## Scientific Programming with R

## Stephen Eglen

Laurent Gatto

September 6, 2015

## Aims of course

This course aims to teach $R$ as a general-purpose programming language. Issues specific to Computational Biology (e.g. Bioconductor packages) are covered in other course modules.
In part 1, topics to be mastered in this course include:

- Interactive use of R.
- Basic data types: vector, matrix, list, data.frame, factor, character.
- Writing scripts.
- Graphical facilities.
- Writing your own functions.
- File input/output.
- Control-flow statements, looping.
- Vectorization.
- Numerics issues.
- Debugging.

Books and online help

- Introductory Statistics with R (Springer, Dalgaard).
- A first course in statistical programming with R (CUP, Braun and Murdoch).
- Computational Genome Analysis: An Introduction (Springer, Deonier, Tavaré and Waterman).
- S programming (Springer, Venables and Ripley).
- R programming for Bioinformatics (CRC Press, Gentleman).
- Scientific programming and simulation using R (CRC, Jones, Maillardet and Robinson).
- The art of R programming (No Starch Press, Matloff).
- Writing Scientific Software (WSS) (CUP, Oliveira and Stewart).
- www.r-project.org, www.rseek.org, www.r-bloggers.com
- $R$-help mailing list.
- Eglen (2009) http://www.ploscompbiol.org/doi/pcbi. 1000482


## Part 2: Scientific computing issues

In part 2 of the course ${ }^{1}$, we will explore various other topics, building on core knowledge of $R$.

- Numerical integration
- Phase plane analysis
- Handling large files/data bases
- String processing (e.g. for genomic data)
- Advanced graphing / presenting results
- Reproducible research
- Future directions ( R and generally)
- (Object-oriented programming)
- (Package development)
- (R profiling)

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## Outline

## Introduction

Vectors
Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping
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## Scientific programming/software

- Different from software engineering (but should try to adhere to SE best practice, of course).
- Moving target.
- Domain scientists write the code (some argue this is a weakness).
- Has of course to be accurate, user-friendly (no GUI vs CLI ranting here), usable and useful, flexible, efficient and open, owned by the community, facilitate reproducible research.
- Contribute to users education (i.e not be a black box), in terms data requirements, the data processing and result interpret.
Importance of documentation.
See Gentleman et al. (2004) Genome Biology (Bioconductor paper) for an example of successful scientific software.


## What is R ?

- Statistical computing environment, and programming language.
- Very popular in many areas of statistics, computational biology.
- "Programming with data" (Chambers)
- Approach: command-line for one-liners; interactive usage; write scripts/functions for larger work (edit/run cycle); develop package for consolidation and distribution.
- S language came from Bell Labs (Becker, Chambers and Wilks). Commercial version S-plus (1988).
- Remerged as a combination of S and Scheme: Ross Ihaka and Robert Gentleman (NZ).
- 1993: first announcement.
- 1995: 0.60 release, now under GPL.
- 2014-10-07: release 3.1.1. Stable, multi-platform. Major release every March/April.
- R-core now 20 people, key academics in field, including John Chambers.


## Graphics example




Jean YH Yang; gpQuality
http://bioinf.wehi.edu.au/marray/ibc2004/lect1b-quality.pdf

- GPL'd, available on many platforms.
- Excellent development team with yearly release cycle.
- Source always available to examine/edit.
- Fast for vectorized calculations.
- Foreign-language interface (C/Fortran) when speed crucial, or for interfacing with existing code
- Good collection of numerical/statistical routines.
- Comprehensive R Archive Network (CRAN) ~ 5913 packages [2014-10-07] (cf 1000 in April 2007).
- On-line doc, with examples.
- High-quality graphics (pdf, postscript, quartz, $\times 11$, bitmaps). Often used just for plotting.


## Weaknesses of $R$

- Loops are slow. Learn how to vectorize solutions.
- No fast compiler yet, and unlikely to happen due to nature of language. Byte compiler available in compiler package.
- No (decent) endorsed GUI built-in to R. Tk is available within base R, and packages for other graphical tooklits (e.g. Gtk2, Qt) are also available.
"Programming Graphical User Interfaces with R ", M. F. Lawrence and J. Verzani
- Flexible language, similar to matlab, but definitely not "everything is a matrix". Frames, lists, vectors ...
- From matlab to R :
http://cran.r-project.org/doc/contrib/R-and-octave.txt
- Comprehensive matlab and R guide: http:
//www.math.umaine.edu/faculty/hiebeler/comp/matlabR.html
- Use $\mathrm{x}[\mathrm{i}]$ not $\mathrm{x}(\mathrm{i})$ for indexing vectors.
- Making vectors: $\mathrm{x}<-\mathrm{c}(10,9,5,1)$
- Assignment: best to use $<-$ rather than $=$.


## My very first $R$ session

- Start-up: type 'R' at command line.
- Type commands interactively, and get results.
- Type commands into a file; source(' myfile.R'); edit file ...
- Mac/Win has a GUI for interactive use, with internal editors.
- All platforms have a command-line interface
- Many external editors have support for $R$, including
- Emacs through ESS (http://ess.r-project.org),
- Eclipse IDE (http://www.walware.de/goto/statet),
- Rstudio (http://www.rstudio.org),
- ...


## Interacting with R

- Can use up/down arrow keys to go through command history. Within a command, use left/right arrow keys to edit.
- History can be saved over sessions (?history).
- Multiple commands can be put onto one line, using ; as separator between lines, e.g. $x<-10 ; y<-3$; $a<-5$.
- TAB can do object/file completion.

R manipulates objects. Each object has a name and a type (vector, matrix, list,...)
Name of an object: letters (upper/lower case are distinct), digits, period. Start with a letter.
Objects set by way of assignement. Use the <- assignment operator rather than $=$ wherever possible. (Does $\mathrm{i}=\mathrm{i}+1$ make sense?)

What's up with the assignment and underscore?
(Advanced)

Historically, underscore was used in S for assignment (because an old system keyboard had a key equivalent to the ASCII underscore that generated a back arrow). Hence underscore was not used within variables.
More recently, = is now available as an assignment operator (similar to languages like C), but is frowned upon as it can be confusing.
What does $i=i+1$ imply mathematically?
Better to stick to i <- i + 1 and use equals just within calls to functions, e.g. runif (max=3).

Note also that assignments return values:

```
>y<-1 + (x<- 9)
> a<- b <- 0
```

http://developer.r-project.org/equalAssign.html

## Objects and Functions

```
> x <- 200
```

> x <- 200
> half.x <- x/2
> half.x <- x/2
> threshold <- 95.0
> threshold <- 95.0
> age <- c(15, 19, 30)
> age <- c(15, 19, 30)
> age[2] \#\# [] for accessing element.
> age[2] \#\# [] for accessing element.
[1] 19
> length(age) \#\# () for calling function.
[1] 3

```

\section*{Outline}

\section*{Introduction}

\section*{Vectors}

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Matrices
Boolean logic
Lists
Factors
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\section*{Vectors}

Vectors are a fundamental object for R. Scalars are treated as vector of length 1.
\(>y<-c(10,20,40)\)
\(>y[2]\)
[1] 20
> length (y)
[1] 3
> \(\mathrm{x}<-5\)
\(>\) length \((x)\)
[1] 1

\section*{Generating vectors}

Many short hand methods for regular sequences; c() for irregular.
```

> x <- seq(from=1, to=9, by=2)
> y <- seq(from=2, by=7, length=3)
> z<- 4:8
> a <- seq.int(5) \#\# fast for integers
> b <- c(3, 9, 2)
> d <- c(a, 10, b)
>e <- rep(c(1,2), 3)
> <- integer(7)

```

\section*{Vectors}

Some operations work element by element, others on the whole vector, compare:
```

> y <- c(20, 49, 16, 60, 100)
min(y)
[1] 16
> range(y)
[1] 16 100
> sqrt(y)
[1] 4.472136 7.000000 4.000000 7.745967 10.000000
> log(y)

```
[1] 2.9957323 .8918202 .7725894 .0943454 .605170

\section*{Accessing and setting elements}
```

> x <- seq(from=100, by=1, length=20)
> x[3] \#\# just element 3.
[1] }10
> x[c(12,14)] \#\# element 12 and 14
[1] 111 113
> x[1:5]
[1]}100100101 102 103 104
> bad <- 1:4
> x[-bad] \#\# exclude elements
[1] 104 105 106 107 108 109 110 111 112 113 114 115 116 117
[15] 118}11

## Accessing and setting elements

Accessing and setting elements

Can also provide a logical vector of same length as vector (logical values explained later).

```
> x <- c(5, 2, 9, 4)
> v <- c(TRUE, FALSE, FALSE, TRUE)
> x[v]
```

[1] 54

## Elements can be set in several ways

```
> x <- rep (0,10)
> x[1:3] <- 2
> x[5:6] <- c(-5,NA)
> x[7:10] <- c(1,9) ## recycling.
```


## Recycling rule (Advanced)

Recycling is convenient, but dangerous; when vectors are of different lengths, the shorter one is often recycled to make a vector of the same length.

```
> a <- c(1,5) + 2
> x <- c(1,2); y <- c(5,3,9,2)
> x + y
[1] }
> x + c(y,1) ## odd recycling, warning.
Warning in x + c(y, 1): longer object length is not a
multiple of shorter object length
[1] }
```

```
> x <- 1:10
> y<- x * 2
> z <- x^2
> y + z
>x+1:2
> x + 1:3
```

Recycling rule (Advareces)

```
> x <- 1:10
y y <- x * 2
> < <- x^2
> y + z
[1] \begin{tabular}{lllllllllll}
3 & 8 & 15 & 24 & 35 & 48 & 63 & 80 & 99 & 120
\end{tabular}
```

```
> x + 1:2
```

[1] $\begin{array}{llllllllll}2 & 4 & 4 & 6 & 6 & 8 & 8 & 10 & 10 & 12\end{array}$
$>x+1: 3$
Warning in $x+1: 3:$ longer object length is not a multiple
of shorter object length
[1] $\begin{array}{lllllllllll}2 & 4 & 6 & 5 & 7 & 9 & 8 & 10 & 12 & 11\end{array}$

Naming indexes of a vector

```
> joe["height"] == joe[2]
height
    TRUE
```

Refering to index by name rather than by position can make code more readable, and flexible. Cannot do things like $x[1: 4]$ easily though, since you need to name all four elements you want.
Although extremly useful, names have a cost when processing large objects.
Note: in second use of names () above, we are actually using the replacement function names<-, see later.

Naming indexes of a vector

```
> joe <- c(24, 1.70)
> joe
[1] 24.0 1.7
> names(joe)
NULL
> names(joe) <- c("age", "height") ## replacement function
> joe
    age height
24.0 1.7
```


## Common functions for vectors

- length ()
- rev()
- sum(), cumsum(), prod(), cumprod()
- mean(), sd(), var(), median()
- min(), max (), range(), summary ()
- $\exp (), \log (), \sin (), \cos (), \tan ()$ (radians, not degrees)
- round(), ceiling(), floor(), signif()
- sort(), order(), rank()
- which(), which.max()
- any(), all()


## Functions as function args

Functions can be called within function calls; the following are equivalent:

```
> x <- c(3, 2, 9, 4)
>
> y <- exp(x); z1 <- which(y > 20) ## case 1
> z2 <- which ( exp(x) > 20) ## case 2
>
> all.equal(z1, z2)
```

[1] TRUE

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Default values for function arguments

A function will error if not all required arguments are provided. Some functions have both required and optional arguments. If the optional arguments are not provided, they are either ignored, or they take a default value.

Usage:
round $(x$, digits $=0)$

Default values for function arguments

```
> x <- c(2.091, 4.126, 7.925)
> round() ## required arg is missing
Error in eval(expr, envir, enclos): 0 arguments passed to
'round'which requires 1 or 2 arguments
> round(x)
[1] 248
> round(x, digits = 2)
[1] 2.09 4.13 7.92
```

Let's see how this works in more detail.

## Argument matching

$R$ has a flexible method for specifying arguments to function. We can either provide an actual value for a formal argument, or give arguments as key=value (or formal=actual).
As an example, let's look at help for seq:

```
seq(from = 1, to = 1, by = ((to - from)/(length.out - 1)),
```

    length.out = NULL, along.with = NULL, ...)
    (NB: in seq(from=x), from is the formal argument of the function, and here x is the actual value.)

The . . . notation allows for other arguments to be passed, which are not used by this function.

## Argument matching Typical calls are as follows:

```
> seq(1, 3, 0.5) ## positional matching
[1] 1.0 1.5 2.0 2.5 3.0
> seq(1, 5,length.out=3) ## can skip args (e.g. by)
[1] 1 3 5
> seq(to=5) ## order not important.
[1] 1 2 3 4 5
> seq(f=5,t=1) ## abbrev tags
[1] 543 2 1
> seq(len=5, 1,2) ## tags removed before positional matching
```

[1] $1.001 .251 .501 .75 \quad 2.00$

Replacement functions (Adanecea)

```
x}<-1:
> X
[1] 1 2 3 4 5
```

> length (x)
[1] 5
> length $(x)$ <- 2
$>\mathrm{X}$
... in function calls (Advanced)
Why do some functions, like sqrt, require only one argument, yet others take many arguments?
Functions like c, cbind, have . . . in the arguments:
Usage:

```
    c(..., recursive=FALSE)
```


## Arguments:

...: objects to be concatenated.

The ... indicate any number of objects may be passed, not just (say) one or two.
The result of $c()$ is to combine them all into one long vector, taking into account if the keyword "recursive" is provided (when args are first flattened). The ... can also indicate that other arguments can be provided which are not processed directly by this function, but may be useful for other functions (e.g. popular when plotting).

Normally length ( x ) would return a value, rather than you assigning a value to the function! These are replacement functions, see help page:

## Usage:

length ( x )
length (x) <- value

- help(hist) to see help file (or ?hist).
- $\operatorname{args}$ (hist) to see arguments of a function.
- example(boxplot) run examples in help page.
- options(help_type="html") will then use web-browser for help.
- help.search("histogram")
- demo() to list all demos, e.g. demo(graphics)

NB: In the R terminal ?command works as shorthand for help("command") except for a small number of commands, e.g. if, while. Use the longhand for these.

- What you can expect to find:
- Description - one line summary
- Usage - formal arguments
- Arguments - interpretation of arguments
- Details - what the function does
- Value - return value.
- References - documentation
- See also - helps you find related pages
- Examples - guaranteed to run: example(hist)


## Numbers and special values

- numeric (floating-point, double): 12, 4.92, 1.5e3 - is.numeric() (integers converted to f.p.)
- integers 1 L - is.integer()
- complex: 3+2i - is.complex()
> typeof(1)
[1] "double"
> typeof (1L)
[1] "integer"
> is.integer(1)
[1] FALSE
> is.integer(1L)
[1] TRUE

Numbers and special values
Numbers and special values

```
> typeof(NA)
```


## Special values:

- NA: not available. (Often used to represent missing data point) is.na()
- NaN: not a number. e.g. $0 / 0-$ is.nan()
- Inf, -Inf: $\pm \infty$-is.finite()

You will also meet:

- NULL: often, list of zero length - is.null()
[1] "logical"
> typeof(NaN)
[1] "double"
> typeof(Inf)
[1] "double"
> typeof(NULL)
[1] "NULL"

Operator precedence ?Syntax
$>3 * 4+2!=3 *(4+2)$
[1] TRUE
$>2^{\wedge} 3+1 \quad!=2^{\wedge}(3+1)$
[1] TRUE
> 1:5-1
[1] 01234

## Operator precedence ?Syntax

Subset taken from ?Syntax, see that page for full list. Highest precedence at top.

| ' [ [ ${ }^{1}$ | indexing |
| :---: | :---: |
| '\$ @ ' | component / slot extraction |
| $1 \sim 1$ | exponentiation (right to left) |
| $1-+1$ | unary minus and plus |
| ' : ' | sequence operator |
| $1 \%$ any ${ }^{\prime}$ | special operators |
| '* /' | multiply, divide |
| $1+-1$ | (binary) add, subtract |
| $1<><=>==1=1$ | ordering and comparison |
| '! ' | negation |
| '\& \&\&' | and |
| '\| ||' | or |
| $1<-\ll-1$ | assignment (right to left) |
| '?' | help (unary and binary) |

## Operators

## Most operators will be familiar, but some may not:

```
x <- 10
x == 4 ## test for equality
x != 10 ## not equal?
%/% 2 ## division, ignoring remainder. (3)
7%% ## remainder (1)
x <- 9 ## assignment
<<- 9# assign }x\mathrm{ to }9\mathrm{ in the global env. (BAD)
## Raising to a power can be done in two ways.
all.equal(10.1**2.5, 10.1^2.5 )
```

Bottom line: use parentheses to order preference.

## When things go wrong

Syntax errors are those where you've just made a typing mistake.
Logical errors are harder to find!

## Common problems:

- missing close bracket leads to continuation line.

$$
>\times<-(1+(2 * 3)
$$

$+$
Hit Ctrl C (below) or keep typing!

- too many parens:
$2+(2 * 3))$
- wrong/mismatched brackets (see next slide).
- Likewise, do not mix double quotes and single quotes. (Unless you need to quote within quotes.)
- ...
- wrong variable name (not syntax error)
- When things seem to take too long, try C-c [Ctrl and C, together]
- Typing in commands interactively is good for one-liners, but soon you will want to switch to putting your sequence of commands into a script file, and then ask $R$ to run (source) those commands.
- This leaves to a rapid edit-run-edit cycle.
- e.g. type these commands into a file trig.R:

```
x <- seq(from=0, to=2*pi, length=100)
y <- sin(x)
z <- cos(2*x)
z ## will not appear when source'd
print(y[1:10]) ## should use print()
plot(x, y, type='l')
lines(x, z, type='l', col='red')
```

- Eval within $R$ using source('trig.R').


## Scripts

- Use source('trig.R', echo=TRUE) to see commands and output. Or use print ( x ) to print an object within a script.
- Keep your code open in the editor in one window, and keep R running in another window.
- Are you in the right directory? Check that you can see your script file in the same directory as where $R$ is currently. Check $\operatorname{dir}()$, and setwd, see later.
- On unix, the initial directory is the directory from where you started $R$ On windows, the initial directory might be "My Documents". You may need to change directory (setwd) first.
- Use a good editor that helps you spot mistakes (e.g. paren matching, syntax highlighting). Examples: Emacs/ESS, gedit, Rstudio.
- Use . R or .r as the filename suffix. Avoid any temptation to put spaces (although $R$ does not mind) in your filenames!


## Outline

ntroduction

## Vectors

## Calling functions

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Matrices
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- You don't have to remember what commands you ran, they are saved in the file.
- This corresponds to the "source is real" philosophy of using S and R .
- You can easily give your work to others, by passing them the file.
- You can eventually run your scripts in BATCH, i.e. non-interactively. Good for long jobs which you can leave overnight.
- At the command line, type R CMD BATCH trig.R. R will start up, process your commands and then quit.
- Output is stored in the file trig.Rout
- If there were no errors, the last line of the output file shows the time taken to run the script.
- Any output is not shown on the screen but sent to a PDF called Rplots.pdf.
- This is a GREAT way of testing your scripts, since $R$ starts with an empty workspace, you will see if you have all the steps needed.
- Aim to always leave your scripts in a working state at the end of a session, so that a few days later you don't have to remember why it wasn't working!


## Rscript

- Rscript works as an interpreter. It is quicker to start than $R$ (fewer packages are loaded?).
- Can use interactively or for standalone scripts.

```
$ Rscript -e 'round(runif(10))'
    [1] 0 1 0 0 1 0 1 1 0 0
```

```
#!/usr/bin/env Rscript
args <- commandArgs(TRUE)
```

stopifnot(length(args)==3)
args = as.numeric(args)
$\mathrm{n}=\operatorname{args}[1] ;$ mean $=\operatorname{args}[2] ;$ sd $=\operatorname{args}[3]$
rnorm(n, mean, sd)
\$ ./simple_rnorm.R 5102
[1] $8.65180713 .37209410 .06334710 .703155 \quad 6.351886$

## Commenting your work

- Do not be shy when putting comments into your code.
- Meaningful variable names help, but do document. At a bare minimum, each file should state at the top what the purpose of the file. Important variables and functions should be clearly documented.
- You may think it obvious how your code works, but try looking at it a week or a month later and then see if you clearly understand it. If in doubt, document it.
- Describe what your code is doing, not how it is doing it (WSS, p79). Compare the following two:

```
> s <- s + 1 ## prepare to process next subject
> j<- j + 1 ## increment j by 1.
```

- Comments can be put before commands, if you temporarily do not want to run that command; remove the comments when you want to run the command again, or delete the line.

```
> ## x <- c(x,c(1,2,3))
```

- Line-wrapping. Do not write beyond around column 72, for readability. You can break long expressions at suitable points.
- End of line shold not look like end of an expression. Compare:

```
x <- sqrt( c(100, 200, 300, 400, 500) ) + 10
x <- sqrt( c(100, 200, 300, 400, 500) )
+ 10
x <- sqrt( c(100, 200, 300, 400, 500) ) +
10
x <- sqrt( c(100, 200, 300, 400, 500) ) +
    10
```


## Line wrapping

> \#\# 1: ok - all fits onto one line, just
$>(\mathrm{x}<-\operatorname{sqrt}(c(100,200,300,400,500))+10)$
[1] 20.0000024 .1421427 .3205130 .0000032 .36068
> \#\# 2: not okay -- first line is seen as complete.
$>x<-\operatorname{sqrt}(c(100,200,300,400,500))$
$>+10$
[1] 10
$>\mathrm{X}$
[1] $10.0000014 .1421417 .32051 \quad 20.00000 \quad 22.36068$

## Outline

introduction
Vectors
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## Matrices

Boolean logic
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Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping
Vectorization
Random number generation
Debugging

## Line wrapping

```
> ## 3: solved, by moving the operator (+) up
> x <- sqre( c(100, 200, 300, 400, 500) ) +
+ 10
> X
[1] 20.00000 24.14214 27.32051 30.00000 32.36068
> ## 4: as 3, but indentation makes it clearer.
> x <- sqrt( c(100, 200, 300, 400, 500) ) +
+ 10
>
[1] 20.00000 24.14214 27.32051 30.00000 32.36068
```


## Matrices

A matrix is just a vector with some additional markup to reformat it. Matrix stored in column-major order (like fortran, unlike C)

```
x <- 1:6
> is.matrix(x)
1] FALSE
> dim(x) <- c(2,3)
> is.matrix(x)
[1] TRUE
> x
[r,1] [,2] [,3]
[2,] 2 4 6
> dim(x)
[1] 2 3
```


## Matrices

```
> x[2,2] ## extracting a value
```

[1] 4
> $\mathrm{x}[1$,$] \#\# extracting row$
[1] 135
$>x[1: 2,2: 3]$
[,1] [,2]
$[1] \quad 3 \quad$,
$[2] \quad 4 \quad$,
> $\mathrm{x}[, 2]$ \#\# not column vector!
[1] 34

```
> x[, 2, drop=TRUE] ## default
1] }3
> x[, 2, drop=FALSE] ## gotcha!
[,1]
[2,] 4
> x[, 2, drop=TRUE] \#\# default
[1] 34
> x[, 2, drop=FALSE] \#\# gotcha!
[,1]
\([2] \quad\),
```

Typical matrix construction methods

```
> matrix( floor(runif(6, max = 50)), nrow = 3 ) ## ncol = 2
```

    [,1] [,2]
    ```
    [,1] [,2]
[1,] \(43 \quad 37\)
[1,] \(43 \quad 37\)
\([2] \quad 6 \quad\),
\([2] \quad 6 \quad\),
[3,] 47 48
```

[3,] 47 48

```
\begin{tabular}{lrr} 
& {\([, 1]\)} & {\([, 2]\)} \\
{\([1]\),} & 43 & 37 \\
{\([2]\),} & 6 & 8 \\
{\([3]\),} & 47 & 48
\end{tabular}
```

- matrix()
- cbind()
- rbind()
$>\mathrm{m}<-\operatorname{matrix}(\mathrm{floor}($ runif $(6, \max =50)$ ), nrow=3)
$>\mathrm{x}<-\operatorname{rbind}(\mathrm{c}(1,4,9), c(2,6,8), c(3,2,1))$
$>y<-\operatorname{cbind}(c(1,2,3), 5, c(4,5,6))$
$>$ rbind $(c(1,4,9), c(2,6,8), c(3,2,1))$

|  | $[, 1]$ | $[, 2]$ | $[, 3]$ |
| :--- | ---: | ---: | ---: |
| $[1]$, | 1 | 4 | 9 |
| $[2]$, | 2 | 6 | 8 |
| $[3]$, | 3 | 2 | 1 |

## Typical matrix construction methods

Note that matrix indices can also be named:

```
> dimnames(m) <- list(student = c("ann", "bob", "joe"),
+ exam = c("math", "french"))
> m
student math french
\begin{tabular}{llr} 
ann & 19 & 21 \\
bob & 35 & 3 \\
joe & 12 & 12
\end{tabular}
> m["bob",] ## get bob's scores
math french
    35 3
```

```
> cbind( c(1,2,3), 5, c(4,5,6) ) ## recycling again
\begin{tabular}{lrrr} 
& {\([, 1]\)} & {\([, 2]\)} & {\([, 3]\)} \\
{\([1]\),} & 1 & 5 & 4 \\
{\([2]\),} & 2 & 5 & 5 \\
{\([3]\),} & 3 & 5 & 6
\end{tabular}
> cbind( c(1,2,3), 5, c(4,5,6) ) \#\# recycling again
```


## Common matrix operations

- diagonal: diag(x) - watch if $x$ matrix or scalar!
- matrix multiplication: $\% * \%$ vs * (element-wise)

```
> x <- matrix(1:4, 2,2)
> i <- diag(2) ## 2 x 2 identity matrix
> x %*% i ## should be x
```

|  | [, 1] | [,2] |
| :---: | :---: | :---: |
| [1, ] | 1 | - 3 |
| [2,] | 2 | 2 |
| > x | * i |  |
|  | [,1] | [,2] |
| [1, ] | 1 | - 0 |
| [2, ] | 0 |  |

## Common matrix operations

## Arrays

Arrays as extension of matrices to multiple dimensions.

- transpose: $\mathrm{t}(\mathrm{x})$
- dim, row, ncol
- inverse: solve(x),
> (x \%*\% solve(x)) == diag(nrow(x))
[,1] [,2]
[1,] TRUE TRUE
[2,] TRUE TRUE


## Outline

introduction
Vectors
Calling functions
Scripts

## Matrices

## Boolean logic

Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping
Vectorization
Random number generation
Debugging

$$
>\operatorname{array}(1: 12, c(2,2,3))
$$

, , 1
$\begin{array}{lrr} & {[, 1]} & {[, 2]} \\ {[1,]} & 1 & 3 \\ {[2,]} & 2 & 4\end{array}$
, 2
$\begin{array}{lrr} & {[, 1]} & {[, 2]} \\ {[1,]} & 5 & 7\end{array}$
$[2] \quad 6 \quad$,
, 3
$\begin{array}{ll} & {[, 1]}\end{array} \begin{array}{r}{[, 2]} \\ {[1,]} \\ 9\end{array} 11$
$\begin{array}{lrr}{[2,]} & 10 & 12\end{array}$

Boolean values - ?logical
Logical values TRUE/FALSE (avoid abbrev to $\mathrm{T} / \mathrm{F}$ ).
TRUE/FALSE equivalent to $1 / 0$
> as.integer (TRUE)
[1] 1
> as.integer (FALSE)
[1] 0
> as.logical(1)
[1] TRUE
> as.logical(0)
[1] FALSE
$>\mathrm{d}<-\mathrm{c}(3.2,1.0,4.0,9.2,2.3,8.1,6.3)$
$>\mathrm{d}>5.0$
[1] FALSE FALSE FALSE TRUE FALSE TRUE TRUE
$>d[d>5.0]$
[1] $9.28 .1 \quad 6.3$
$>$ which ( $\mathrm{d}>5.0$ )
[1] 467

## Boolean values - ?logical

Key operators for handling boolean values:
> ! TRUE
\#\# negation: swap $T$-- $F$.
[1] FALSE
> TRUE \& FALSE
\#\# and: both must be true.
[1] FALSE
> FALSE | TRUE
\#\# or: one must be true.
[1] TRUE
> xor(TRUE, TRUE)
\#\# xor: only one is true.
[1] FALSE

## Boolean values - ?logical

```
> d[which(d>5.0)]
```

[1] 9.28 .16 .3
$>$ medium.sized <- (d > 3.0) \& ( $\mathrm{d}<\mathrm{5} .0$ )
> d[medium.sized]
[1] 3.24 .0
> d[!medium.sized]
[1] 1.09 .22 .38 .16 .3
> ifelse(d > 3.0, 1.0, 0.0) \#\# same as as.numeric(d > 3)
[1] 1011011

## Boolean logic: issues

$\mathrm{a} \& \mathrm{~b}$ (same for $\mathrm{a} \mid \mathrm{b}$ ) is an elementwise operation, with a result the same length as the longer of $\mathrm{a}, \mathrm{b}$ (recycling is used if one vector is shorter).
$\mathrm{a} \& \& \mathrm{~b}$ examines only the first element of a and b , returning one logical value. Lazy evaluation is used: we calculate only what's needed to determine result.
> TRUE \| $\mid$ some.long.computation()
[1] TRUE
> TRUE \&\& stop("no")
Error in eval(expr, envir, enclos): no
> FALSE \&\& stop("no")
[1] FALSE

Lazy (Advancea)

R uses Lazy evaluation, which delays the evaluation of an expression (here the argument) until its value is actually required [ ${ }^{2}$ ]:

```
> f <- function(x) { 10 }
> system.time(f(Sys.sleep(3)))
    user system elapsed
            0 0 0
> f <- function(x) { force(x); 10 }
> system.time(f(Sys.sleep(3)))
    user system elapsed
    0.000 0.000 3.001
```

${ }^{2}$ example from Hadley Wickham's devtools

## Outline

## introduction

Vectors
Calling functions
Scripts
Matrices
Boolean logic

## Lists

Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping
Vectorization
Random number generation
Debugging

Boolean logic: issues

Comparing numbers: When testing numbers for equality, can use $\mathrm{x}==\mathrm{y}$ when $\mathrm{x}, \mathrm{y}$ are integers, otherwise use all.equal ( $\mathrm{x}, \mathrm{y}$ ). See later on numerics.

Avoid using F
$>\mathrm{F}<-3$
> F == FALSE
[1] FALSE

## What is a list?

A list is used to collect a group of objects of different sizes and types. Very flexible. Often returned as the result of a complex function (e.g. model fit) to return all relevant information in one object.

```
> l <- list(who='joe', height=1.70, dob=c(1960, 12, 1))
> l
$who
[1] "joe"
```

\$height
[1] 1.7
\$dob
[1] $1960 \quad 12 \quad 1$

## What is a list?

```
> length(l)
```

[1] 3
> names(l) \#\# show components
[1] "who" "height" "dob"
> l\$height \#\# access an element.
[1] 1.7

What is a list?

```
> unlist(l) ## opposite of list
    who height dob1 dob2 dob3
    "joe" "1.7" "1960" "12" "1"
> str(l) ## structure of l
List of 3
    $ who : chr "joe"
    $ height: num 1.7
    $ dob : num [1:3] 1960 12 1
```

What is a list?

List elements can either be accessed by name (e.g. l\$height or l['height'] - if named list) or by position (l[ [2]]).

When using numbers to index list, compare 1[2] (a list with one element) with $1[[2]]$. You can therefore do $1[2: 3]$ but not $1[[2: 3]]$.

## Modifying lists (Aduancea)

We can append new items to list either by making a new list from the old one (e.g. 1), or directly by assigning new element (e.g. 2):

```
> l1 <- list(who="fred")
```

$>11<-c(11$, height=1.8) \#\# e.g. 1
> $11[[" d o b "]]<-c(1965,10,17)$ \#\# e.g. 2

## Modifying lists (Advanced)

## Deleting list items:

```
> l1["height"] <- NULL
> str(l1)
List of 2
    $ who: chr "fred"
    $ dob: num [1:3] 1965 10 17
```


## Initialisation

Initialising R object (list, vectors, ...) is much faster than creating and extending these objects on the fly

```
> n <- 1e4
> l <- list()
> system.time(for (i in 1:n) l[[i]] <- rnorm(1e3) )
    user system elapsed
    2.728 0.045 2.776
> l <- vector("list", n)
> system.time(for (i in 1:n) l[[i]] <- rnorm(1e3) )
    user system elapsed
1.211 0.005 1.215
```

Finally, for completeness, here is a way to predefine a list of given length and gradually fill it in:

```
> empty <- vector("list", 3) ## Prealloc to given length.
> names(empty) <- c("who", "height", "dob")
> empty[["height"]] <- 1.8
```


## Data frames

A data frame is like a matrix, but each column can be of a different type.
Data frame is stored as a list, with each element a vector of same length. Useful for reading in tabular data from a file (see read.csv).

```
> nms <- c("joe", "fred", "harry")
> a <- c(24, 19, 30)
> ht <- c(1.7, 1.8, 1.75)
> s <- c(TRUE, FALSE, TRUE)
```


## Data frames

```
> cbind(a, ht) ## matrix, preserves
```

|  | a | ht |
| :--- | ---: | ---: |
| $[1]$, | 24 | 1.70 |
| $[2]$, | 19 | 1.80 |
| $[3]$, | 30 | 1.75 |

> class(cbind(a, ht))
[1] "matrix"

## Data frames

```
> d <- data.frame(name = nms,
+ age = a,
+ height = ht,
+ student = s)
> class(d)
[1] "data.frame"
```


## Data frames

```
> cbind(nms, ht) ## matrix, ht col becomes strings.
        nms ht
[1,] "joe" "1.7"
[2,] "fred" "1.8"
[3,] "harry" "1.75"
> class(cbind(nms, ht))
[1] "matrix"
```


## Data frames

```
> d$age ## same as d[, "age"]
```

[1] 241930
$>$ names (d)
[1] "name" "age" "height" "student"
> d[2,] \#\# access 2nd row.
name age height student
2 fred 19 1.8 FALSE

Data frames
Compare how a data frame (d) is printed, compared to printing as.list(d).
$>d$

\$height
1] 1.701 .801 .75
\$student
[1] TRUE FALSE TRUE
(Mostly seen when reading in data frames)
Factors internally code categorical variables with a number. e.g. 1=Sunday, $2=$ Monday, $\ldots 7=$ Saturday. For large vectors, this is more efficient storage, especially when character strings repeat. Can also make code more readable.

Also useful in many statistical functions, using the formula interface.

## Outline

ntroduction
Vectors
Calling functions
Scripts
Matrices
Boolean logic
Lists

## Factors

Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping
Vectorization
Random number generation
Debugging

## Factors (Advanced)

```
> scores1 <- c('good', 'poor', 'bad', 'poor',
+ 'bad', 'bad', 'good')
> scores <- factor(scores1)
> scores
[1] good poor bad poor bad bad good
Levels: bad good poor
> levels(scores)
```

[1] "bad" "good" "poor"
> as.integer(scores) \#\# integer representation
[1] $2 \begin{array}{lllllll}2 & 1 & 3 & 1 & 1 & 2\end{array}$
> as.character(scores) \#\# show strings
[1] "good" "poor" "bad" "poor" "bad" "bad" "good"

## Factors (Advanced)

```
> which(scores1 == 'bad')
```

[1] 356
> \#\# Can do further comparisons with an ordered factor.
> \#\# Levels are now ordered, as shown by "<" in levels.
> s2 <- factor(scores1,

+ levels = c('poor', 'bad', 'good'),
$+\quad$ ordered $=$ TRUE)
> s2[1] > s2[2]
[1] TRUE


## Strings / character arrays

## Character arrays are vectors of strings.

Use single (') or double (") quotes to mark strings, but don't mix

```
> x <- 'good'
> z <- "no' # need to match"
> z <- "it's working"
```


## Outline

ntroduction
Vectors
Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors

## Character arrays

Objects in your environment
Basic plotting
Reading/writing clata to file system
Writing functions
Conditionals and looping
Vectorization
Random number generation
Debugging

## Strings / character arrays

Within a script, easy way to generate output:

```
> cat("Now computing the steady-state\n")
Now computing the steady-state
> x <- 134
> cat("sqrt of", x, "is", sqrt(x), "\n")
sqrt of 134 is 11.57584
> cat("sqrt of", x, "is", sqrt(x), "\n", sep='__')
sqrt of__134__is__11.57584__
```

See also message, warning and stop for communicating diagnostic messages - cat and print are generally used when displaying an object.

Strings / character arrays
blackslash characters allow you to generate control characters, importantly: newline: $\backslash n$, tab: $\backslash t$.

```
> cat("5\t9\n_")
```

59
-
paste() returns a string, e.g. for assignment.
$>x<-1: 5 ; ~ e x p . d i r ~<-~ ' / h o m e / s t e p h e n / r e s ' ~$
> file <- paste(exp.dir, '/expt_res', x, '.dat', sep='')
> file
[1] "/home/stephen/res/expt_res1.dat"
[2] "/home/stephen/res/expt_res2.dat"
[3] "/home/stephen/res/expt_res3.dat"
[4] "/home/stephen/res/expt_res4.dat"
[5] "/home/stephen/res/expt_res5.dat"

## Don't forget paste0() where sep=".

(Humour): http://simplystatistics.org/2013/01/31/
paste0-is-statistical-computings-most-influential-contributior

## Strings

- Just as R stores vectors of numbers, it also stores vectors of strings.
- Pattern matching facilities are available, based on Unix terms (grep, regular expressions). These are worth learning:
> s <- c('apple', 'bee', 'cars', 'danish', 'egg')
$>$ nchar(s)
[1] 53463
> substr(s, 2,3)
[1] "pp" "ee" "ar" "an" "gg"


## Strings

```
> grep('e', s)
[1] 1 2 5
> grep(1^e', s) ## regexps...
[1] 5
> sub('e', '_', s)
[1] "appl_" "b_e" "cars" "danish" "_gg"
> gsub('e', '_', s) ## global sub, watch "bee"
[1] "appl_" "b__" "cars" "danish" "_gg"
```


## Strings

> toupper (s)
[1] "APPLE" "BEE" "CARS" "DANISH" "EGG"
> sprintf('name \%s len \%d', s, nchar(s)) \#\# C users!
[1] "name apple len 5" "name bee len 3"
[3] "name cars len 4" "name danish len 6"
[5] "name egg len 3"

## Outline

ntroduction
Vectors
Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping
Vectorization
Random number generation
Debugging
109 / 191

## Environments (Advanced)

An environment is a frame, or collection of named objects (variables), and a pointer to an enclosing environment.

The working environment of your interactive $R$ session is the <R_GlobalEnv>.

```
> x <- 2
\(>\operatorname{ls}()\)
```

[1] "x"
> \#\# current environment
> environment()
<environment: R_GlobalEnv>
[1] TRUE

```
> e <- new.env()
```

> e <- new.env()
> e
> e
<environment: 0x7ff9e22741f0>
<environment: 0x7ff9e22741f0>
> ls(e)
> ls(e)
character(0)
character(0)
> parent.env(e)
> parent.env(e)
<environment: R_GlobalEnv>
<environment: R_GlobalEnv>
>e$x <- 1
>e$x <- 1
> ls(e)
> ls(e)
[1] "x"
[1] "x"
> e\$x != x

```
> e$x != x
```

Inspecting variables and the environment

```
x <- 9
y<- c(2,4,5)
m <- matrix(2:5, 12,10)
objects() ## what vars do I have?
ls() ## shorthand for objects.
str(m) ## Display compact representation
head(m)
tail(m)
rm(list = ls()) ## clear up the working environment
ls()
```


## What is an object?

- An object is typically either a variable or a function.
- You can use the same name for a function and a variable, and $R$ uses context to decide which you mean:

```
sum <- 3 + 4 + 5
> total <- sum(1:4)
> total
```

> sum
> sum(sum) \#\# can get confusing!

Some ways of finding out what kind of object you have and converting.

```
mode(y)
typeof(y)
class(y)
object.size(y)
is.vector(y) ## is.xyz() check if type XYZ
is.matrix(y)
as.vector(m) ## convert object to type XYZ
l <- list (x=c(1,4), y=c(6,9,12))
unlist(l)
```


## OO programming in $R$

Data abstraction; object manipluation becomes independent of the implementation details.

There are several OO programming frameworks in R . The main ones, supported in base R are S3, S4 and S4 ReferenceClasses.

Bioconductor provides many ad hoc classes to store, manipulate and process microarray, RNA Seq, proteomics, flow cytometry, ...data.

It there is interest, we could have a session about OO programming at the end of the course.

## Outline

Vectors
Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping

## Vectorization

Random number generation
Debugging

## Basic plotting

- Basic x,y plots (scatter plots)
- Multiple plots in one figure
- Saving your plots

This section will just introduce the mechanics of making basic plots, rather than worry about interpreting them.

## Basic plotting

$>\mathrm{x}<-\operatorname{seq}($ from $=0$, to $=2 * \mathrm{pi}$, len=1000)
$>y<-\cos (2 * x)$
> plot (x,y) \#\# just provide data; sensible labelling

## Basic plotting

> \#\# Expand on previous plot
> plot(x,y, main='cos(2x)', type='l', lty=1, bty='n')
$>$ y2 <- $\sin (2 * x)$
$>$ lines(x,y2, type='l', lty=2)
> same <- which( abs (y - y2) < 0.01)
$>$ points(x[same], y[same], pch=19, col='red', cex=3)
> legend('bottomleft', c("cos(2x)", "sin(2x)"), bty='n', lty=c(1,2))

par() outputs the (long) list of options that control plotting behaviour. Read ?par for all the details!
Common options to explore:

- mfrow, mfcol: multiple plots in figure
- mar, oma: margins around plot and figure.
- ask: whether to hit RETURN between pages of figures.


## Multiple plots in one figure

mfrow and mfcol are useful parameters within par(), but margins often need to be changed to maximise space.

```
pdf(file = 'mfrow_eg.pdf'
    width = 4, height = 6
par(mar = c(3.5, 3.5, 1.5, 0.5),
mgp =c(2.5, 1, 0))
    <- seq(from = 0, to = 2*pi,
len = 100)
plot(x, sin(x), main = "sin (x)",
&type = 'l'),
plot(x, sin (2*x), main = "sin (2x)"
plot(x, sin(3*x), main = "sin (3x)",
plot(x, cos(x), main = "cos (x)",
type = '1')
plot(x, cos(2*x), main = " cos (2x)",
type = '1')
plot(x, cos(3*x), main = "cos (3x)",
dev.off()
```


## Multiple data sources on one plot

When you wish to have multiple data sources on one plot (e.g. two time-series plots), the approach is to draw the first using plot and then draw subsequent features using lines or points. Axes are not rescaled, so draw the bigger plot first.
$>x<-1: 30$
$>y<-\operatorname{sqrt}(x) ; z<-\log (x)$
$>\operatorname{plot}(x, y)$; lines $(x, z, c o l=' r e d ')$
> plot(x,z, type $=$ "l", col = "red"); points(x,y) \#\# some data missing

122 / 191


## Saving your plots

R can save plots in many formats, including PDF, postscript, PNG, JPEG. Best to use vector formats (PDF, postscript) for graphs and bitmap formats (png, jpeg) for images.
$R$ has output devices, only one of which is active, dev.cur().

```
> dev.list()
> pdf(file='hist.pdf', width=7, height=7) ## inch
> dev.list()
> hist( rnorm(9999) )
> dev.off() ## close device
> png(file='hist.png', w=600, h=600) ## pixels
> hist( rnorm(9999) )
> dev.off()
```

Zoom in on text of PNG to see limitations of this format.

Next steps with plotting (Advanced)
$R$ has a vast range of functions for plotting particular data types. You may read about different packages for plotting:

- base graphics (or "traditional")
- lattice/grid (lattice is built upon grid)
- ggplot2 - http://had.co.nz/ggplot2/

ractor(cy)
$: 4$
$: 6$
08

Here are some starting points to explore:

- demo(graphics) to see diversity of plots.
- low-level functions: symbols(), rect(), segments(), abline().
- curve( $\mathrm{x} * \log (\mathrm{x})$ )
- R graphics gallery
http://gallery.r-enthusiasts.com/,
http://rgm2.lab.nig.ac.jp/RGM2/images.php

Outline
Introduction
Vectors
Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping
Vectorization
Random number generation
Debugging

- where am I currently? getwd()
- change me to a new directory: setwd("/tmp")
(GUls have chooser for interactively changing directory.)
-What files are in my [current] directory?

```
> dir()
> dir("/tmp")
> dir(pattern="\\.R$") ## regexps, see later.
```


## Scan, write, readLines

For basic reading/writing of data, use scan/write. Filenames are specified relative to current directory. Can even give URL as a file. Files often have a header which can be skipped over.

```
> x <- scan('files/ages.dat', skip=1)
summary(x)
Min. 1st Qu. Median Mean 3rd Qu. Max.
\[
\begin{array}{llllll}
14.00 & 19.25 & 27.25 & 31.58 & 39.25 & 65.00
\end{array}
\]
> s1 <- readLines('files/ages.dat') ## treats as strings
> summary(s1)
    Length Class Mode
        5 character character
```

> h <- scan('http://damtp.cam.ac.uk/user/sje30/r/heights.dat') \#\# heights.dat
$>$ summary $(\mathrm{h})$
Min. 1st Qu. Median Mean 3rd Qu. Max.
$\begin{array}{rrrrrr}136.2 & 158.3 & 164.9 & 163.0 & 169.2 & 178.7\end{array}$

Scan, write, readLines

```
> rand.vals <- round( runif(100, min=5, max=10), 2)
> tf <- tempfile()
> tf
[1] "/var/folders/rw/2g0_whns55x580hlgf1vjs8m0000gn/T//RtmpwUm0lm/fi
> write(rand.vals, tf)
> s <- scan(tf)
> all.equal(s, rand.vals)
[1] TRUE
```


## read.table et al.

If data are tabular, read.table or read.csv is often useful. (Useful for importing spreadsheets; just save as a comma separated value file, CSV.)

```
x <- read.table('./files/players.dat', sep = '\t',
    header = TRUE)
names(x)
head(x)
x[2,]
x$Goals
is.data.frame(x)
tf <- tempfile()
write.csv(x, tf, row.names=FALSE)
## sort by goals scored.
x[order(x$Goals, decreasing = TRUE), ]
```

See ?read.table.

## Rdata files

Text files are useful for portably storing data, so that they can be read across applications. R has its own format for efficiently storing objects. Files much smaller than text files. However, this format is not universally known.

```
n <- 99999; x <- rnorm(n)
```

n <- 99999; x <- rnorm(n)
txt.file <- tempfile()
txt.file <- tempfile()
> rda.file <- tempfile()
> rda.file <- tempfile()
> write(x, n, file = txt.file)
> write(x, n, file = txt.file)
save(x, n, file = rda.file)
save(x, n, file = rda.file)

## Compare sizes of files with the object.

## Compare sizes of files with the object.

object.size(x)
object.size(x)
file.info(txt.file)$size
file.info(txt.file)$size
file.info(rda.file)$size ## compressed
file.info(rda.file)$size \#\# compressed
rm(x,n)
rm(x,n)
load(rda.file) \#\# reload data.

```
load(rda.file) ## reload data.
```

Saving your workspace with .RData files
When you quit $R$, you are asked:
$>\mathrm{q}()$
Save workspace image? [y/n/c]:
If you answer y , all objects (variables and functions) in your global environment are saved for future use, using save.image. From ?save:
'save.image()' is just a short-cut for "save my current
workspace", i.e., 'save(list = ls(all=TRUE), file = ".RData")' It is also what happens with 'q("yes")'.
Warning: If an .RData file is present in the current directory when starting $R$, it is silently loaded. I think it can be dangerous, as you may not realise what values have been silently loaded. Better to instead be explicit:

```
> save.image(file='all.Rda') ## keep everything
> save(x, y, z, file='key.Rda') ## or just key objects
> ##
> load('all.Rda') ## reload objects
```


## Further I/O functions (Adraneces)

$R$ has many facilities for I/O. See for example the following help topics.

- ?connections - interface to files, pipes, sockets, compressed files ...
- ?sink - divert R output to a connectin
- ?dget / ?dput - read/write ASCII representation of an R object.

Also: XML, DBMS, SQLite, netCDF, hdf5, ...
or start R using R --no-restore

## Outline

ntroduction
Vectors
Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system

## Writing functions

Conditionals and looping
Vectorization
Random number generation
Debugging

Writing functions: overview

- Why bother?
- How to write (local args, return value; cannot change value)
- Example: computing std. deviation
- Local variables within functions
- Recursion.


## Functions

- Black-box approach; given inputs, what output should I expect? This requires good documentation of what your function does. Can it be described without having to look at the code?
- Finding the right level of definition for a function is hard, and how to modularise comes with experience. Typically rewrite many times before getting final solution
- How to define a new function:

```
> my.fun <- function(arg1, arg2) {
+ ## Doc string here.
+ x <- arg1 * 2
+ y <- sqrt(arg2) + 5
+ z <- x * y
+ ## last value is the return value of the function.
+ ## Use a list to return several items.
+ z ## same as return(x)
+ }
```


## Example of writing a new function

Compute the standard deviation of a vector of numbers:

$$
\text { std.dev }=\sqrt{\frac{\sum_{i=1}^{n}\left(x_{i}-\bar{x}\right)^{2}}{n-1}} \quad \text { where } \quad \bar{x}=\frac{\sum_{i=1}^{n} x_{i}}{n}
$$

```
std.dev <- function(x) {
+ ## Return std dev of X.
+ n <- length(x)
+ xbar <- sum(x)/n
+ diff <- x - xbar
+ sum.sq <- sum( diff^2)
+ var <- sum.sq / (n-1)
+ ## last value calculated is return value.
+ sqrt(var)
+ }
```

- Any change to formal args within a function does not change value of actual argument outside the function:

```
> sum.sq <- function(x) {
+ x <- x^2 ## change internally
+ sum(x)
+ }
> y <- c(4, 5, 6)
> sum.sq(y)
[1] 77
> y
[1] 4 5 6
```

- In std.dev, x is the name of a formal argument. In the following, y is called the actual argument (doesn't have to be named x - can be named however you wish)

```
> <- 5
> y <- c(9, 2, 7, 10)
> std.dev(y)
[1] 3.559026
> print(n) ## should still be 5, not 4
[1] 5
```

- Local variables within function are not available outside of function.


## Handling unbound variables

Variables created by assignment within a function are known as local variables (e.g. y below). If a variable is not local, or a formal argument, it is an unbound variable. It may be found in the enclosing environment (typically the global workspace), or an error is generated - this is bad practice!

```
> fn1 <- function(x) {
| y <- x^2
+ res <- sum( (y - thresh)^2 )
res
+ }
dat <- 1:5
> fn1(dat) ## case 1
Error in fn1(dat): object 'thresh' not found
> thresh <- 10
> fn1(dat) ## case 2
```

Handling unbound variables (2)
The
. . . arguments

In this case, better to define thresh as an argument of the function, and provide a default value:

```
> fn1 <- function(x, thresh=10) {
+ y<- x^2
+ res <- sum( (y - thresh)^2 )
+ res
+ }
> fn1(dat) ## case 3
```

[1] 379

Advanced: use codetools: :checkUsage() to find unbound vars; codetools: :findGlobals() for globals.

## The ... arguments

```
> p <- function(x, y, ...) {
+ if (diff(range(x)) > diff(range(y))) {
    plot(x, y, ...)
    } else {
        plot(y, x, ...)
    }
+ }
> p(rnorm(10, 0, 1), rnorm(10, 0, 5), main = "Expect swap",
+ col = "red")
> p(rnorm(10, 0, 5), rnorm(10, 0, 1), main = "No swap", pch = 19)
```



From R language definition
The . . . argument is special and can contain any number of arguments. It is generally used if the number of arguments is unknown or in cases where the arguments will be passed on to another function.
See ?cat for an example of number of arguments is unknown.

## Writing a replacement function (Adaneace)

Convention for a replacement function is that the name should end with <-. The last argument of the replacement function must be called value and is the RHS of the assignment.

```
> "threshold<-" <- function(x, value) {
    ## X is the object to update
    ## VALUE is the value on the RHS.
+ y <- ifelse(x > value, 1, 0)
return(y)
+ }
> x <- c(0.3, 0.1, 0.6,0.7, 0.9, 0.2)
> threshold(x) <- 0.4
> x
[1] 0 0 1 1 1 0
```

In general, replacement functions are used to set variable attributes or object slots.

- if
- switch
- for
- while
- Vectorization
- simple applications - numerics


## Tips for writing functions

- Can you think of a way to break down the problem so that a team can work on the problem, with each person assigned to a independent piece? "Divide + conquer".
- Each function should be easy to test, then you can "freeze" it. Write test cases, which can be automatically checked.

```
> all.equal(my.fun(100,200), 300)
```

- Rule of thumb: each function should be no more than a page or two of code.
- For large projects, avoid mixing computation and plotting in the same function - separate the two jobs; this makes it easier to run in batch.

```
> res <- some.computation(par1, par2, par3)
> plot.results(res)
```


## Outline

Introduction

## Vectors

Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping
Vectorization
Random number generation
Debugging
149 / 191

## Control-flow constructs

if / if ... else ...

```
x <- 8;
>
if (x > 10) {
+ ## condition was true
cat("x is bigger than 10\n")
} else {
cat("x is 10 or less\n")
+ }
x is 10 or less
```

Notes:
"else ..." can be omitted if you do not need it.
if returns a value, which can be assigned, e.g. y <- if ( $\mathrm{x}<10$ ) 40
else 20. A better solution in this case however is the vectorized form y <- ifelse( $\mathrm{x}<10,40,20$ )

## Braces in conditional constructs

Curly braces not needed if there is only one expression in the if clause:

```
if ( x > 10 ) {
```

```
if ( x > 10 )
```

if ( x > 10 )
y <- 1
y <- 1
y<-1 y \#

```

But braces are needed in multiline if/else statement:
```

if ( x > 10 ) {
y <- 1
} else {
y <- 0 \#\# OK
}

```
```

if ( x > 10 )
y<-1
else
y <- 0 \#\# NOT OK

## 

```

From ?Control: Note that it is a common mistake to forget to put braces ('\{ .. \}') around your statements, e.g., after 'if(..)' or 'for(....)'. In particular, you should not have a newline between '\}' and 'else' to avoid a syntax error in entering a 'if ... else' construct at the keyboard or via 'source'. For that reason, one (somewhat extreme) attitude of defensive programming is to always use braces, e.g., for 'if' clauses.

\section*{Switch (Advanced)}

Nested if ... else commands can get a bit messy. Like other languages, \(R\) has a switch construct. From ?switch:

\section*{Recursive functions}

Here is an example of using conditionals with a divide and conquer approach; quicksort in a few lines (albeit not very efficient).
+ ## Sort data into ascending order.
+ n <- length(data)
+ if (n <= 1) {
    data
+ } else {
        + pivot <- data[floor(n/2)]
+ less <- data[which(data < pivot)]
+ equal <- data[which(data == pivot)]
+ greater <- data[which(data > pivot)]
+ greater <- data[which(data > pivot)]
+ }
+ }
> all(replicate(99, {
+ data <- runif(2000, max=10)
+ all.equal(qsort(data), sort(data)) }))
```

```
```

> qsort <- function(data) {

```
```

```
> qsort <- function(data) {
```

[1] TRUE
[1] -2.801097
> centre(x, "median")

```
centre <- function(x, type) {
```

centre <- function(x, type) {
switch(type,
switch(type,
mean = mean(x),
mean = mean(x),
median = median(x),
median = median(x),
trimmed = mean(x, trim = .1))
trimmed = mean(x, trim = .1))
}
}
x <- rcauchy(10)
x <- rcauchy(10)
centre(x, "mean")
centre(x, "mean")
[1] 0.08003974
> centre(x, "trimmed")
[1] -0.05755348
centre(x, "median")
centre(x, "median")
[1] 0.08003974
centre(x, "trimmed")
[1] -0.05755348

```

\section*{Looping constructs}

Looping constructs allow you to repeat calculations as many times as you wish. This is why computers are so useful - it is just as easy (usually) to repeat something 1000 times as 10 times.
e.g. if you want to simulate flipping a (biased) coin 100 times, and counting the number of heads, no problem. If you want to repeat this process 1000 times, no problem. (See later.)

\section*{for (var in seq) command}
seq is a vector; var is set in turn to each value in the vector, and then command executed. Multiple commands can be given within braces. e.g.
```

> x <- 6
> for (i in 1:3) {

+ res <- x * i
+ cat(x, "*", i, "=", res, "\n")
+ }
6*1 = 6
6*2 = 12
6*3 = 18

```
while loops
```

while (condition) {
command
command
}

```

So the commands are executed until the condition is no longer true.
Typically then one of the commands will change the condition.
e.g. print all the Fibonacci numbers ( \(f[i]=f[i-1]+f[i-2]\) ) less than 100.
```

> n1 <- 0; n2 <- 1
> while (n2 < 100) {

+ print(n2)
+ old <- n2
+ n2 <- n2 + n1
+ n1 <- old
+ }

```

\section*{Breaking out of loops}
repeat expr will repeatedly execute expr until you break out of the loop.
```

> i <- 3
> repeat {

+ if (i == 5) {
+ break
+ } else {
+ cat("i is", i, "\n")
+ i<- i+1
}
+ }
i is 3
i is 4

```

\section*{Breaking out of loops}
next allows you to skip to next iteration of a loop. Both next and break can be used within other loops (while, for).
```

> for (i in 1:10) {

+ if ((i %% 2) == 0)
+ next
+ print(i)
+ }

```
[1] 1
[1] 3
[1] 5
[1] 7
[1] 9

\section*{A word on indentation}

Indentation helps you see the flow of the logic, rather than flattened version. (Use tab key to indent). Reformatting tools are available (e.g. within Emacs).
```


## version 1.

i <- 3
repeat {
if (i==10) {
break
} else {
cat("i is", i, "\n")
i<- i+1
}
}

```

Indentation helps to show structure, and match braces.

\section*{Vectorization}

When possible, operate on vectors, rather than using for loops.
Rewrite code, but beware sometimes not possible (Fibonacci).
Compute difference between times of events, e. Given \(n\) events, there will be \(n-1\) inter-event times. interval[i] <- e[i+1] - e[i]
> diff1 <- function(e) \{
\(+\mathrm{n}<-\) length \((\mathrm{e})\)
interval <- rep(0, n-1) \#\# good to pre-alloc!
for (i in 1:(n-1)) \{
\} interval[i] <- e[i+1] - e[i]
\(+\quad\) interval
+ + \(\quad\) diff2 <- function(e)
\(+\mathrm{n}<-\) length \((\mathrm{e})\)
\(+{ }_{+}^{+}\)
\(>\mathrm{e}<-\mathrm{c}(2,5,10.2,12,19)\)
> diff2(e)
[1] 3.05 .21 .87 .0
> all.equal(diff1(e), diff2(e))
[1] TRUE

\section*{Outline}
ntroduction

\author{
Vectors
}

Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping

\section*{Vectorization}

Random number generation
Debugging

\section*{Vectorization example}

Q: Flip a biased coin [ \(\mathrm{p}=0.6\) of heads] 100 times; how many heads do you get? Repeat this for 1000 trials.
```

> n <- 100 \#\# number of coin flips in trial
> p<-0.6 \#\# prob of getting heads
> ntrials <- 1000

```
> trial1 <- function(n, p.heads) {
```

> trial1 <- function(n, p.heads) {

+ count <- 0
+ count <- 0
for (i in 1:n) {
for (i in 1:n) {
if (runif(1) < p.heads)
if (runif(1) < p.heads)
count <- count +1
count <- count +1
    + }
    + }
count
count
+ }
+ }
>
>
> res <- rep(0, ntrials)
> res <- rep(0, ntrials)
> for (j in 1:ntrials) {
> for (j in 1:ntrials) {
+ res[j] <- trial1(n, p)
+ res[j] <- trial1(n, p)
+ }
+ }
> hist(res)

```
```

> hist(res)

```
```


## Vectorization example

Q: Flip a biased coin [ $\mathrm{p}=0.6$ of heads] 100 times; how many heads do you get? Repeat this for 1000 trials.

```
> n <- 100 ## number of coin flips in trial
> p <- 0.6 ## prob of getting heads
> ntrials <- 1000
```

> trial1 <- function(n, p.heads)

+ count <- 0
> trial2 <- function(n, p.heads) \{
for (i in 1:n) \{
if (runif(1) < p.heads)
$+\quad$ rand.vals <- runif(n)
$+\quad$ sum( rand.vals < p.heads)
if (runif(1) < p.heads)
$+\}$
$>$
count <- count +1
$+\quad\}$
count
$+\}$
> res <- replicate(ntrials,
$+\quad$ hist(res)
> res <- rep(0, ntrials)
$>$ for (j in 1:ntrials) \{
$+\quad$ res $[j]<-$ triall (n,
$+\operatorname{res}[j]<-\operatorname{triall}(n, p)$
$+\}$
> hist(res)


## apply family

e.g. how to compute sum of each row of a matrix? sum(A) will normally return the sum of all elements of $A$.
apply(X, MARGIN, FUN, ...)
MARGIN = 1 for row, 2 for cols.
FUN = function to apply
... = extra args to function.

```
> A <- matrix(1:6, 2,3)
> row.means <- apply(A, 1, mean) ## or rowMeans
> col.sums <- apply(A, 2, sum, na.rm=TRUE) ## colSums
```


## Vectorization example

Q: Flip a biased coin [ $p=0.6$ of heads] 100 times; how many heads do you get? Repeat this for 1000 trials.

```
> n <- 100 ## number of coin flips in trial
> p<- 0.6 ## prob of getting heads
> ntrials <- }100
```

> trial1 <- function(n, p.heads)
> trial2 <- function(n, p.heads) \{
$+\quad$ rand.vals <- runif(n)
count <- 0
for (i in 1:n)
$+\quad$ sum( rand.vals < p.heads)
$+\}$
$>$
if (runif(1) < p.heads) apply family

Other functions: lapply (apply to list), sapply (simplify), replicate, ...

```
> lapply(ls(), object.size)
> sapply(ls(), object.size)
> tapply(1:20, factor(rep(letters[1:5], each = 4)), sum)
> ## see also ?by
> mapply(sum, 1:5, 1:5, 1:5)
> hist( replicate(200, mean(rnorm(100)) ))
```

How to *apply - http://stackoverflow.com/questions/3505701

Exercise: Why is this better than writing a for loop?
(Eglen, 2009); parallel package: mclapply(), parLapply(),....

Sometimes you don't want to pollute name space by defining a new function, so just use an "anonymous function", i.e. a function without a name. Particularly useful e.g. in an apply call.

```
> my.mat <- matrix(1:10, ncol=5)
> apply(my.mat, 2, function(x) { sum(x^2) + 10 })
```

Since functions are just objects, anonymous functions are just objects without names, similar to 'anonymous numbers' like $a+b$ in an expression $a+b+c$.

## Efficiency

"premature optimization is the root of all evil" (Knuth, 1974).

Examples adopted from www.mathworks.com/res/code_segments $f 1$ is bad; should pre-allocate vector, rather than rely on $R$ to allocate memory repeatedly ${ }^{3}$.

```
> f1 <- function() {
+ n<- 1e4; decay <- 0.9995
+ out <- 1.0
+ for (i in 2:n)
+ for (i in 2:n)
    out
+ }
> system.time(o1 <- f1())
```

user system elapsed
$0.396 \quad 0.128 \quad 0.525$

```
> f2 <- function() {
```

> f2 <- function() {

+ n <- 1e4; decay <- 0.99995
+ n <- 1e4; decay <- 0.99995
+ out <- rep(0, n) \#\#pre-alloc
+ out <- rep(0, n) \#\#pre-alloc
+ out[1] <- 1.0
+ out[1] <- 1.0
+ for (i in 2:n)
+ for (i in 2:n)
out[i] <- out[i-1] * decay
out[i] <- out[i-1] * decay
out
out
+     + }
+     + }
> system.time(o2 <- f2())
> system.time(o2 <- f2())
user system elapsed
user system elapsed
0.036 0.001 0.037

```
    0.036 0.001 0.037
```

$0,1,1,2,3,5,8,13,21, \ldots$

$$
f[n]=f[n-1]+f[n-2]
$$

How to vectorize?
Exercise: write a function, fibonnaci(n) that returns the $n^{\text {th }}$ element of the sequence. Assume that fibonnaci(1) $==0$, $f$ ibonacci(2) $==1$. Exercise: use fibonacci() to estimate the golden ratio.

## A final word on efficiency

- Rule 1 of optimization: don't bother (Kernighan).
- For loops are not always a bad thing. See last example in Help Desk article (May 2008).
http://cran.r-project.org/doc/Rnews/Rnews_2008-1.pdf

[^1]
## Numerics issues

Although integer arithmic is reliable, floating-point arithmetic is to be treated with care! (All $R$ 's calculations are in what $C$ programmers call "double precision".) Compare:
> $1.0+2.0=3.0$
[1] True
$>0.1+0.2=0.3$
[1] False
From FAQ (7.31?)
> a <- sqrt(2)
> a * a == 2
[1] FALSE
>a*a-2
[1] $4.440892 \mathrm{e}-16$

How big is infinity?
Use while loop to estimate it:
$>x<-1$
$>$
$>$
while (is.finite(x*2) )
${ }_{>}^{+}$
> x \#\# 8.988466e+307
[1] $8.988466 \mathrm{e}+307$
> x*2 \#\# Inf
[1] Inf
> (x*2)/2 \#\# Inf
[1] $\operatorname{Inf}$
> .Machine\$double.xmax
[1] $1.797693 \mathrm{e}+308$

## Solution - testing near equality

```
> all.equal( 0.1 + 0.2, 0.3)
```

[1] TRUE

How small is epsilon?
How big can $\epsilon$ be such that $1+\epsilon=1$ ? (Taken from Goldberg (1991) ACM article, p220)

```
> eps <- 1
* whie (eps + 1 > 1)
+ + 'e
eps ##1.110223e-16
```

[1] 1.110223e-16
$>1+$ eps \#\# 1
[1] 1
$>(1+$ eps $==1)$ \#\# TRUE
[1] TRUE
$>1+(2 * e p s) \# \# 1$
[1] 1
> $(1+(2 * e p s)==1)$ \#\# FALSE

## Outline

introduction

## Vectors

Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping

## Vectorization

Random number generation

## Debugging

## Random number generation

Computers usaully generate "pseudo-random numbers". They are generated based on some iterative formula:

$$
x_{\text {new }}=f\left(x_{\text {old }}\right) \quad \bmod N
$$

where modulo operation provides the "remainder" division. To generate the first random number, you need a seed. Setting the seed allows you to reliably generate the same sequence of numbers, which can be useful when debugging programs.
$R$ has many routines for generating random samples from various distributions, but for now we will just use runif(), (and maybe rnorm()). Exercise: write a random number generator. See: "Randu: a bad random number generator". http://physics.ucsc.edu/~peter/115/randu.pdf Exercise: Apply the central limit theorem to generate samples from a normal distribution by adding together samples from a uniform distribution.

## Outline

Introduction
Vectors
Calling functions
Scripts
Matrices
Boolean logic
Lists
Factors
Character arrays
Objects in your environment
Basic plotting
Reading/writing data to file system
Writing functions
Conditionals and looping
Vectorization
Random number generation

## Warnings and errors

- A warning is softer than an error; if a warning is generated your program will still continue, whereas an error will stop the program.
$>\log (c(2,1,0,-1,2))$
Warning in $\log (c(2,1,0,-1,2)):$ NaNs produced
[1] 0.69314720 .0000000 -Inf NaN 0.6931472
> xor( c(TRUE, FALSE))
Error in xor(c(TRUE, FALSE)): argument "y" is missing, with no default
- If you try to isolate warnings, you can change warnings to errors: options (warn=2). See ?options for further details.
- Add warnings and errors to your code using warning(), stop().
- Can add "assertions" into your code to check that certain values hold.
$>$ stopifnot $(x>0)$
- Other useful safety checks: all ( $x>0$ ), any ( $x>0$ )


## Single-stepping through your code

Use browser() to single-step through your code. Place it within your function at the point you want to examine (e.g.) local variables.

Can use debug (function.name) to step through entire function. undebug() will remove that debug call.
Within the browser, you can enter expressions as normal, or you can give a few debug commands:

- n : single-step
- c: exit browser and continue
- Q: exit browser and abort, return to top-level.
- where: show stack trace.

Debug on stddev.R

## Traceback

When your program generates an error, use traceback() to find out where it went wrong:

```
> start <- function() { go( sqrt(10)) }
go <- function(x) { inner(x, '-13')}
inner <- function(a, b) {
c <- sqrt(b)
+ a * log(c)
+ }
start() ## error
traceback() ## postmortem debugging
```

Here's a possible usage of browser() that I have in my code:
find. high <- function ( $x, t$ ) \{
\#\# Return samples in $x$ bigger than $t$
\#\# (Better to use $x[x>t]$ in real-life!)
max.length <- 100 \#\# should be upper limit...
results $<-$ rep ( 0 , max.length)
counter <- 0
for ( i in x )
if $(i>t)\{$
counter <- counter + 1
if (counter $>$ max.length ) \{
browser()
\} else \{
results [counter] $<-$ i
\}
\} $\}$
\}
results [1: counter]
\}
$x<-\operatorname{rnorm}(100)$
find.high (x, 0.7)
$x<-$ rnorm (1000)
(1- pnorm (0.7)) * length $(x) \# \#$ expected.
find.high (x, 0.7)
recover() is like browser(), except you can choose which level to inspect, rather than the level at which browser was called.
Following allows recover() to be launched when you hit an error:
options(error=recover)
Here we simply tell $R$ that when an error is generated, we call the function "recover". The default is NULL, in which case stop is called.

## From ?options:

Note that these need to specified as e.g.
options(error=utils::recover) in startup files such as .Rprofile.

## CRAN: Comprehensive R Archive Network

CRAN: Site(s) for downloading R, and also its many contributed *packages*.
Mac/Win have a GUI for installing packages, or it can be done on the command line:

```
install.packages(c("splancs", "sp"))
R CMD INSTALL mypackage.tar.gz ## from shell
```

If asked to selected a CRAN mirror, in UK use:
http://www.stats.bris.ac.uk/R.
For Bioconductor

```
> source("http://www.bioconductor.org/biocLite.R")
> library("BiocInstaller")
> biocLite("Biobase")
```

- R has a packaging system for external code.
- A package is loaded from a library using library("pkg.name").
- Beware: don't call a package a library! A library is a group of folders where packages are stored...


## Packages

```
> library()
```

> library()
library(help="cluster") \#\# what's in this?
library(help="cluster") \#\# what's in this?
> library("cluster") \#\# load package
> library("cluster") \#\# load package
> example(pam) \#\# can use pam and friends.
> example(pam) \#\# can use pam and friends.
> detach("package:cluster") \#\# remove pkg.
> detach("package:cluster") \#\# remove pkg.

## view available packages

```
## view available packages
```


## Managing libraries

If you do not have write access to the default library, $R$ will create a local one in your home directory. Multiple libraries are supported:

```
> .libPaths()
[1] "/Users/stephen/NOBACKUP/RLIB"
[2] "/Library/Frameworks/R.framework/Versions/3.2/Resources/li
```

You can also set $R$ libraries as a global shell environment in your . bashrc file (e.g. PWF linux)
export R_LIBS=\$HOME/NOBACKUP/RLIB
(Be careful! Check that you are not overwriting an existing R_LIBS setting.)

A success story of R. Started 2001 with aims to:

- provide access to stat/graphical methods for analysis of genomic data.
- link seamlessly to on-line databases (PubMed/GenBank).
- allow rapid development of extensible software.
- provide training in methods (short courses).
- promote software with high quality docs and reproducible research (vignettes) ...
- Gentleman et al. (2004) Genome Biology 5:R80. http://genomebiology.com/2004/5/10/R80
- S3, S4 and S4 Reference classes for OOP.
- Building your own packages. Useful for packaging up your code, data sets and documentation. You may wish to do this for large projects that you wish to share with others. Read Writing $R$ Extensions manual and see package.skeleton to get started.
- Access to databases. Computational Biology datasets are often quite large, and you might wish to access data via databases. R package DBI provides common interface to SQLite, MySQL, Oracle. See Gentleman (2008), Chapter 8.

Reproducible research: Sweave, knitr and vignettes

- Use one file to store code and document. Best shown by way of example... estimate.Rnw
- Vignettes often used in Bioconductor to document packages.
- Interactively explore vignettes.

```
```

> library("tkWidgets")

```
```

```
```

> library("tkWidgets")

```
```

> vExplorer()

- knitr is the natural successor to Sweave. It rocks!


[^0]:    ${ }^{1}$ tentative

[^1]:    ${ }^{3}$ This is not optimisation, but rather good coding practice.

