

10. Interfacing R

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With Bioconductor, R can do a **huge** proportion of the analyses you'll want – but not everything

- Intensive (or anachronistic) C++, FORTRAN work, e.g. for pedigrees
- 'Speciality' analyses; some need different computing architecture
- Fancy interactive graphics

R can be used to 'manage' other software. Today we'll illustrate some favorite examples

Starting other software

NB these commands are for Windows only; see help files for e.g. Unix versions

- system is the equivalent of a DOS-style command
- system("notepad") starts the Notepad editor
- If the command takes arguments, put them in the same string; system("notepad myfile.txt")

The shell command does much the same thing.

Some more options for system;

- wait; R 'hangs' until completion
- minimized; new program only appears in TaskBar
- show.output.on.console

Paths for files can be a little messy; system starts in your working directory (getwd). Outside of this, give the full pathway.

paste is useful, if you need to do a lot of this sort of thing. source may also help

Examples

Code for a really mundane job;

```
for(i in 1:100){
     infile <- paste("gene",i,"data.txt", sep="")
     outfile <- paste("gene",i,"phase.out", sep="")
     system(paste("PHASE",infile,outfile))
}
```
... this will churn away for hours, although with no error-control.

Why did we use wait=TRUE here? (the default)

Examples

- WinBUGS implements Bayesian analyse; it's not super-fast but is very flexible
- It needs special (& clever) architecture to achieve this
- WinBUGS' input, output, graphics are all rather clunky
- R is better; so R2WinBUGS calls WinBUGS for the difficult bits, and does all the 'translation' itself
- This is done with (repeated) use of system()

Outline

Many programs already exist to do useful analyses. It is more convenient to call them from R than to rewrite them in R.

Sometimes this involves calling the C code directly, sometimes just involves using R to write input files for another program

Examples:

- Graphviz: drawing networks
- PMF: input files for ancient Fortran software
- Google Earth: displaying outliers in context.

GraphViz (<http://www.graphviz.org>) is a free program for drawing networks, written by AT&T researchers.

Its input format looks like

```
"15" [shape= box,regular=1 ,height= 0.5 ,width= 0.75 ,style=filled,color= grey ] ;
"16" [shape= circle ,height= 0.5 ,width= 0.75 ,style=filled,color= grey ] ;
"2x3" [shape=diamond,style=filled,label="",height=.1,width=.1] ;
"2" -> "2x3" [dir=none, weight=1] ;
"3" -> "2x3" [dir=none,weight=1] ;
"2x3" -> "1" [dir=none, weight=2] ;
"2x3" -> "4" [dir=none, weight=2] ;
"2x3" -> "5" [dir=none,weight=2] ;
"2x3" -> "6" [dir=none, weight=2] ;
```
The sem package uses GraphViz to display path diagrams for structural equation models and the gap package uses it to draw pedigrees.

Drawing networks

In gap the pedtodot() function writes a GraphViz input file from a pedigree in GAS or LINKAGE format.

First the code prints nodes for each individual, with sex and affectedness information

```
for (s in 1:n) cat(paste("\"", id.j[s], "\" [shape=",
    sep = ""), shape.j[s], ",height=", height, ",width=",
   width, ",style=filled,color=", shade.j[s], "] ;\n")
```
giving output like

"16" [shape= circle ,height= 0.5 ,width= 0.75 ,style=filled,color= grey] ;

It then works out all the matings and creates small nodes for each mating and lines connecting the parents to these nodes

```
mating \leq paste("\"", s1, "x", s2, "\"", sep = "")
cat(mating, "[shape=diamond,style=filled,label=\"\",height=.1,width=.1] ;\n")
cat(paste("\\"1, s1, "\\"1, sep = ""); " -> ", mating,paste(" [dir=", dir, ",weight=1]", sep = ""),
     " : \n\langle n" \ranglecat(paste("\\"1", s2, "\\"1", sep = ""), " -> ", mating,
    paste(" [dir=", dir, ",weight=1]", sep = ""),
    " ; \n")
```
Drawing networks

giving output like

```
"2x3" [shape=diamond,style=filled,label="",height=.1,width=.1] ;
"2" -> "2x3" [dir=none,weight=1] ;
"3" -> "2x3" [dir=none,weight=1] ;
```
and then connects children to parents.

Drawing networks

[Bioconductor also has GraphViz more integrated with R in the RGraphViz package]

PMF: factor analysis

PMF is a program for constrained factor analysis in analytic chemistry. It is controlled by an ugly text input file:

```
pmfini<-c(" ##PMF2 .ini file for: Simulations from R",
" ## Monitor code M: if M>1, PMF2 writes output every Mth step",
" ## For finding errors, use M<1 to output debug information",
" ## M PMF2 version number",
\frac{1}{4.2}",
" ## Dimensions: Rows, Columns, Factors. Number of \"Repeats\"",
" @nt@ @ns@ @sources@ 1",
" ## \"FPEAK\" (>0.0 for large values and zeroes on F side)",
" @FPEAK@",
" ## Mode(T:robust, F:non-robust) Outlier-distance (T=True F=False)",
" @isrobust@ @outlier@",
" ## Codes C1 C2 C3 for X_std-dev, Errormodel EM=[-10 ... -14]",
" \t 0.0100 \t 0.0000 \t 0.0100 \t -12"." ## G Background fit: Components Pullup_strength",
\frac{1}{10} 0.0000",
" ## Pseudorandom numbers: Seed Initially skipped",
" and " @seed@ \qquad \qquad 0"," ## Iteration control table for 3 levels of limit repulsion \"lims\"",
```
The @value@ are places where we want to substitute in a value.

PMF: factor analysis

R code for the substitutions looks like

```
temp<-gsub("@FPEAK@",formatC(fpeak,digits=4,format="f"),pmfini)
temp<-gsub("@isrobust@",isrobust,temp)
seed<-as.character(as.integer(seed))
temp<-gsub("@seed@", seed, temp)
```
We can write data files needed by PMF, and then write the control file, then call PMF with the system() function. After PMF finishes we read in the results.

```
write.table(cX,file=xfile, quote=FALSE, col.names=FALSE, row.names=FALSE)
write.table(cU,file=efile, quote=FALSE, col.names=FALSE, row.names=FALSE)
if (!debug)
    on.exit(unlink(c(xfile,efile,inifile)))
writeLines(temp,inifile)
```

```
sysval<-system(paste(pmf,inifile), intern=TRUE,invisible=!debug)
```
ffactor<-read.table(outfiles\$f,row.names=sourcenames,col.names=species) gfactor<-read.table(outfiles\$g,row.names=times,col.names=sourcenames)

From the user's viewpoint it looks as though everything was done in R.

SVG (Scalable Vector Graphics) is a non-bitmap graphics format for the web.

The RSvgDevice and RSVGTipsDevice packages allow R output to SVG format.

We can use this to create graphs with links and tooltips. For example, a funnelplot showing associations between a large number of SNPs and VTE.

Point at a dot to see the SNP it represents, and click to go to information about the gene.

SVG+tooltips

```
for(i in 1:length(or)) {
     setSVGShapeToolTip(title=gene[i],
          desc1=snp[i],
          desc2=if(abs(lor[i]/se[i])>qnorm(0.5/n,lower.tail=FALSE))
                                  qvals[i] else NULL
          )
      setSVGShapeURL(paste("http://pga.gs.washington.edu/data",
                                            tolower(gene[i]),
                                             sep="1")
      points(prec[i],lor[i], cex=1, pch=19, col='grey')
  }
```
Google Earth is controlled by KML files specifying locations. KML is another plain text format.

```
We can write a KML file
<?xml version="1.0" encoding="UTF-8"?>
         <kml xmlns="http://earth.google.com/kml/2.1">
         <Placemark>
         \langlename\rangle 1 \langle/name\rangle<Point> <coordinates>-118.0256,34.11619,400</coordinates>
  </Point>
 </Placemark>
\langle/kml\rangle
```
and then send it to Google Earth with the shell.exec(filename) function, which opens a file using whatever is the appropriate program.

Google Earth

The identify() function lets the user select a point on a scatterplot.

In this example the points are locations where air pollution was measured, and we can call Google Earth to look at the location.