

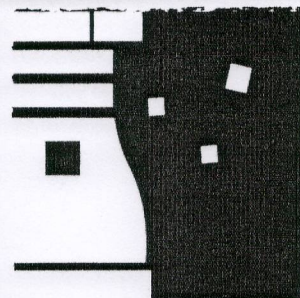
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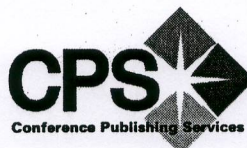
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## A Scientific Rapid Prototyping Model for the Haskell Language

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### Abstract

*A simple methodology for computational and scientific rapid prototyping in Haskell is proposed. The methodology is based on identifying abstractions that are defined over entire list data structures combined with identifying a set of processes that initially construct a list data structure and subsequently modify that data structure in iterative steps leading to a final post modification process to output the data structure. A case study exemplifies and details an application of the proposed model. The results indicate that the proposed methodology can be easily implemented for developers or coders with little in-depth knowledge of functional programming capabilities; thereby, enabling applicability for a wide range of users.*

### 1. Introduction

Nowadays, the computing infrastructure for computational sciences relies heavily on networked computers with on-demand access to high performance computers if and when needed. In addition, such computing infrastructures also include networked clusters, server-based clusters and in some cases, peer-to-peer clusters. There are many available options for the programming of these systems, depending on user choice and computing needs, for example, Java and related languages are commonly used for networked cluster based applications whereas FORTRAN and C are also commonly used for high-ended performance computations. Moreover, there are many libraries and packages that also can be incorporated into applications, for example, LAPACK, BLAS, VECLIB and MPI.

There are many computer resource, programming language, and library models that are available choices for the development and coding phases of computational applications. Many of these have individual conceptual models for procedure and data representations. In some cases, the models can integrate together; for example, typical

MPI-based parallel programming often includes master-slave farming (master-slave) and Single Program Multiple Data (SPMD) models combined and integrated with application code. Moreover, the models may have various abstractions, which in some cases, may have varying conceptual layouts, for example, OCC is well known to provide low-level point-to-point communication specification, the latter at the language level and the latter at the library level. However, models based on parallel programming based languages (e.g. FORTRAN-D system) or languages such as Linda or Haskell provide high-level specification requirements and more high-level 'implicit' concepts. A number of models are based on the idea of program skeletons, that is, parallel fragments. The complexity of choice is illustrated in Figure 1.

This paper follows from the earlier work where we have defined the term: *Programming models* to describe the viewpoint of "The parallel program is guided by at least one model, or guided by several." This paper continues the direction. Here, we are motivated by allowing choice upon the user for scientific and computational model development. This scientific rapid prototyping needs to be supported by model development fast development while allowing the developer to concentrate on core ideas. Hence, this paper is about the development time for research prototyping, allowing developers to more quickly produce computational models in the business, science and ubiquitous computing.

This rest of this paper is structured as follows. Section 2 describes the scientific rapid prototyping requirements. Section 3 presents the Haskell structures that we propose as suitable for computational and scientific rapid prototyping. Section 4 presents a case study application and conclusions in Section 5.



structure patterns that support rapid prototyping; in particular, that allow the developer to concentrate on *solution* as opposed to algorithm. Therefore, despite the general purpose capability and the many special functional capabilities of the language, we suggest simpler structures that we feel may be easier for the developer to program in, especially for those who do not have in-depth knowledge of functional languages. In this paper, we concentrate on the basic model of such structure patterns that is widely applicable to many programmed solutions, however, which might not be universally applicable.

Haskell's features of recursion and outer-layer pattern matching provide a nice compact semantic and grammar notation for tail recursion. Here, we introduce the term *take 1, drop 1* recursive processing of a list data structure to describe this pattern structure. There are three purposes that such a recursive procedure abstraction serves: 1) reduce a list to a scalar, 2) modify each element in a list (i.e. construct a new list), or 3) filter each element in a list (i.e. construct a sub-list). Note that any particular code fragment could combine the latter two. This structure uses the pattern specification ( $x:xs$ ) to implement the take 1, drop 1 concept; recursion terminating upon the pattern  $[]$ ; and explicit list construction using the  $:$  operator. Such a pattern requires processing individual data elements. The time required to code is comparative large.

Haskell also provides a list comprehension abstraction whereby a list construction can be specified by including three parts: the list element, the set of data inputs that collectively determine the output set and various filters over the inputs. This syntax is much more compact than the take 1, drop 1 structure, and can capture the latter two purposes of the take 1, drop 1 structure. Semantically, this structure still deals with element by element, however, in an overall single list manner.

Lastly, Haskell also provides a number of high level abstractions (and more experienced coders can implement user-defined abstractions). Examples of pre-existing functions in this category include: *fold* (and its variants), *map*, *zip* and its variants, and *filter*. Collectively such abstractions apply over an entire data structure.

Figure 3 illustrates the suggested relative abstractions of the above structures. Due to the preference of defining operations over a data structure as opposed to element by element, we determine the order of preference for coding for rapid prototyping: high level, then mid level, then low-level. High level functions can be implemented by low level abstractions; ultimately, all code can be implemented via the recursive structure.

The semantics of using higher level functions influences the way programs are developed. This leads us to propose a simple three stage algorithm methodology: 1) create a basic data structure representing the input, modify the basic data

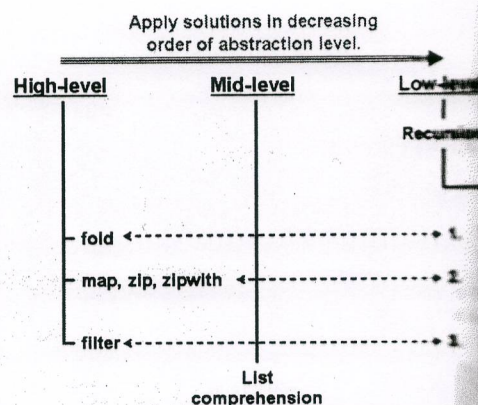


Figure 3. Abstraction levels of identified structure patterns

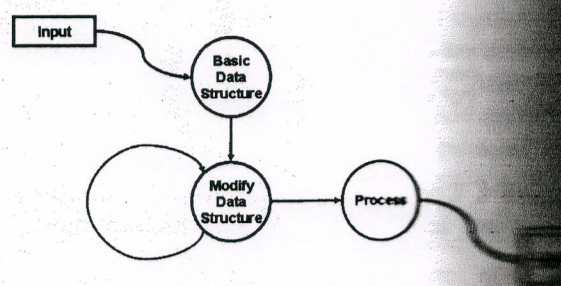


Figure 4.

structure, in some cases, by applying an iteration abstraction steps, 3) final process of the modified structure leading to the final output. This methodology is illustrated in Figure 4.

## 4 Application

We have applied the proposed model for the problem of palindrome identification in bioinformatics. The results indicate the potential benefit of the proposed model for rapid computational and scientific prototyping. In this section, we briefly show the application to the palindrome identification problem.

A palindrome in an RNA or DNA sequence is a cleotide sequence consisting of  $pgp'$  where  $p$  is a subsequence of length  $n$  and  $p'$  is the reverse complement of  $p$  also of length  $n$  and  $g$  is a subsequence gap between  $p$  and  $p'$ . The significance of identifying such patterns in a cleotide sequence is varied, for example, as motifs in secondary structure formations in RNA. In this paper, we use the proposed model to give a solution to the palindrome identification. The input sequence is represented as a list. The final output of the Haskell program is a list of palindromes.



Wiring

Second Substring

ctggac	253	258	gtccag
caagtc	223	228	gacttg
agtccc	194	199	gggact
ttacca	135	140	tggtaa
ctgcgg	96	101	ccgcag
ctgcgg	182	187	ccgcag
tgcggc	181	186	gccgca
gcgggt	180	185	agccgc
actcat	263	268	acgaac

```
build :: Int -> String -> [String]
build _ [] = []
build n (x:xs)
    | (length xs) >= (n-1) =
        (x:take (n-1) xs) : build n xs
    | otherwise             = []
```

The main part of `palindrome` is a list comprehension which implements both a modification and a filtration of the previous modified data structure. This further modifies the data structure to contain appropriate information necessary to the required output (see Figure 5) as well as filters the data structure to contain only identified palindromes (without duplication). At this point, in reference to Figure 4, this part of the process effects two transitions from ‘Modify Data Structure’ to ‘Modify Data Structure’ (i.e., there are two modifications when considered individually).

Lastly, as part of the palindrome identification filter in `palindrome`, the function `reverseComplement` is referenced which itself uses high level functions (e.g. `reverse` and `map`) to define  $p'$  as a single operation over the input subsequence string.

```
complement :: Char -> Char
complement c
  | c=='c'  = 'g'
  | c=='g'  = 'c'
  | c=='a'  = 't'
  | c=='t'  = 'a'
```

```
reverseComplement :: String -> String
reverseComplement s =
    (reverse . map complement) s
```

## 5 Conclusions

In this paper, we propose a simple methodology for computational and scientific rapid prototyping in Haskell. The methodology is based on identifying abstractions that are defined over entire list data structures combined with identifying a set of processes that initially construct a list data

**Final output of the Haskell program on the proposed model: the report of palindromes of length 6 for a portion of the hepatitis C virus NS5 gene, sequence obtained from the NCBI Nucleotide database, U0769711, and consists of 269 bases.**

is shown in Figure 5.

High level code is given below, in particular the functions `palindromeFormat` and `palindromeHeader` represent the final process of the methodology.

```
Print :: Int->String->IO()
```

```
Print n s = putStr
```

```
fromHeader ++
```

```
fromFormat (palindrome n s))
```

palindrome contains a reference to the `build` and `reverseComplement` as well as `seq`. This function is further discussed below.

```

:: Int->String->[PalindromeInfo]

```

ns =

```

- zip (i n s) (build n s)

```

$$((fst\ x, fst\ x+n-1),$$

```
fst y, fst y+n-1), snd x, snd y) |
```

```

1  x <- ls,

```

← ls,

 $x < y,$ 

(snd x) ==

```
(reverseComplement (snd y)) ]
```

```
Int -> String -> [Int]
```

```
s = take (length s - n + 1) [1..]
```

function build represents the construction of the data structure from the string representation of the original sequence. This data structure is a list of strings, where each element represents a possible nucleotide subsequence. The data structure contains all such prospective subsequences at successive indices in the original sequence. The use of the length function in the guard



structure and subsequently modify that data structure in iterative steps leading to a final post modification process to output the data structure. We have detailed an application of the proposed model in a case study of a palindrome identification algorithm for bioinformatics. The application described here illustrates how the proposed methodology can be easily implemented for developers or coders with little in-depth knowledge of functional programming capabilities; thereby, enabling our simple approach for a wide range of users.

The approach and results in this paper, although both encouraging and preliminary, should be developed further. In particular, user experiments to establish quantitatively the benefit for rapid prototyping should be conducted. Also, we would like to more qualitatively determine the range of algorithms that can be realized by our proposed model. In line with our earlier work both in model identification of program structures and in the visualization of the same, we are motivated to re-consider our work in this paper for application in software visualization.

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