultCfg {xmax = 15, verbose = False, labels = ["john","doe"]}
Cfg {verbose = False, path = "./fig", title = "plot", rate = 1.0e-2, xmax = 15.0, labels = ["
john","doe"], fire = True, other = True}

```
silentCfg : : Config

Silent configuration: does not fire any program or print our unnecessary info. Check source for settings.
```

noJunkCfg :: Config

```

Clean configuration: verbose, does not dump more than necessary, fire whatever program needed. Check source for settings.

\section*{Data preparation}

\section*{prepare}
:: Plot a
\(\Rightarrow\) Config configuration settings
-> a plottable data type
-> PlotData structure ready for dumping
Prepares a single plottable data structure to be dumped and/or plotted.
```

prepareL :: Plot a => Config -> [a] -> PlotData

```

Prepares a list of plottable data structures to be dumped and/or plotted. See prepare.
```

prepareV :: Plot a => Config -> Vector a -> PlotData

```

Prepares a vector of plottable data structures to be dumped and/or plotted. See prepare.

\section*{Dumping and plotting data}
showDat :: PlotData -> IO ()
Prints out the sampled contents of a prepared data set.
dumpDat :: PlotData -> IO [String]
Dumps the sampled contents of a prepared data set into separate .dat files.
plotGnu :: PlotData -> IO ()
Generates a GNUplot script and . dat files for plotting the sampled contents of a prepared data set. Depending on the configuration settings, it also dumps LaTeX and PDF plots, and fires the script.
OBS: needless to say that GNUplot needs to be installed in order to use this command. Also, in order to fire GNUplot from a ghci session you might need to install gnuplot-x11.
```

heatmapGnu :: PlotData -> IO ()

```

Similar to plotGnu but creates a heatmap plot using the GNUplot engine. For this, the input needs to contain at least two columns of data, otherwise the plot does not show anything, i.e. the samples need to be lists or vectors of two or more elements.

OBS: same dependencies are needed as for plotGnu.
showLatex : : PlotData -> IO ()
Prints out a LaTeX environment from a prepared data set. This environment should be paste inside a tikzpicture in a document title which imports the ForSyDe-LaTeX package.
```

dumpLatex :: PlotData -> IO [String]

```

Dumps a set of formatted data files with the extension .flx that can be imported by a LaTeX document which uses the ForSyDe-LaTeX package.
```

plotLatex :: PlotData -> IO ()

```

Creates a standalone LaTeX document which uses the ForSyDe-LaTeX package, plotting a prepared data set. Depending on the configuration settings, the command pdflatex may also be invoked to compile a pdf image.
OBS: A LaTeX compiler is required to run the pdflatex command. The ForSyDe-LaTeX package also needs to be installed according to the instructions on the project web page.

\subsection*{4.15.2 The data types}

Below the data types involved are shown and the plottable structures are documented.

\section*{class Plottable a where}

This class gathers types which can be sampled and converted to a numerical string which can be read and interpreted by a plotter engine.

\section*{Methods}
toCoord : : a -> String
Transforms the input type into a coordinate string.
instance (Show a, Real a) => Plottable a
Real numbers that can be converted to a floating point representation
```

instance Plottable TimeStamp

```

Time stamps
instance Plottable a => Plottable (Vector a)
Vectors of plottable types
instance (Show a, Plottable a) => Plottable (AbstExt a)
Absent-extended plottable types
class Plot a where
This class gathers all ForSyDe-Atom structures that can be plotted.

\section*{Methods}
```

    sample :: Float -> a -> Samples
    ```

Samples the data according to a given step size.
sample :: a -> Samples ,
Samples the data according to the internal structure.
takeUntil :: Float -> a -> a
Takes the first samples until a given tag.
getInfo :: a -> PInfo
Returns static information about the data type.
```

instance Plottable a => Plot (Vector a)

```
    vectors of coordinates
instance Plottable a => Plot (Signal a)
    SY signals.
instance Plottable a => Plot (Signal a)
    SDF signals.
instance Plottable a => Plot (Signal a)
    DE signals.
instance Plottable a => Plot (Signal a)
    CT signals.

\section*{data PInfo}
\(=\) Info
> typeid : : String id used usually in implicit tags
\(>\) command :: String LaTeX identifier
> measure :: String unit of measure
> style :: String style tweaking in the GNUplot script
\(>\) stacking : : Bool if the plot is stacking
sparse :: Bool if the sampled data is sparse instead of dense
Static information of each plottable data type.
instance Show PInfo
type Samples = [(String, String)]
Alias for sampled data
type PlotData \(=(\) Config, PInfo, [(String, Samples)] \()\)
Alias for a data set prepared to be plotted.

\subsection*{4.16 ForSyDe.Atom.Utility.Tuple}
```

module ForSyDe.Atom.Utility.Tuple (

    at22, (||<), (<>), ($$)
    ) where
    ```

This module implements general purpose utility functions. It mainly hosts functions dealing with tuples. Utility are provided for up until 9-tuples. Follow the examples in the source code in case it does not suffice.

\section*{Reminder}

Make sure to consult naming conventions in section 4.1.1 in order to interpret the names and type signatures correctly.
at22 :: (a1, a2) -> a2
The at \(x y\) functions return the \(y\)-th element of an \(x\)-tuple.
ForSyDe.Atom.Utility exports the constructors below. Please follow the examples in the source code if they do not suffice:
```

at21, at22,
at31, at32, at33,
at41, at42, at43, at44,
at51, at52, at53, at54, at55,
at61, at62, at63, at64, at65, at66,
at71, at72, at73, at74, at75, at76, at77,
at81, at82, at83, at84, at85, at86, at87, at88,
at91, at92, at93, at94, at95, at96, at97, at98, at99,

```

\section*{Example:}
```

|> at53 (1,2,3,4,5)

```
3
(||<) :: (Functor f1, Functor f2) =>
    f1 (f2 (a1, a2)) -> (f1 (f2 a1), f1 (f2 a2))

This set of utility functions "unzip" nested \(n\)-tuples, provided as postfix operators. They are crucial for reconstructing data types from higher-order functions which input functions with multiple outputs. It relies on the nested types being instances of Functor.
The operator convention is \((1+<+)\), where the number of \(\mid\) represent the number of layers the n-tuple is lifted, while the number of \(<+1\) is the order \(n\) of the n-tuple.
ForSyDe.Atom.Utility exports the constructors below. Please follow the examples in the source code if they do not suffice:

\(||<, \quad|| \ll, \quad| | \lll, \quad| | \lll<, \quad| | \lll \ll, \quad| | \lll \lll, \quad| | \lll \lll<, \quad| | \lll \lll \lll\),



\section*{Example:}
\(\mid \lambda>\) : set -XPostfixOperators
\(\lambda>\) ([Just (1,2,3), Nothing, Just (4,5,6)] ||<<)
([Just 1, Nothing, Just 4], [Just 2, Nothing, Just 5], [Just 3, Nothing, Just 6])
(<>) :: (a1 -> a2 -> b1) -> (a1, a2) -> b1
Infix currying operators used for convenience.
The operator convention is (<>+), where the number of \(>+1\) is the order \(n\) of the \(n\)-tuple.
ForSyDe.Atom.Utility exports the constructors below. Please follow the examples in the source code if they do not suffice:
```

<<>, <>>, <>>>, <>>>>>, <>>>>>>, <>>>>>>>, <>>>>>>>>, <>>>>>>>>>>

```

Example:
```

|> (+) <> (1,2)

```
3
(\$\$) :: (a1 -> b1, a2 -> b2) -> (a1, a2) -> (b1, b2)
Infix function application operator for tuples.
The operator convention is (\$+), where the number of \(\$\) is the order \(n\) of the n-tuple. For Applying a function on nontuples we rely on \(\$\) provided by Prelude.
ForSyDe.Atom.Utility exports the constructors below. Please follow the examples in the source code if they do not suffice:
| \(\$\), \(\$ \$ \$\), \(\$ \$ \$ \$, \$ \$ \$ \$, \$ \$ \$ \$ \$, \$ \$ \$ \$ \$ \$, \$ \$ \$ \$ \$ \$ \$, \quad \$ \$ \$ \$ \$ \$ \$\)
Example:
```

|> ((+),(-)) $$
(1,1)
$$ (2,2)
(3,-1)
```

### 4.17 References

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[^0]:    ${ }^{1}$ currently available online at https://forsyde.github.io/forsyde-atom/
    ${ }^{2}$ available online at https://github.com/forsyde/forsyde-atom-examples

[^1]:    ${ }^{1}$ ForSyDe-Аtom v0.2.1

[^2]:    ${ }^{2} \Gamma_{\text {SDF }}=($ consumption rate for first input $[$, consumption rate for second input] $)$, production rate

[^3]:    ${ }^{1}$ Lee argues that the cost we pay in the loss of determinism is a property of the chosen modeling framework, and not of the model itself.

[^4]:    ${ }^{2}$ theoretical implications will be analyzed in the upcoming journal publications
    ${ }^{3}$ version 0.2.2

