Overview of the *intervals* package

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1 Introduction

The *intervals* packages defines two S4 classes which represent collections of intervals over either the integers (\mathbb{Z}) or the real number line (\mathbb{R}) . An instance of either class consists of a two-column matrix of endpoints, plus additional slots describing endpoint closure and whether the intervals are to be thought of as being over \mathbb{Z} or \mathbb{R} .

```
> library( intervals )
> x <- Intervals( matrix( 1:6, ncol = 2 ) )
> x

Object of class Intervals
3 intervals over R:
[1, 4]
[2, 5]
[3, 6]
> x[2,2] <- NA
> x[3,1] <- 6
> x

Object of class Intervals
3 intervals over R:
[1, 4]
[2, NA]
[6, 6]
```

Objects of class Intervals represent collections of intervals with common endpoint closure, e.g., all left-closed, right-open. More control over endpoints is permitted with the Intervals_full class. (Both classes are derived from Intervals_virtual, which is not intended for use by package users.)

```
> y <- as( x, "Intervals_full" )
> closed(y)[2:3,1] <- FALSE
> y

Object of class Intervals_full
3 intervals over R:
[1, 4]
(2, NA]
(6, 6]
```

The size method gives measure — counting measure over \mathbb{Z} or Lebesgue measure over \mathbb{R} — for each interval represented in an object. The empty method identifies intervals that are in fact empty sets, which over \mathbb{R} is not the same thing as having size 0. (Valid objects must have each right endpoint no less than the corresponding left endpoint. When one or both endpoints are open, however, valid intervals may be empty.)

```
> size(x)
[1] 3 NA 0
> empty(x)
[1] FALSE NA FALSE
> empty(y)
[1] FALSE NA TRUE
```

2 Interpretation of objects

An Intervals or Intervals_full object can be thought of in two different modes, each of which is useful in certain contexts:

- 1. As a (non-unique) representation of a subset of \mathbb{Z} or \mathbb{R} .
- 2. As a collection of (possibly overlapping) intervals, each of which has a meaningful identity.

2.1 As a subset of \mathbb{Z} or \mathbb{R}

The *intervals* package provides a number of basic tools for working in the first mode, where an object represents a subset of \mathbb{Z} or \mathbb{R} but the rows of the endpoint matrix do not have any external identity. Basic tools include reduce, which returns a sorted minimal representation equivalent to the original (dropping any intervals with NA endpoints), as well as interval_union, interval_complement, and interval_intersection.

```
> reduce( y )
Object of class Intervals_full
1 interval over R:
[1, 4]
> interval_intersection( x, x + 2 )
Object of class Intervals
2 intervals over R:
[3, 4]
[6, 6]
> interval_complement( x )
```

```
Object of class Intervals
3 intervals over R:
(-Inf, 1)
(4, 6)
(6, Inf)
```

Note that combining x and its complement in order to compute a union requires mixing endpoint closure types; coercion to Intervals_full is automatic.

```
> interval_union( x, interval_complement( x ) )
Object of class Intervals_full
1 interval over R:
(-Inf, Inf)
```

The distance_to_nearest function treats its to argument in the first mode, as just a subset of \mathbb{Z} or \mathbb{R} ; it treats its from argument, however, in the second mode, returning one distance for every row of the from object. In the example below, we also look at performance for large data sets (less than one second on a 2 GHz Intel Core 2 Duo Macintosh, although the time shown below will likely differ). A histogram of d is given in Figure 1.

```
> B <- 100000
> left <- runif( B, 0, 1e8 )
> right <- left + rexp( B, rate = 1/10 )
> v <- Intervals( cbind( left, right ) )
> head( v )
Object of class Intervals
6 intervals over R:
[23416442.0748129, 23416442.3044362]
[35034340.0845304, 35034350.4605233]
[76282386.877574, 76282390.2322924]
[89128359.0346575, 89128359.2160263]
[37480682.9728186, 37480688.2968107]
[65373180.3176925, 65373202.0375136]
> mean( size( v ) )
[1] 10.01945
> dim( reduce( v ) )
[1] 99005
> system.time( d <- distance_to_nearest( sample( 1e8, B ), v ) )
   user
         system elapsed
  0.787
          0.896
                  5.428
```

Distance to nearest interval

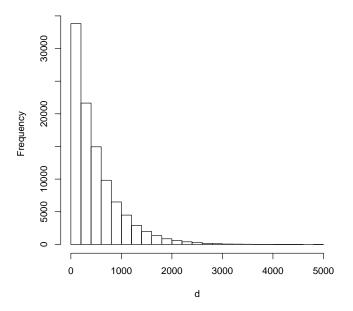


Figure 1: Histogram of distances from a random set of points to the nearest point in v. There is also a distance_to_nearest method for comparing two sets of intervals.

2.2 As a set of meaningful, possibly overlapping intervals

In some applications, each row of an object's endpoint matrix has a meaningful identity, and particular points from \mathbb{Z} or \mathbb{R} may be found in more than one row. To support this mode, objects may be given row names, which are propagated through calculations when appropriate. The **combine** methods simply stack objects (like **rbind**), preserving row names and retaining redundancy, if any.

The interval_overlap method works in this mode. In the next example we use it to identify rows of v which are at least partially redundant, i.e., which intersect at least one other row of v. All rows overlap themselves, so we look for rows that overlap at least two rows:

```
> rownames(v) <- sprintf( "%06i", 1:nrow(v) )
> io <- interval_overlap( v, v )
> head( io, n = 3 )

$`000001`
[1] 1

$`000002`
[1] 2

$`000003`
[1] 3
```

```
> n <- sapply( io, length )
> sum( n > 1 )
[1] 1981
> k \leftarrow which.max(n)
> io[ k ]
$`011631`
[1] 11631 61305 79184
> v[ k, ]
Object of class Intervals
1 interval over R:
011631 [27565230.8482677, 27565265.7762516]
> v[ io[[ k ]], ]
Object of class Intervals
3 intervals over R:
011631 [27565230.8482677, 27565265.7762516]
061305 [27565238.927491, 27565247.9397463]
079184 [27565254.2943135, 27565261.0125249]
```

The which_nearest method (v. 0.9.0, not yet implemented) also respects row identity, returning a list with one vector of indices into the to object for each row of the from object.

3 Floating point and intervals over \mathbb{R}

When type == "R", interval endpoints are not truly in \mathbb{R} , but rather, in the subset which can be represented by floating point arithmetic. (For the moment, this is also true when type == "Z". See Section 4.) This limits the endpoint values which can be represented; more importantly, if computations are performed on interval endpoints, it means that floating point error can affect whether or not endpoints coincide, whether intervals which meet at or near endpoints overlap one another, etc.

In spite of this potentially serious limitation, it is still often convenient to work with intervals with non-integer endpoints, including data where adjacent intervals exactly meet at a non-integer endpoint. To address this, the *intervals* package takes the following approach:

- Floating point representations of interval endpoints are assumed to be exactly equal (in the sense of == in R or C++) if and only if the user intends the real values corresponding to these representations to be exactly equal.
- For cases where floating point error and approximate equality are a concern, tools are provided to permit distinguishing between ambiguous and unambiguous intersection, union, etc.

In the next example, y1 does not literally overlap y2[2,], although R's all.equal function asserts that the gap between them is smaller than the default tolerance for equivalence up to floating point precision.

```
> delta <- .Machine[[ "double.eps" ]]^0.5</pre>
> y1 <- Intervals( c( .5, 1 - delta / 2 ) )</pre>
> y2 <- Intervals( c( .25, 1, .75, 2 ) )
> y1
Object of class Intervals
1 interval over R:
[0.5, 0.9999999254942]
> y2
Object of class Intervals
2 intervals over R:
[0.25, 0.75]
[1, 2]
> all.equal( y1[1,2], y2[2,1] )
[1] TRUE
> interval_intersection( y1, y2 )
Object of class Intervals
1 interval over R:
[0.5, 0.75]
```

The expand and contract methods, used with type = "relative", permit consideration of the maximal and minimal interval sets which are consistent with the nominal endpoints — from the point of view of endpoint relative difference. The contract method, for example, contracts each interval in a collection so that the relative difference between original and contracted endpoints is equal to tolerance delta. Thus, if a relative difference less than or equal to delta is our criterion for approximate floating point equality, the contract object has endpoints which are approximately equal to those of the original — even though the contract object is a proper subset of the original. The expand methods is similar, but generates a proper superset.

```
> contract( y1, delta, "relative" )
Object of class Intervals
1 interval over R:
[0.500000007450581, 0.999999977648258]
```

We compute two separate intersections which bound the nominal intersection:

Finally, we identify points which may or may not be in the intersection, depending on whether we make a conservative, literal, or anti-conservative interpretation of the nominal endpoints.

```
> interval_difference( outer, inner )
Object of class Intervals_full
3 intervals over R:
[0.499999992549419, 0.500000007450581)
(0.749999988824129, 0.750000011175871]
[0.999999985098839, 1.00000000745058]
```

The expand and contract methods have other uses as well. Here, we eliminate gaps of size 2 or smaller:

```
> x <- Intervals( c(1,10,100,8,50,200), type = "Z" )
> x

Object of class Intervals
3 intervals over Z:
[1, 8]
[10, 50]
[100, 200]
> w <- 2
> close_intervals( contract( reduce( expand(x, w/2) ), w/2 ) )

Object of class Intervals
2 intervals over Z:
[1, 50]
[100, 200]
```

4 Notes on implementation

ullet Endpoint representation. For the moment, interval endpoints are always stored using R's numeric data type. Although this is wasteful from an

memory and speed point of view, we do it for two reasons. First, use of R's Inf and -Inf — not possible with the *integer* type — is very convenient when computing complements. Second, the range of integers which can be represented using the *numeric* data type is considerably greater:

> .Machine\$integer.max

[1] 2147483647

```
> numeric_max <- with( .Machine, double.base^double.digits )
> options( digits = ceiling( log10( numeric_max ) ) )
> numeric_max
```

[1] 9007199254740992

• Efficiency. All computations are accomplished by treating intervals as pairs of tagged endpoints, sorting these endpoints (along with their tags), and then making a single pass through the results. Computational complexity for set operations is therefore $O(n \log n)$, where input object i contains n_i rows and $n = \sum_i n_i$. The same sorting approach is also used for interval_overlap, although if every interval in a query object of m rows overlaps every intervals in a target object of n rows, generating output alone must of necessity be O(mn).

Sorted endpoint vectors are not retained in memory. If one wishes to query a particular object over and over, repeated sorting would be inefficient; in practice so far, however, such repeated querying has not been needed.

5 Session information

- R version 2.7.2 (2008-08-25), i386-apple-darwin8.11.1
- Base packages: base, datasets, grDevices, graphics, methods, stats, tools, utils
- Other packages: intervals 0.9.5