An Introduction to *IRanges*

Patrick Aboyoun, Michael Lawrence, Hervé Pagès

December 11, 2015

1 Introduction

The *IRanges* package is designed to represent sequences, ranges representing indices along those sequences, and data related to those ranges. In this vignette, we will rely on simple, illustrative example datasets, rather than large, real-world data, so that each data structure and algorithm can be explained in an intuitive, graphical manner. We expect that packages that apply *IRanges* to a particular problem domain will provide vignettes with relevant, realistic examples.

The IRanges package is available at bioconductor.org and can be downloaded via biocLite:

```
> source("http://bioconductor.org/biocLite.R")
> biocLite("IRanges")
> library(IRanges)
```

2 Vector objects

In the context of the *IRanges* package, a sequence is an ordered finite collection of elements. The *IRanges* packages represents two types of objects as sequences: (1) atomic sequences and (2) lists (or non-atomic sequences). The following subsections describe each in turn. All *IRanges*-derived sequences inherit from the *Vector* virtual class.

2.1 Atomic Vectors

In R, atomic sequences are typically stored in atomic vectors. The *IRanges* package includes an additional atomic sequence object type, *Rle*, which compresses an atomic sequence through run-length encoding.

We begin our discussion of atomic sequences using two Rle vectors.

All atomic sequences in R have three main properties: (1) a notion of length or number of elements, (2) the ability to extract elements to create new atomic sequences, and (3) the ability to be combined with one or more atomic sequences to form larger atomic sequences. The main functions for these three operations are length, [, and c.

```
> length(xVector)
[1] 10000000
```

> xVector[1]

```
integer-Rle of length 1 with 1 run
  Lengths: 1
  Values : 0
> zVector <- c(xVector, yVector)</pre>
```

While these three methods may seem trivial, they provide a great deal of power and many atomic sequence manipulations can be constructed using them.

2.1.1 Vector Subsetting

As with ordinary R atomic vectors, it is often necessary to subset one sequence from another. When this subsetting does not duplicate or reorder the elements being extracted, the result is called a *subsequence*. In general, the [function can be used to construct a new sequence or extract a subsequence, but its interface is often inconvenient and not amenable to optimization. To compensate for this, the *IRanges* package supports seven additional functions for sequence extraction:

- 1. window Extracts a subsequence over a specified region.
- 2. subset Extracts the subsequence specified by a logical vector.
- 3. head Extracts a consecutive subsequence containing the first n elements.
- 4. tail Extracts a consecutive subsequence containing the last n elements.
- 5. rev Creates a new sequence with the elements in the reverse order.
- 6. rep Creates a new sequence by repeating sequence elements.

The following code illustrates how these functions are used on an ordinary R integer vector:

```
> xSnippet <- xVector[IRanges(4751, 4760)]</pre>
> xSnippet
integer-Rle of length 10 with 9 runs
 Lengths: 1 1 1 1 1 1 1 2
 Values: 4 6 5 4 6 2 6 7 5
> head(xSnippet)
integer-Rle of length 6 with 6 runs
 Lengths: 1 1 1 1 1 1
 Values: 4 6 5 4 6 2
> tail(xSnippet)
integer-Rle of length 6 with 5 runs
 Lengths: 1 1 1 1 2
 Values : 6 2 6 7 5
> rev(xSnippet)
integer-Rle of length 10 with 9 runs
 Lengths: 2 1 1 1 1 1 1 1 1
 Values: 5 7 6 2 6 4 5 6 4
```

```
> rep(xSnippet, 2)
integer-Rle of length 20 with 18 runs
  Lengths: 1 1 1 1 1 1 1 1 2 1 1 1 1 1 1 1 1 2
  Values : 4 6 5 4 6 2 6 7 5 4 6 5 4 6 2 6 7 5
> subset(xSnippet, xSnippet >= 5L)
integer-Rle of length 7 with 5 runs
  Lengths: 1 1 2 1 2
  Values : 6 5 6 7 5
```

2.1.2 Combining Vectors

The *IRanges* package uses two generic functions, c and append, for combining two *Vector* objects. The methods for *Vector* objects follow the definition that these two functions are given the the *base* package.

```
> c(xSnippet, rev(xSnippet))
integer-Rle of length 20 with 17 runs
  Lengths: 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
  Values : 4 6 5 4 6 2 6 7 5 7 6 2 6 4 5 6 4
> append(xSnippet, xSnippet, after = 3)
integer-Rle of length 20 with 18 runs
  Lengths: 1 1 1 1 1 1 1 1 1 1 1 2 1 1 1 1 1 2
  Values : 4 6 5 4 6 5 4 6 2 6 7 5 4 6 2 6 7 5
```

2.1.3 Looping over Vectors and Vector subsets

In R, for looping can be an expensive operation. To compensate for this, *IRanges* uses three generics, endoapply, lapply, and sapply, for looping over sequences and two generics, aggregate and shiftApply, to perform calculations over subsequences. The lapply and sapply functions are familiar to many R users since they are the standard functions for looping over the elements of an R *list* object. The endoapply function performs an endomorphism equivalent to lapply, i.e. returns a *Vector* object of the same class as the input rather than a *list* object. More will be given on these three functions in the Lists subsection.

The aggregate function combines sequence extraction functionality of the window function with looping capabilities of the sapply function. For example, here is some code to compute medians across a moving window of width 3 using the function aggregate:

```
> xSnippet
```

```
integer-Rle of length 10 with 9 runs
  Lengths: 1 1 1 1 1 1 1 2
  Values : 4 6 5 4 6 2 6 7 5

> aggregate(xSnippet, start = 1:8, width = 3, FUN = median)
[1] 5 5 5 4 6 6 6 5
```

The shiftApply function is a looping operation involving two sequences whose elements are lined up via a positional shift operation. For example, the elements of xVector and yVector were simulated from Poisson distributions with the mean of element i from yVector being equivalent to the mean of element i + 250 from xVector. If we did not know the size of the shift, we could estimate it by finding the shift that maximizes the correlation between xVector and yVector.



Figure 1: Correlation between xVector and yVector for various shifts.

```
> cor(xVector, yVector)
[1] 0.5739224
> shifts <- seq(235, 265, by=3)
> corrs <- shiftApply(shifts, yVector, xVector, FUN = cor)
> plot(shifts, corrs)
   The result is shown in Fig. 1.
```

2.1.4 Run Length Encoding

Up until this point we have used R atomic vectors to represent atomic sequences, but there are times when these object become too large to manage in memory. When there are lots of consecutive repeats in the sequence, the data can be compressed and managed in memory through a run-length encoding where a data value is paired with a run length. For example, the sequence $\{1, 1, 1, 2, 3, 3\}$ can be represented as values $= \{1, 2, 3\}$, run lengths $= \{3, 1, 2\}$.

The *Rle* class in *IRanges* is used to represent a run-length encoded (compressed) sequence of *logical*, *integer*, *numeric*, *complex*, *character*, or *raw* values. One way to construct an *Rle* object is through the *Rle* constructor function:

```
> xRle <- Rle(xVector)
> yRle <- Rle(yVector)</pre>
> xRle
integer-Rle of length 10000000 with 1510219 runs
                                1 1599
 Lengths:
            780
                        208
                                                     5
                                                               91
                                                                        927
                    1
                                           1 ...
                                                          1
  Values:
                    1
                          0
                                1
                                           1 ...
                                                     0
                                                          1
                                                                     1
```

> yRle

Values:

```
integer-Rle of length 10000000 with 1511351 runs
Lengths: 1003  1 413  1 896  1 ... 3  1 845  1
```

0

1

When there are lots of consecutive repeats, the memory savings through an RLE can be quite dramatic. For example, the xRle object occupies less than one quarter of the space of the original xVector object, while storing the same information:

0

1

1 ...

419

1

> as.vector(object.size(xRle) / object.size(xVector))

[1] 1

> identical(as.vector(xRle), xVector)

[1] FALSE

The functions runValue and runLength extract the run values and run lengths from an *Rle* object respectively:

> head(runValue(xRle))

[1] 0 1 0 1 0 1

> head(runLength(xRle))

[1] 780 1 208 1 1599 1

The *Rle* class supports many of the basic methods associated with R atomic vectors including the Ops, Math, Math2, Summary, and Complex group generics. Here is a example of manipulating *Rle* objects using methods from the Ops group:

> xRle > 0

```
logical-Rle of length 10000000 with 197127 runs
```

Lengths: 780 1 208 1 1599 ... 1 91 1 927
Values: FALSE TRUE FALSE TRUE FALSE ... TRUE FALSE TRUE FALSE

> xRle + yRle

integer-Rle of length 10000000 with 1957707 runs

1 208 1 413 ... 5 1 507 1 419 Lengths: 780 13 1 91 Values: 1 0 1 0 1 0 ... 0 1 0 1 1

> xRle > 0 | yRle > 0

logical-Rle of length 10000000 with 210711 runs

Lengths: 780 1 208 1 13 ... 1 507 1 419 Values: FALSE TRUE FALSE TRUE FALSE ... TRUE FALSE TRUE FALSE

Here are some from the Summary group:

> range(xRle)

[1] 0 26

[1] 0.8519138 0.8517324 0.8517725

For more information on the methods supported by the Rle class, consult the Rle man page.

> shiftApply(249:251, yRle, xRle, FUN = function(x, y) var(x, y) / (sd(x) * sd(y)))

2.2 Lists

In many data analysis situation there is a desire to organize and manipulate multiple objects simultaneously. Typically this is done in R through the usage of a list. While a list serves as a generic container, it does not confer any information about the specific class of its elements, provides no infrastructure to ensure type safety, and the S3 and S4 method dispatch mechanisms do not support method selection for lists with homogeneous object types. The *List* virtual class defined in the *IRanges* package addresses these issues. *List* is a direct extension of *Vector*.

2.2.1 Lists of Atomic Vectors

The first type of lists we consider are those containing atomic sequences such as *integer* vectors or *Rle* objects. We may wish to define a method that retrieves the length of each atomic sequence element, without special type checking. To enable this, we define collection classes such as *IntegerList* and *RleList*, which inherit from the *List* virtual class, for representing lists of *integer* vectors and *Rle* objects respectively.

```
> getClassDef("RleList")
Virtual Class "RleList" [package "IRanges"]
Slots:
Name: elementType elementMetadata metadata
Class: character DataTableORNULL list

Extends:
Class "AtomicList", directly
Class "List", by class "AtomicList", distance 2
Class "Vector", by class "AtomicList", distance 3
Class "Annotated", by class "AtomicList", distance 4
Known Subclasses: "RleViews", "CompressedRleList", "SimpleRleList"
```

As the class definition above shows, the *RleList* class is virtual with subclasses *SimpleRleList* and *CompressedRleList*. A *SimpleRleList* class uses a regular R list to store the underlying elements and the *CompressedRleList* class stores the elements in an unlisted form and keeps track of where the element breaks are. The former "simple list" class is useful when the Rle elements are long and the latter "compressed list" class is useful when the list is long and/or sparse (i.e. a number of the list elements have length 0).

In fact, all of the atomic vector types (raw, logical, integer, numeric, complex, and character) have similar list classes that derive from the *List* virtual class. For example, there is an *IntegerList* virtual class with subclasses *SimpleIntegerList* and *CompressedIntegerList*.

Each of the list classes for atomic sequences, be they stored as vectors or Rle objects, have a constructor function with a name of the appropriate list virtual class, such as IntegerList, and an optional argument compress that takes an argument to specify whether or not to create the simple list object type or the compressed list object type. The default is to create the compressed list object type.

```
> args(IntegerList)
function (..., compress = TRUE)
> cIntList1 <- IntegerList(x = xVector, y = yVector)</pre>
> cIntList1
IntegerList of length 2
> sIntList2 <- IntegerList(x = xVector, y = yVector, compress = FALSE)
> sIntList2
IntegerList of length 2
> ## sparse integer list
> xExploded <- lapply(xVector[1:5000], function(x) seq_len(x))</pre>
> cIntList2 <- IntegerList(xExploded)</pre>
> sIntList2 <- IntegerList(xExploded, compress = FALSE)
> object.size(cIntList2)
33080 bytes
> object.size(sIntList2)
253952 bytes
```

The length function returns the number of elements in a *Vector*-derived object and, for a *List*-derived object like "simple list" or "compressed list", the elementLengths function returns an integer vector containing the lengths of each of the elements:

```
> length(cIntList2)
[1] 5000
> Rle(elementLengths(cIntList2))
```

```
integer-Rle of length 5000 with 427 runs
                              1 1599
 Lengths: 780
                   1
                      208
                                         1 ...
                                                  1
                                                             1
                                                                  1
                                                                       1
 Values :
                   1
                         0
                              1
                                   0
                                         1 ...
                                                 10
```

Just as with ordinary R list objects, List-derived object support the [[for element extraction, c for combining, and lapply/sapply for looping. When looping over sparse lists, the "compressed list" classes can be much faster during computations since only the non-empty elements are looped over during the lapply/sapply computation and all the empty elements are assigned the appropriate value based on their status.

```
> system.time(sapply(xExploded, mean))
         system elapsed
   user
  0.062
         0.000
                  0.061
> system.time(sapply(sIntList2, mean))
   user
         system elapsed
  0.056
          0.000
                  0.056
> system.time(sapply(cIntList2, mean))
   user system elapsed
  0.057
         0.000
                  0.056
> identical(sapply(xExploded, mean), sapply(sIntList2, mean))
[1] TRUE
> identical(sapply(xExploded, mean), sapply(cIntList2, mean))
[1] TRUE
   Unlist ordinary R list objects, AtomicList objects support the Ops (e.g. +, ==, &), Math (e.g. log, sqrt),
Math2 (e.g. round, signif), Summary (e.g. min, max, sum), and Complex (e.g. Re, Im) group generics.
> xRleList <- RleList(xRle, 2L * rev(xRle))
> yRleList <- RleList(yRle, 2L * rev(yRle))
> xRleList > 0
RleList of length 2
[[1]]
logical-Rle of length 10000000 with 197127 runs
             780
                     1
                         208
                                  1 1599 ...
 Values: FALSE TRUE FALSE TRUE FALSE ... TRUE FALSE TRUE FALSE
[[2]]
logical-Rle of length 10000000 with 197127 runs
 Lengths:
             927
                     1
                           91
                                  1
                                        5 ...
                                                   1
                                                       208
                                                               1
                                                                    780
 Values : FALSE TRUE FALSE TRUE FALSE ... TRUE FALSE TRUE FALSE
> xRleList + yRleList
```

```
RleList of length 2
[[1]]
integer-Rle of length 10000000 with 1957707 runs
                                 1 413 ... 5 1 91 1 507
 Lengths: 780
                1 208
                         1 13
                                                                 1 419
 Values: 0
                1
                     0
                         1
                             0
                                 1
                                     0 ...
                                             0 1 0 1 0
                                                                 1 0
[[2]]
integer-Rle of length 10000000 with 1957707 runs
 Lengths: 419
                 1 507
                                     5 ... 413
                                                         1 208
                                                                 1 780
                         1 91
                                 1
                                                 1 13
  Values :
                2 0
                         2
                             0
                                 2
                                     0 ... 0
                                                 2 0
                                                         2 0
                                                                 2 0
> sum(xRleList > 0 | yRleList > 0)
[1] 2105185 2105185
   Since these atomic lists inherit from List, they can also use the looping function endoapply to perform
endomorphisms.
> safe.max <- function(x) { if(length(x)) max(x) else integer(0) }</pre>
> endoapply(sIntList2, safe.max)
IntegerList of length 5000
[[1]] integer(0)
[[2]] integer(0)
[[3]] integer(0)
[[4]] integer(0)
[[5]] integer(0)
[[6]] integer(0)
[[7]] integer(0)
[[8]] integer(0)
[[9]] integer(0)
[[10]] integer(0)
<4990 more elements>
> endoapply(cIntList2, safe.max)
IntegerList of length 5000
[[1]] integer(0)
[[2]] integer(0)
[[3]] integer(0)
[[4]] integer(0)
[[5]] integer(0)
[[6]] integer(0)
[[7]] integer(0)
[[8]] integer(0)
[[9]] integer(0)
[[10]] integer(0)
<4990 more elements>
> endoapply(sIntList2, safe.max)[[1]]
integer(0)
```

3 Data Tables

 ${\bf To\ Do:}\ DataTable,\ DataFrame,\ DataFrameList,\ SplitDataFrameList$

4 Vector Annotations

Often when one has a collection of objects, there is a need to attach metadata that describes the collection in some way. Two kinds of metadata can be attached to a *Vector* object:

- Metadata about the object as a whole: this metadata is accessed via the metadata accessor and is represented as an ordinary list;
- 2. Metadata about the individual elements of the object: this metadata is accessed via the mcols accessor (mcols stands for metadata columns) and is represented as a DataTable object (i.e. as an instance of a concrete subclass of DataTable, e.g. a DataFrame object). This DataTable object can be thought of as the result of binding together one or several vector-like objects (the metadata columns) of the same length as the Vector object. Each row of the DataTable object annotates the corresponding element of the Vector object.

5 Vector Ranges

When analyzing sequences, we are often interested in particular consecutive subsequences. For example, the a vector could be considered a sequence of lower-case letters, in alphabetical order. We would call the first five letters (a to e) a consecutive subsequence, while the subsequence containing only the vowels would not be consecutive. It is not uncommon for an analysis task to focus only on the geometry of the regions, while ignoring the underlying sequence values. A list of indices would be a simple way to select a subsequence. However, a sparser representation for consecutive subsequences would be a range, a pairing of a start position and a width, as used when extracting sequences with window.

When analyzing subsequences in *IRanges*, each range is treated as an observation. The virtual *Ranges* class represents lists of ranges, or, equivalently and as a derivative *IntegerList*, sequences of consecutive integers. The most commonly used implementation of *Ranges* is *IRanges*, which stores the starts and widths as ordinary integer vectors. To construct an *IRanges* instance, we call the *IRanges* constructor. Ranges are normally specified by passing two out of the three parameters: start, end and width (see help(IRanges) for more information).

```
> ir1 <- IRanges(start = 1:10, width = 10:1)
> ir2 <- IRanges(start = 1:10, end = 11)
> ir3 <- IRanges(end = 11, width = 10:1)
> identical(ir1, ir2) & identical(ir2, ir3)

[1] FALSE
> ir <- IRanges(c(1, 8, 14, 15, 19, 34, 40), # width = c(12, 6, 6, 15, 6, 2, 7))</pre>
```

All of the above calls construct an *IRanges* instance with the same ranges, using different combinations of the start, end and width parameters.

Accessing the starts, widths and ends is supported by every Ranges implementation.

```
> start(ir)
[1] 1 8 14 15 19 34 40
```

```
> end(ir)
[1] 12 13 19 29 24 35 46
> width(ir)
[1] 12 6 6 15 6 2 7
```

For IRanges and some other Ranges derivatives, subsetting is also supported, by numeric and logical indices.

> ir[1:4]

```
IRanges of length 4
    start end width
[1]
        1 12
[2]
        8 13
                  6
[3]
       14 19
                  6
[4]
       15
          29
                 15
> ir[start(ir) <= 15]
IRanges of length 4
    start end width
[1]
        1 12
                 12
[2]
        8 13
                  6
[3]
       14 19
                  6
[4]
       15 29
```

One may think of each range as a sequence of integer ranges, and Ranges is, in fact, derived from IntegerList.

> ir[[1]]

[1] 1 2 3 4 5 6 7 8 9 10 11 12

In order to illustrate range operations, we'll create a function to plot ranges.

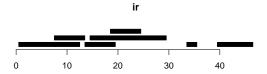


Figure 2: Plot of original ranges.

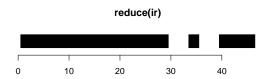


Figure 3: Plot of reduced ranges.

5.1 Normality

Sometimes, it is necessary to formally represent a subsequence, where no elements are repeated and order is preserved. Also, it is occasionally useful to think of a *Ranges* object as a set, where no elements are repeated and order does not matter. While every *Ranges* object, as a *Vector* derivative, has an implicit ordering, one can enforce the same ordering for all such objects, so that ordering becomes inconsequential within that context.

The *NormalIRanges* class formally represents either a subsequence encoding or a set of integers. By definition a Ranges object is said to be *normal* when its ranges are: (a) not empty (i.e. they have a non-null width); (b) not overlapping; (c) ordered from left to right; (d) not even adjacent (i.e. there must be a non empty gap between 2 consecutive ranges).

There are three main advantages of using a *normal Ranges* object: (1) it guarantees a subsequence encoding or set of integers, (2) it is compact in terms of the number of ranges, and (3) it uniquely identifies its information, which simplifies comparisons.

The reduce function reduces any Ranges object to a NormalIRanges by merging redundant ranges.

> reduce(ir)

IRanges of length 3
start end width
[1] 1 29 29
[2] 34 35 2
[3] 40 46 7

> plotRanges(reduce(ir))

5.2 Lists of Ranges objects

It is common to manipulate collections of *Ranges* objects during an analysis. Thus, the *IRanges* package defines some specific classes for working with multiple *Ranges* objects.

The RangesList class asserts that each element is a Ranges object and provides convenience methods, such as start, end and width accessors that return IntegerList objects, aligning with the RangesList object. To explicitly construct a RangesList, use the RangesList function.

```
> rl <- RangesList(ir, rev(ir))
> start(rl)
IntegerList of length 2
[[1]] 1 8 14 15 19 34 40
[[2]] 40 34 19 15 14 8 1
```

5.3 Vector Extraction

As the elements of a *Ranges* object encode consecutive subsequences, they may be used directly in sequence extraction. Note that when a *normal Ranges* is given as the index, the result is a subsequence, as no elements are repeated or reordered. If the sequence is a *Vector* subclass (i.e. not an ordinary *vector*), the canonical [function accepts a *Ranges* instance.

5.4 Finding Overlapping Ranges

The function findOverlaps detects overlaps between two Ranges objects.

```
> ol <- findOverlaps(ir, reduce(ir))</pre>
> as.matrix(ol)
     queryHits subjectHits
[1,]
              1
                             1
[2,]
               2
                             1
[3,]
              3
                             1
              4
[4,]
                            1
[5,]
              5
                             1
[6,]
               6
                            2
[7,]
              7
                            3
```

5.5 Counting Overlapping Ranges

The function coverage counts the number of ranges over each position.

```
> cov <- coverage(ir)
> plotRanges(ir)
> cov <- as.vector(cov)
> mat <- cbind(seq_along(cov)-0.5, cov)
> d <- diff(cov) != 0
> mat <- rbind(cbind(mat[d,1]+1, mat[d,2]), mat)
> mat <- mat[order(mat[,1]),]
> lines(mat, col="red", lwd=4)
> axis(2)
```

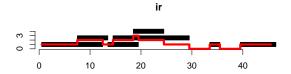


Figure 4: Plot of ranges with accumulated coverage.

5.6 Finding Neighboring Ranges

The nearest function finds the nearest neighbor ranges (overlapping is zero distance). The precede and follow functions find the non-overlapping nearest neighbors on a specific side.

5.7 Transforming Ranges

Utilities are available for transforming a *Ranges* object in a variety of ways. Some transformations, like reduce introduced above, can be dramatic, while others are simple per-range adjustments of the starts, ends or widths.

5.7.1 Adjusting starts, ends and widths

Perhaps the simplest transformation is to adjust the start values by a scalar offset, as performed by the shift function. Below, we shift all ranges forward 10 positions.

> shift(ir, 10)

There are several other ways to transform ranges. These include narrow, resize, flank, reflect, restrict, and threebands. For example narrow supports the adjustment of start, end and width values, which should be relative to each range. These adjustments are vectorized over the ranges. As its name suggests, the ranges can only be narrowed.

> narrow(ir, start=1:5, width=2)

IRar	iges o	ρf	ler	igth	7
	start	; (end	widt	:h
[1]	-	L	2		2
[2]	Ş)	10		2
[3]	16	3	17		2
[4]	18	3	19		2
[5]	23	3	24		2

```
[6] 34 35 2
[7] 41 42 2
```

The restrict function ensures every range falls within a set of bounds. Ranges are contracted as necessary, and the ranges that fall completely outside of but not adjacent to the bounds are dropped, by default.

```
> restrict(ir, start=2, end=3)
```

```
IRanges of length 1 start end width [1] 2 3 2
```

The threebands function extends narrow so that the remaining left and right regions adjacent to the narrowed region are also returned in separate *Ranges* objects.

```
> threebands(ir, start=1:5, width=2)
```

\$left

IRar	nges of	fler	ngth 7
	start	end	width
[1]	1	0	0
[2]	8	8	1
[3]	14	15	2
[4]	15	17	3
[5]	19	22	4
[6]	34	33	0
[7]	40	40	1

\$middle

IRar	nges	01	Ler	ngth	7
	staı	ct	end	widt	h
[1]		1	2		2
[2]		9	10		2
[3]	-	16	17		2
[4]	-	18	19		2
[5]	2	23	24		2
[6]	3	34	35		2
[7]	4	11	42		2

\$right

```
IRanges of length 7
    start end width
[1]
         3
            12
                   10
[2]
        11
            13
                     3
                    2
[3]
        18
            19
[4]
        20
            29
                   10
        25
            24
                    0
[5]
[6]
        36
            35
                    0
                     4
[7]
        43
            46
```

The arithmetic operators +, - and * change both the start and the end/width by symmetrically expanding or contracting each range. Adding or subtracting a numeric (integer) vector to a Ranges causes each range to be expanded or contracted on each side by the corresponding value in the numeric vector.

> ir + seq_len(length(ir))

```
IRanges of length 7
    start end width
[1]
          13
        0
[2]
           15
                  10
        6
[3]
       11
           22
                  12
[4]
       11 33
                  23
[5]
       14
           29
                  16
[6]
       28 41
                  14
[7]
       33 53
                  21
```

The * operator symmetrically magnifies a *Ranges* object by a factor, where positive contracts (zooms in) and negative expands (zooms out).

> ir * -2 # double the width

IRar	nges o	f ler	ngth 7
	start	end	width
[1]	-5	18	24
[2]	5	16	12
[3]	11	22	12
[4]	7	36	30
[5]	16	27	12
[6]	33	36	4
[7]	36	49	14

WARNING: The semantic of these arithmetic operators might be revisited at some point. Please restrict their use to the context of interactive visualization (where they arguably provide some convenience) but avoid to use them programmatically.

5.7.2 Making ranges disjoint

A more complex type of operation is making a set of ranges disjoint, *i.e.* non-overlapping. For example, threebands returns a disjoint set of three ranges for each input range.

The disjoin function makes a *Ranges* object disjoint by fragmenting it into the widest ranges where the set of overlapping ranges is the same.

> disjoin(ir)

IRanges	of	leng	gth	10
sta	art	end	wid	lth
[1]	1	7		7
[2]	8	12		5
[3]	13	13		1
[4]	14	14		1
[5]	15	18		4
[6]	19	19		1
[7]	20	24		5
[8]	25	29		5
[9]	34	35		2
[10]	40	46		7

> plotRanges(disjoin(ir))

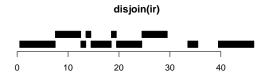


Figure 5: Plot of disjoined ranges.

A variant of disjoin is disjointBins, which divides the ranges into bins, such that the ranges in each bin are disjoint. The return value is an integer vector of the bins.

> disjointBins(ir)

[1] 1 2 1 2 3 1 1

5.7.3 Other transformations

Other transformations include reflect and flank. The former "flips" each range within a set of common reference bounds.

> reflect(ir, IRanges(start(ir), width=width(ir)*2))

IRanges of length 7 start end width [1] 13 24 12 [2] 14 19 6 [3] 20 25 6 [4] 30 44 15 [5] 25 30 6 36 2 [6] 37 7 [7] 47 53

The flank returns ranges of a specified width that flank, to the left (default) or right, each input range. One use case of this is forming promoter regions for a set of genes.

> flank(ir, width = seq_len(length(ir)))

IRar	ıges	of	ler	ngth	7
	staı	rt	end	widt	h
[1]		0	0		1
[2]		6	7		2
[3]	-	11	13		3
[4]	-	11	14		4
[5]	-	14	18		5
[6]	2	28	33		6
[7]	3	33	39		7

gaps(ir, start = 1, end = 50) 10 20 30 40 50

Figure 6: Plot of gaps from ranges.

5.8 Set Operations

Sometimes, it is useful to consider a *Ranges* object as a set of integers, although there is always an implicit ordering. This is formalized by *NormalIRanges*, above, and we now present versions of the traditional mathematical set operations *complement*, *union*, *intersect*, and *difference* for *Ranges* objects. There are two variants for each operation. The first treats each *Ranges* object as a set and returns a *normal* value, while the other has a "parallel" semantic like pmin/pmax and performs the operation for each range pairing separately.

The *complement* operation is implemented by the gaps and pgap functions. By default, gaps will return the ranges that fall between the ranges in the (normalized) input. It is possible to specify a set of bounds, so that flanking ranges are included.

```
> gaps(ir, start=1, end=50)
```

```
IRanges of length 3
start end width
[1] 30 33 4
[2] 36 39 4
[3] 47 50 4
```

> plotRanges(gaps(ir, start=1, end=50), c(1,50))

pgap considers each parallel pairing between two *Ranges* objects and finds the range, if any, between them. Note that the function name is singular, suggesting that only one range is returned per range in the input.

The remaining operations, *union*, *intersect* and *difference* are implemented by the [p]union, [p]intersect and [p]setdiff functions, respectively. These are relatively self-explanatory.

6 Vector Views

The *IRanges* package provides the virtual *Views* class, which stores a sequence together with an *IRanges* object defining ranges on the sequence. Each range is said to represent a *view* onto the sequence.

Here, we will demonstrate the *Rle Views* class, where the sequence is of class *Rle*. Other *Views* implementations exist, such as *XString Views* in the *Biostrings* package.

6.1 Creating Views

There are two basic constructors for creating views: the Views function based on indicators and the slice based on numeric boundaries.

```
> xViews <- Views(xRle, xRle >= 1)
> xViews <- slice(xRle, 1)
> xViewsList <- slice(xRleList, 1)</pre>
```

6.2 Aggregating Views

While sapply can be used to loop over each window, the native functions viewMaxs, viewMins, viewSums, and viewMeans provide fast looping to calculate their respective statistical summaries.

```
> head(viewSums(xViews))
[1] 1 1 1 1 1 2
> viewSums(xViewsList)
IntegerList of length 2
[[1]] 1 1 1 1 1 2 1 1 2 3 1 6 1 3 4 ... 12 6 37 10 8 11 6 4 5 1 1 5 1 1
[[2]] 2 2 10 2 2 10 8 12 22 16 20 74 12 ... 2 12 2 6 4 2 2 4 2 2 2 2 2 2
> head(viewMaxs(xViews))
[1] 1 1 1 1 1 2
> viewMaxs(xViewsList)
IntegerList of length 2
[[1]] 1 1 1 1 1 2 1 1 1 2 1 2 1 2 1 2 3 1 ... 3 5 2 5 6 2 8 3 2 2 1 1 2 1 1
[[2]] 2 2 4 2 2 4 4 6 16 4 12 10 4 10 6 ... 4 2 4 2 4 2 2 2 4 2 2 2 2
```

7 Data on Ranges

IMPORTANT NOTE: Starting with BioC 2.12, the use of RangedData and RangedDataList objects is discouraged in favor of GRanges and GRangesList objects (those classes are defined in the GenomicRanges package).

When analyzing ranges, there are often additional variables of interest, besides the geometry (starts, ends and widths). To formally represent a dataset where the ranges are the observations, IRanges defines the RangedData class.

To create a *RangedData* instance, one needs to provide a *Ranges* object and, optionally, any number of variables on those ranges. The variable objects need not be vectors, but they must satisfy the contract of *DataFrame*.

```
> values <- rnorm(length(ir))</pre>
> rd <- RangedData(ir, name = letters[seq_len(length(ir))], values)</pre>
> rd
RangedData with 7 rows and 2 value columns across 1 space
              ranges |
                              name
                                        values
  <factor> <IRanges> | <character>
                                    <numeric>
         1 [1, 12] |
1
                                 a 0.4359675
2
         1 [8, 13] |
                                 b 0.7279741
         1 [14, 19] |
3
                                  c 0.2617505
         1 [15, 29] |
4
                                 d 0.8157306
                                 e 1.2115267
5
         1 [19, 24] |
6
                                 f 1.3045325
         1
           [34, 35] |
         1 [40, 46] |
                                 g -1.2268231
```

One might notice the term "sequence" in the above output. This refers to an important feature of RangedData: the ability to segregate ranges by their sequence (or space). For example, when analyzing genomic data, one is often working with ranges on different chromosomes. In many cases, such as when calculating overlap, one needs to separately treat ranges from different spaces, and RangedData aims to facilitate this mode of operation. The segregation may be performed at construction time.

```
> rd <- RangedData(ir, name = letters[seq_len(length(ir))], values,</pre>
                   space = rep(c("chr1", "chr2"), c(3, length(ir) - 3)))
> rd
RangedData with 7 rows and 2 value columns across 2 spaces
     space
              ranges |
                               name
                                        values
  <factor> <IRanges> | <character>
                                     <numeric>
            [ 1, 12]
                                     0.4359675
1
      chr1
                     2
            [8, 13] |
      chr1
                                  b 0.7279741
3
      chr1
            [14, 19] |
                                  c 0.2617505
      chr2
4
            [15, 29] |
                                  d 0.8157306
5
      chr2
            [19, 24]
                                  е
                                    1.2115267
6
      chr2
            [34, 35]
                                  f 1.3045325
7
      chr2
            [40, 46] |
                                  g -1.2268231
```

With the knowledge that the data is split into spaces, it should not be surprising that the ranges accessor returns a RangesList and values returns a SplitDataFrameList.

> ranges(rd)

IRangesList of length 2

```
$chr1
IRanges of length 3
    start end width
[1]
        1
           12
                  12
[2]
        8
           13
                   6
[3]
       14
           19
                   6
$chr2
IRanges of length 4
    start end width
[1]
       15
            29
                  15
[2]
       19
            24
                   6
[3]
       34
           35
                   2
                   7
       40
[4]
           46
> values(rd)
SplitDataFrameList of length 2
$chr1
DataFrame with 3 rows and 2 columns
                  values
         name
  <character> <numeric>
             a 0.4359675
1
2
             b 0.7279741
3
             c 0.2617505
```

\$chr2

DataFrame with 4 rows and 2 columns
name values
<character> <numeric>
1 d 0.8157306
2 e 1.2115267
3 f 1.3045325

g -1.2268231

To obtain a *RangedData* for a specific set of spaces, one should use the [function, which accepts logical, numeric and character indices.

> rd["chr1"]

[1] TRUE

The names and length functions return the names and number of spaces, respectively.

- > names(rd)
- [1] "chr1" "chr2"
- > length(rd)
- [1] 2

The lapply function operates over the spaces. The object passed to the user function is a subset RangedData.

```
> lapply(rd, names)
```

\$chr1

[1] "chr1"

\$chr2

[1] "chr2"

The above would suggest that RangedData is a sequence of spaces. However, RangedData also inherits from DataTable, so it in some ways behaves like a sequence of columns. For example, one can extract a column via $\$ or $\[\]$

> rd[[2]]

```
[1] 0.4359675 0.7279741 0.2617505 0.8157306 1.2115267 1.3045325 [7] -1.2268231
```

> rd\$values

```
[1] 0.4359675 0.7279741 0.2617505 0.8157306 1.2115267 1.3045325 [7] -1.2268231
```

Note that the extracted columns are "unlisted" over the spaces, which is usually much more convenient than obtaining them as lists. It is important to note that the elements have been sorted by the space factor and thus may not have the same order as the objects passed to the constructor.

The two dimensional matrix-style subsetting is also supported. The rows are indexed globally, independent of space.

> rd[1:3, "name"]

 ${\tt RangedData\ with\ 3\ rows\ and\ 1\ value\ column\ across\ 2\ spaces}$

	space	ranges		name
	<factor></factor>	Ranges		<character></character>
1	chr1	[1, 12]		a
2	chr1	[8,13]	1	Ъ
3	chr1	[14, 19]	1	С

8 IRanges in Biological Sequence Analysis

The *IRanges* packages was primarily designed with biological sequence analysis in mind and Table 1 shows how some biological sequence analysis concepts are represented in the *IRanges* class system.

Biological Entity	Vector Subclass
Genome browser track(s)	GRanges/GRangesList
Coverage across chromosomes/contigs	RleList
Mapped ranges to genome	CompressedIRangesList
Data (sans ranges) across chroms/contigs	SplitDataFrameList

Table 1: Vector subclasses for Biological Sequence Analysis

> toLatex(sessionInfo())

- R version 3.2.2 (2015-08-14), x86_64-pc-linux-gnu
- Locale: LC_CTYPE=en_US.UTF-8, LC_NUMERIC=C, LC_TIME=en_US.UTF-8, LC_COLLATE=C, LC_MONETARY=en_US.UTF-8, LC_MESSAGES=en_US.UTF-8, LC_PAPER=en_US.UTF-8, LC_NAME=C, LC_ADDRESS=C, LC_TELEPHONE=C, LC_MEASUREMENT=en_US.UTF-8, LC_IDENTIFICATION=C
- Base packages: base, datasets, grDevices, graphics, methods, parallel, stats, stats4, utils
- Other packages: BiocGenerics 0.16.1, IRanges 2.4.6, S4Vectors 0.8.5
- Loaded via a namespace (and not attached): tools 3.2.2

Table 2: The output of sessionInfo on the build system after running this vignette.

9 Session Information