## R PROFILING AND OPTIMISATION

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## PENDING CHANGES

Warning: this is under construction.

This vignette contains experimental material which may sink down to the package implementation, or vanish.

## Known issues:

- Control information may be included as special stack in raw format.
- A list of profiles may become default. Only one profiling interval value per profile.
- Nodes may be implemented as factor. Work-around for the R factor handling needs to be added, i.e. factor as a data structure.
- changing timing interval is too expensive, as rle is not transparent to data frames. Implement profiles as a list, with a time interval attribute per list element.

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An R vignette for package sprof.

URL: http://sintro.r-forge.r-project.org/

Private Version

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## Profiling facilities in R

For the impatient: table 1 on page 9 and table 2 on page 9 give you information from profiling 100 runs of a simple linear regression. You get another view of the same information by fig. 15 on page 47. If you want to know more, please have some patience.

R provides the basic instruments for profiling, both for time based samplers as for event based instrumentation. Information on R profiling is in section 3.2 "Profiling R code for speed" and section 3.3. "Profiling R code for memory use" of "Writing R Extensions" http://cran.r-project.org/doc/manuals/R-exts.html. Specific information on memory profiling is in

http://developer.r-project.org/memory-profiling.html.

However this source of information seems to be rarely used.

Maybe the supporting tools are not adequate. The summaries provided by R reduce the information beyond necessity. Additional packages are available, but these are not sufficiently action oriented.

With package **sprof** we want to give a data representation that keeps the full profile information. Tools to answer common questions are provided. The data structure should make it easy to extend the tools as required.

The package is currently distributed at r-forge as part of the sintro material.

```
To install this package directly within R, type install.packages("sprof",repos="http://r-forge.r-project.org")
```

To install the recent package from source directly within R, type install.packages("sprof",repos="http://r-forge.r-project.org",type="source")

## L<sup>A</sup>T<sub>E</sub>X LAYOUT TOOLS AND R SETTINGS

You may want to skip this section, unless you want to modify the vignette for your own purposes, or look at the internals.

This is the main library we are going to use.

```
_ Input _
library(sprof)
 search()
                            _____ Output ____
"package:Rgraphviz"
[1] ".GlobalEnv"
 [3] "package:graph"
                             "package:sna"
 [5] "package:grid"
                             "package:wordcloud"
 [7] "package:RColorBrewer" "package:Rcpp"
[9] "package:xtable"
                             "package:sprof"
[11] "package:stats"
                             "package:graphics"
[13] "package:grDevices"
                             "package:utils"
[15] "package:datasets"
                             "package:methods"
[17] "Autoloads"
                             "package:base"
```

We want immediate warnings, if necessary. Set to level 2 to handle warnings as error.

We want a second chance on errors. So we install an error handler.

```
options(error = recover)

Input ______
```

Print parameters used here:

```
options(width = 72)
options(digits = 6)
```

For *str* output, we generally use these settings:

```
strx <- function(x,
    max.level=2, vec.len=3,
    nchar.max=40,
    list.len=12,
    width=70, strict.width="wrap",...)
{
    cat(paste("##strx:",deparse(substitute(x)),"\n"))</pre>
```

<u>ToDo:</u> add keep3 to keep header, some middle, tail

For larger tables and data frames, we use a kludge to avoid long outputs.

```
Input
xcutdata.frame <- function(df, cut, margin=5)
{
#! keep3, to add: margin top - random center - margin bottom
        if (!is.data.frame(df)) return(df)
        nrow <- nrow(df)</pre>
        # cut a range if it is not empty.
        # Quiet noop else.
        # Does not cut single lines.
         cutrng <- function(cutfrom, cutto){</pre>
                  if (cutfrom<cutto){</pre>
                 df[cutfrom,] <- NA</pre>
                 if (!is.null(rownames(df))) rownames(df)[cutfrom] <- "< cut >"
                 if (!is.null(df$name)) df$name[cutfrom] <- ""</pre>
                 cutfrom <- cutfrom+1</pre>
                 df[-(cutfrom:cutto),]
                 }#if
        if (!missing(cut)) {df <- cutrng(cut[1],cut[2]); return(df)}</pre>
        if (!missing(margin)) { # changed. Set to default 5.
                 if (length(margin)==1) margin <- c(margin,margin)</pre>
        cut <- c(margin[1]+1,nrow-margin[2])</pre>
        df <-cutrng(cut[1],cut[2]);</pre>
        return(df)}
         if (!missing(keep3)) { cut <- c(keep3[1]+1, keep3[1]+1,
                  nrow-keep3[3]-1,nrow-keep3[3]-1)
         if (cut[3]-cut[4] > keep3[2]+2){delta<-(cut[3]-cut[2]) div 2}
         cut[3]<-0
         browser()
         } else df <- cutrng(cut[1],cut[4])</pre>
         cutrng(cut[1],cut[4]) return(df)}
```

We use the R function xtable() for output and LATEX longtable. A convenient wrapper to use this in our Sweave source is given here. Among others, it adds zero.print.

ToDo: remove text vdots from string/-name columns. Note: this may be a factor. Use empty string.

\_ Input \_

```
library(xtable)
prxt <- function (x, digits=2, cut=TRUE, caption=NULL,</pre>
        label=NULL, zero.print=NULL, print.results=TRUE,...) {
        if (cut) {margin <- 10
        if (nrow(x)> 2*margin+3) x <-xcutdata.frame(x, margin=margin)}</pre>
        pr <- print(</pre>
                xtable(x, digits=digits, caption=caption,
                                 label=label, ...),
                floating=FALSE,
                tabular.environment="longtable",
                caption.placement="top",
                zero.print = ".",
                NA.string="\\vdots",
                print.results=FALSE)
#
                 NA.string="", #NA.string="\\vdots",
                pr <- gsub( "$\backslash$vdots","\vdots",x=pr, fixed=TRUE)</pre>
                if(!is.null(zero.print))
                         pr <- gsub( " 0 ",zero.print, x=pr, fixed=TRUE)</pre>
                if (print.results) cat(pr)
                invisible(pr)
        }
```

This is to be used with <<pre>rint=FALSE, results =tex, label=tab:prxx>>=

The graph visualisation family is not friendly. We try to get control by using a wrapper which is at least used to the members of the *graphviz* clan. This will be used in later sections.

```
}
#"lightyellow",
```

#### 1. Profiling

The basic information provided by all profilers is a protocol of sampled stacks. For each recorded event, the protocol has one record, such as a line with a text string showing the sampled stack.

We use profiles to provide hints on the dynamic behaviour of programs. Most often, this is used to improve or even optimise programs. Sometimes, it is even used to understand some algorithm.

Profiles represent the program flow, which is considered to be laid out by the control structure of a program. The control structure is represented by the control graph, and this leads to the common approach to (re)construct the control graph, map the profile to this graph, and used graph based methods for further analysis. The prime example for this strategy is the GNU profiler <code>gprof</code> (see http://sourceware.org/binutils/docs/gprof/) which is used as master plan for many common profilers.

It is only half of the truth that the control graph can serve as a base for the profiled stacks. In R, we have some peculiarities.

lazy evaluation: Arguments to functions can be passed as promises. These are only evaluated when needed, which may be at a later time, and may then lead to insertions in the stack. So we may have information resulting from the data flow, interspersed with the control flow.

**memory management:** Allocation of memory, and garbage collection, may interfere and leave their traces in the stack. While allocation is closely related to the visible control flow, garbage collection is a collective effect largely out of control of the code to execute.

**primitives:** Internal functions may escape the usual stack conventions and execute without leaving any identifiable trace on the stack.

control structures: In R, many control structures are implemented as function. Most notably, the <code>apply()</code> family appears as function calls and can lead to cliques in the graph representation that do not correspond to relevant structures. Since these functions are well know, they can have a special treatment.

So while the stack follows an overall well known dynamics, in R there are exceptions from regularity.

The general approach, by <code>summaryRprof()</code> and others, is to reduce the profile to node information, or to consider single transitions.

We take a different approach. We take the stacks, as recorded in the profiles as our basic information unit. From this, we ask: what are the actions we need to answer our questions? Representation in graphs may come later, if they can help.

If the stacks would come from the control flow only, we could make use of the sequential nature of stacks. But since we have to live with the R specific interferences, we stay with the raw stacks.

ToDo: stacks? order?

Do: rearrange cks? detect

In this presentation, we will use a small list of examples. Since <code>Rprof</code> is not implemented on all systems, and since the profiles tend to get very large, we use some prepared examples that are frozen in this vignette and not included in the distribution, but all the code to generate the examples is provided.

## 1.1. Simple regression example.

```
n <- 10000
x <- runif(n)
err <- rnorm(n)
y <- 2+ 3 * x + err
reg0data <- data.frame(x=x, y=y, err=err)
rm(x,y,err)</pre>
```

We will use this example to illustrate the basics. Of course the immediate questions are the variance between varying samples, and the influence of the sample size n. We keep everything fixed, so the only issue for now is the computational performance under strict iid conditions.

Still we have parameters to choose. We can determine the profiling granularity by setting the timing interval, and we can use repeated measurements to increase precision below the timing interval.

The timing interval should depend on the clock speed. Using for example 1ms amounts to some 1000 steps on a current CPU, per kernel.

If we use repeated samples, the usual rules of statistics applies. So taking 100 runs and taking the mean reduces the standard deviation by a factor 1/10.

By the usual R conventions, seconds are used as time base for parameters. However report will use ms as a time base.

Here is an example how to take a profile, using basic R. See section 1.1.2 on page 10 how to use **sampleRprof** in package **sprof** for an easier solution.

```
profinterval <- 0.001

simruns <- 100

Rprof(filename="RprofsRegressionExpl.out", interval = profinterval)

for (i in 1:simruns) xxx<- summary(lm(y~x, data=reg0data))

Rprof(NULL)
```

We now have the profile data in a file RprofsRegressionExpl.out. For this vignette, we use a frozen version RprofsRegressionExpl01.out.

## 1.1.1. R basic. The basic R functions invite us to get a summary.

ToDo: Can we calibrate times to CPU rate? Introduce cpu clock cycle as a time base

sumRprofRegressionExpl <- summaryRprof("RprofsRegressionExpl01.out")
#str(profile\_nodes\_rle, max.level=2, vec.len=3, nchar.max=40, list.len=6)
strx(sumRprofRegressionExpl)</pre>

```
Output -
##strx: sumRprofRegressionExpl
List of 4
$ by.self :'data.frame': 41 obs. of 4 variables:
..$ self.time : num [1:41] 0.087 0.057 0.051 0.043 0.042 0.04 0.032
..$ self.pct : num [1:41] 16.67 10.92 9.77 8.24 ...
..$ total.time: num [1:41] 0.113 0.099 0.069 0.043 0.474 0.045 0.033
   0.114 ...
..$ total.pct : num [1:41] 21.65 18.97 13.22 8.24 ...
$ by.total :'data.frame': 62 obs. of 4 variables:
..$ total.time: num [1:62] 0.522 0.522 0.521 0.521 0.521 0.521 0.521
   0.521 ...
..$ total.pct : num [1:62] 100 100 99.8 99.8 ...
..$ self.time : num [1:62] 0.006 0 0.001 0 0 0 0 0 ...
..$ self.pct : num [1:62] 1.15 0 0.19 0 0 0 0 0 ...
$ sample.interval: num 0.001
$ sampling.time : num 0.522
```

The summary reduces the information contained in the profile to marginal statistics per node. This is provided in two data frames giving the same information, only in different order.

The file contains several spurious recordings: nodes that have been recorded only few times. It is worth noting these, but then they better be discarded. We use a time limit of 4ms, which given our sampling interval of 1ms means we require more than four observations.

Table 1: summaryRprof result: by.self as final stack entry, all records

	self.time	self.pct	total.time	total.pct
"lm.fit"	0.09	16.67	0.11	21.65
"[.data.frame"	0.06	10.92	0.10	18.97
"model.matrix.default"	0.05	9.77	0.07	13.22
"as.character"	0.04	8.24	0.04	8.24
"lm"	0.04	8.05	0.47	90.80
"summary.lm"	0.04	7.66	0.04	8.62
"structure"	0.03	6.13	0.03	6.32
"na.omit.data.frame"	0.03	4.98	0.11	21.84
"anyDuplicated.default"	0.02	4.21	0.02	4.21
"as.list.data.frame"	0.02	4.21	0.02	4.21
< cut >	:	:	:	:

"FUN"	0.00	0.19	0.01	1.34
"%in%"	0.00	0.19	0.00	0.77
"deparse"	0.00	0.19	0.00	0.38
" <b>\$</b> "	0.00	0.19	0.00	0.19
"as.list.default"	0.00	0.19	0.00	0.19
"as.name"	0.00	0.19	0.00	0.19
"coef"	0.00	0.19	0.00	0.19
"file"	0.00	0.19	0.00	0.19
"NCOL"	0.00	0.19	0.00	0.19
"terms.formula"	0.00	0.19	0.00	0.19

Table 2: summary R<br/>prof result: by.total, total time  $>4\mathrm{ms}$ 

	total.time	total.pct	self.time	self.pct
" <anonymous>"</anonymous>	0.52	100.00	0.01	1.15
"Sweave"	0.52	100.00	0.00	0.00
"eval"	0.52	99.81	0.00	0.19
"doTryCatch"	0.52	99.81	0.00	0.00
"evalFunc"	0.52	99.81	0.00	0.00
"try"	0.52	99.81	0.00	0.00
"tryCatch"	0.52	99.81	0.00	0.00
"tryCatchList"	0.52	99.81	0.00	0.00
"tryCatchOne"	0.52	99.81	0.00	0.00
"withVisible"	0.52	99.81	0.00	0.00
< cut >	:	:	:	:
"as.list"	0.02	4.41	0.00	0.00
"anyDuplicated.default"	0.02	4.21	0.02	4.21
"as.list.data.frame"	0.02	4.21	0.02	4.21
"sapply"	0.01	2.68	0.00	0.19
"match"	0.01	2.11	0.00	0.19
"[[.data.frame"	0.01	1.53	0.00	0.19
"[["	0.01	1.53	0.00	0.00
"rep.int"	0.01	1.34	0.01	1.34
"FUN"	0.01	1.34	0.00	0.19
"list"	0.01	0.96	0.01	0.96

1.1.2. Package sprof. In contrast to the common R packages, in the **sprof** implementation we take a two step approach. First we read in the profile file to an internal representation. Analysis is done in later steps.

```
______ Input ______
sprof01<- readRprof("RprofsRegressionExpl01.out")
```

The data contain identification information for reference. This will be used in the functions of **sprof** and shown in the displays. Here is the summary of this section:

```
_ Input _
str(sprof01$info)
                                   Output
'data.frame':
                    1 obs. of 9 variables:
                 : Factor w/ 1 level "RprofsRegressionExpl01.out 2013-06-13 23:46:04": 1
s id
                 : POSIXct, format: "2013-08-15 22:02:43"
$ date
$ nrnodes
                 : int 62
$ nrstacks
                 : int 50
$ nrrecords
                 : int 522
$ sample.interval: num 0.001
$ sampling.time : num 0.522
                  : Factor w/ 1 level "sample.interval=1000": 1
$ ctllines
$ ctllinenr
                  : num 1
```

For this vignette, we change the id information. So in this context:

```
sprof01$info$id <- "sprof01"
```

We keep this example and use the copy sprof01 of it extensively for illustration.

```
save(sprof01, file="sprof011m.RData")
```

To run the vignette with a different profile, replace **sprof01** by your example. You still have the file for reference.

Package sprof provides a function sampleRprof() to take a sample and create a profile on the fly, as in

```
sprof01temp <- sampleRprof(runif(10000), runs=100)
```

The basic data structure consists of four data frames. The <code>info</code> section collects global information from the input file, such as an identification strings and various global matrix. The <code>nodes</code> section initially gives the same information marginal information as <code>summaryRprof</code>. The <code>stacks</code> section puts the node information into their calling context as found in the input profile file. The <code>profiles</code> section gives the temporal context. It is implemented as a list, but conceptually it is a data frame. Implementing it as a list allows run length encoding of variables, which unfortunately is not allowed by R in data frames.

ToDo: add sampling.interval, sampling.time for backward compatibility

```
strx(sprof01) Input _____
```

```
Output -
##strx: sprof01
List of 4
$ info :'data.frame': 1 obs. of 9 variables:
..$ id : chr "sprof01"
..$ date : POSIXct[1:1], format: "2013-08-15 22:02:43"
..$ nrnodes : int 62
..$ nrstacks : int 50
..$ nrrecords : int 522
..$ sample.interval: num 0.001
..$ sampling.time : num 0.522
..$ ctllines : Factor w/ 1 level "sample.interval=1000": 1
..$ ctllinenr : num 1
$ nodes :'data.frame': 62 obs. of 5 variables:
..$ name : Factor w/ 62 levels "!","..getNamespace",..: 1 2 3 4 5 6 7
  8 ...
..$ self.time : num [1:62] 2 0 2 0 0 57 0 1 ...
..$ self.pct : num [1:62] 0.38 0 0.38 0 ...
..$ total.time: num [1:62] 2 1 4 26 99 99 8 8 ...
..$ total.pct : num [1:62] 0.03 0.01 0.05 0.34 1.29 1.29 0.1 0.1 ...
$ stacks :'data.frame': 50 obs. of 7 variables:
..$ nodes :List of 50
.. .. [list output truncated]
..$ shortname : Factor w/ 50 levels
   "S<A>eFttCtCLtCOdTCwVeesleem.m..n.n...[["| __truncated__,..: 27 17
   19 1 35 36 37 30 ...
..$ refcount : num [1:50] 1 5 26 55 13 43 51 87 ...
..$ stacklength : int [1:50] 19 20 19 21 14 15 15 14 ...
..$ stackheadnodes: int [1:50] 52 52 52 52 52 52 52 ...
..$ stackleafnodes: int [1:50] 27 28 41 6 39 14 38 30 ...
..$ stackssrc : Factor w/ 50 levels "! [.data.frame [
  na.omit.data.frame na."| __truncated__,..: 27 28 39 5 37 13 36 30
profiles:List of 4
..$ data : int [1:522] 1 2 2 3 4 4 5 5 ...
..$ mem : NULL
.. $ malloc : NULL
..$ timesRLE:List of 2
.. ..- attr(*, "class")= chr "rle"
- attr(*, "class")= chr [1:2] "sprof" "list"
```

The nodes do not come in a specific order. Access via a permutation vector is preferred. This allows different views on the same data set. For example, table 4 on page 13 uses a permutation by total time, and a selection (compare to table 2 on page 9). The only difference is that **sprof** works uses a millisecond (ms) base, whereas R in general uses seconds as a base.

Table 3: splot result: by.self, self time > 4ms

-	name	self.time	self.pct	total.time	total.pct
30	lm.fit	87.00	16.67	113.00	1.47
6	[.data.frame	57.00	10.92	99.00	1.29
38	model.matrix.default	51.00	9.77	69.00	0.90
14	as.character	43.00	8.24	43.00	0.56
29	lm	42.00	8.05	474.00	6.16
51	$\operatorname{summary.lm}$	40.00	7.66	45.00	0.59
49	structure	32.00	6.13	33.00	0.43
41	na.omit.data.frame	26.00	4.98	114.00	1.48
13	anyDuplicated.default	22.00	4.21	22.00	0.29
16	as.list.data.frame	22.00	4.21	22.00	0.29
40	na.omit	20.00	3.83	134.00	1.74
39	model.response	13.00	2.49	56.00	0.73
36	model. frame. default	12.00	2.30	168.00	2.18
46	rep.int	7.00	1.34	7.00	0.09
10	<anonymous></anonymous>	6.00	1.15	522.00	6.79
28	list	5.00	0.96	5.00	0.07

At this level, it is helpful to note the expectations, and only then inspect the timing results. Since we are using a linear model, we are not surprised to see functions related to linear models on the top of the list. We may however be surprised to see functions related to data access and to character conversion very high on the list. The sizeable amount of time spent on NA handling is another aspect that is surprising.

Table 4: splot result: by total, total time > 4ms

	name	self.time	self.pct	total.time	total.pct
10	<anonymous></anonymous>	6.00	1.15	522.00	6.79
52	Sweave	0.00	0.00	522.00	6.79
21	doTryCatch	0.00	0.00	521.00	6.78
22	eval	1.00	0.19	521.00	6.78
23	evalFunc	0.00	0.00	521.00	6.78
55	try	0.00	0.00	521.00	6.78
56	tryCatch	0.00	0.00	521.00	6.78
57	tryCatchList	0.00	0.00	521.00	6.78
58	tryCatchOne	0.00	0.00	521.00	6.78

62	withVisible	0.00	0.00	521.00	6.78
< cut >	\vdots	:	:	:	:
61	vapply	3.00	0.57	23.00	0.30
13	anyDuplicated.default	22.00	4.21	22.00	0.29
16	as.list.data.frame	22.00	4.21	22.00	0.29
47	sapply	1.00	0.19	14.00	0.18
31	match	1.00	0.19	11.00	0.14
7	[[	0.00	0.00	8.00	0.10
8	[[.data.frame	1.00	0.19	8.00	0.10
25	FUN	1.00	0.19	7.00	0.09
46	rep.int	7.00	1.34	7.00	0.09
28	list	5.00	0.96	5.00	0.07

sampling sampling runctions e cleared vdots from string/-name columns

**ToDo:** remove text

Given the sampling structure of the profiles, two aspect are common. The sampling picks up scaffold functions with a high, nearly constant frequency. And the sampling will pick up rare recordings that are near to detection range. The display functions hide these effects by default. In our example, about half of the nodes are cleared by this garbage collector.

Common rearrangements as by total time and by self time are supplied by the display functions.

plot\_nodes(), for example, currently gives a choice of four displays for nodes, and supports trimming by default. Our profile starts with 62 nodes. The defaults cut off 34 nodes as uninformative, either because they are too rare, or ubiquitous.

```
#8
oldpar <- par(mfrow=c(2,2))
plot_nodes(sprof01)
par(oldpar)
```

Basic information on node level: see fig. 1 on the following page.

Information in the time scatterplots may sometimes more accessible when using a logarithmic scale, so this is added.

If you prefer, you can have the bar charts in horizontal layout, giving more space for labels.

```
#8
oldpar <- par(mfrow=c(2,2))
plot_nodes(sprof01, horiz=TRUE)
par(oldpar)
```

Basic information on node level - horizontal bars: see fig. 2 on page 15.

We can add colour. To illustrate this, we encode the frequency of the nodes as colour. As a palette, we choose a heat map here.

ToDo: apply colour to selection?
ToDo: spread colour on displayed part

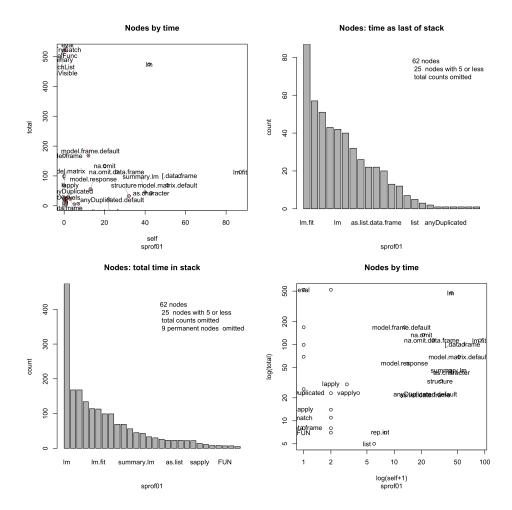


FIGURE 1. Basic information on node level

```
Input
freqrank01 <- rkindex(-sprof01$nodes$total.time, ties.method="random")
freqrankcol01 <- heat.colors(length(freqrank01))</pre>
```

Here is the node view using these choices:

```
#10
sprof01$nodes$icol <- freqrank01
oldpar <- par(mfrow=c(2,2))
plot_nodes(sprof01, col=freqrankcol01)
par(oldpar)
```

Basic information on node level, colour by total time: see fig. 2 on the facing page.

Colour is considered a volatile attribute. So you may need to pay some attention to keep colour indices (and colour palettes) aligned to your context. You may

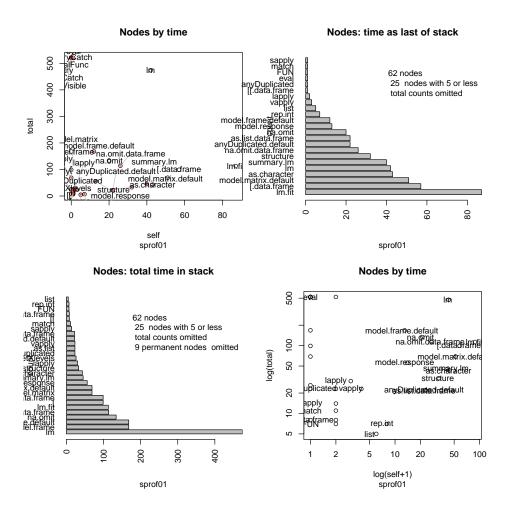


FIGURE 2. Basic information on node level

want to do experiments with colour, trying to find a good solution for your visual preferences. The recommended way is to use some stable colour index (the slot <code>icol</code> is reserved for this) and use this as an index to a choice of colour palettes. So <code>icol</code> becomes a part of the data structure, and the colour palette to be used is passed as a parameter. Package <code>RColorBrewer</code> may be a helpful source for colour palettes. You find more information and references on this in the vignette "Bertin matrices" of package <code>bertin</code>, <code>http://bertin.r-forge.r-project.org</code>.

1.1.3. *Node classes*. We can add attributes to the plots. But we can also add attributes to the nodes, and use these in the plots. In principle, this has been alway available. We are now making explicit use of this possibility.

The attribute *icol* is a special case which we used above. If present, it will be interpreted as an index to a colour table. For example, we can collect special well known functions in groups.

ToDo: improve colour: support colour in a structure

<u>ToDo:</u> colour by class – redo. Bundle colour index with colour?

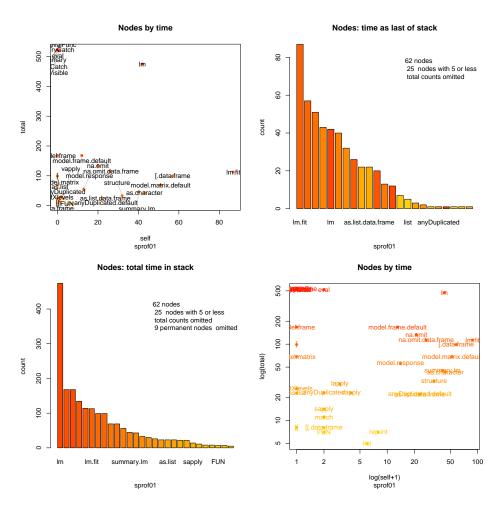


FIGURE 3. Basic information on node level, colour by total time.

The node information is to some part arbitrary. You may achieve the same functionality by different functions, and you will see different load in the profiles. Grouping nodes may be a mean to clarify the picture.

Grouping may also help you to focus your attention. "HOT" and "cold" may be ver helpful tags. These can be used in a flexible way.

ToDo: Move class attributes to package code
ToDo: add class by keyword

```
nodekeyword0 <- function(node)
{
}
```

```
nodepackages <- nodepackage(sprof01$nodes$name)
names(nodepackages) <- sprof01$nodes$name
table(nodepackages)
```

```
nodepackages
<not found> base stats utils
6 41 14 1
```

```
sprof01$nodes$icol <-as.factor(nodepackages)
```

Nodes by package, colours by RColorBrewer: see fig. 4 on the next page.

(Extend as you need it) and then use, as for example:

or use assignments on the fly

```
Input
nodeclass[sprof01$nodes$name %in%
        c("eval", "evalFunc",
                "try", "tryCatch", "tryCatchList", "tryCatchOne",
                "doTryCatch", "withVisible")
                ] <- "x_eval"
nodeclass[sprof01$nodes$name %in%
        c("model.frame", "model.matrix.default", "model.frame.default",
         " model.response", "model.matrix", "model.response")
                ] <- "x_model"</pre>
nodeclass[sprof01$nodes$name %in%
        c("lm", "lm.fit", "summary.lm")
                ] <- "x_1m"
nodeclass[sprof01$nodes$name =="<Anonymous>"] <- "x_Anon"
nodeclass[sprof01$nodes$name == "Sweave"] <- "x_Sweave"</pre>
nodeclass[sprof01$nodes$name %in% c( "summay" , "summay.lm")] <- "x_summary"
```

```
_______Input ______sprof01$nodes$icol <-as.factor(nodeclass)
```

adds a sticky colour attribute. To interpret, you should choose your preferred colour palette, for example

# gold cyan4 aquamarine pink violet orchid hotpink salmon turquoise1

<u>ToDo:</u> Defaults by class

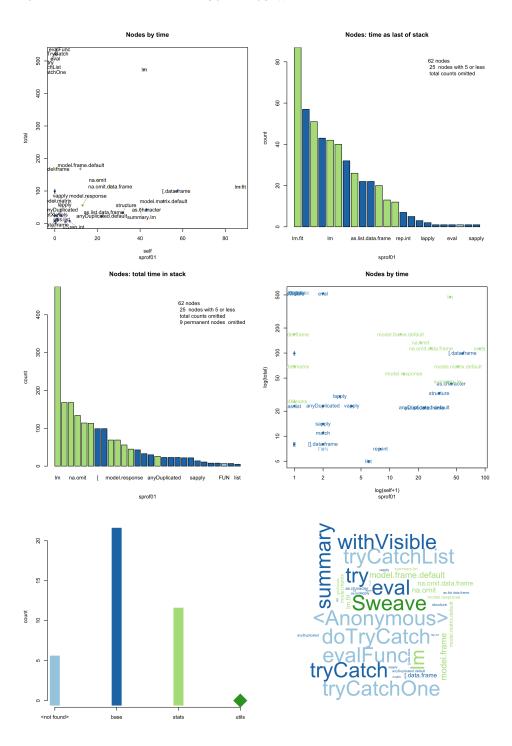


FIGURE 4. Nodes by package, colours by RColorBrewer.

```
#8 12
oldpar <- par(mfrow=c(3,2))
plot_nodes(sprof01, which=1:6, col=classcol)
par(oldpar)
```

\_ Input \_

Nodes by class, user defined colors: see fig. 5 on the following page.

<u>ToDo:</u> classes need separate colour palette

```
#8 12

oldpar <- par(mfrow=c(3,2))

plot_nodes(sprof01, which=1:6)

par(oldpar)
```

Nodes by class: default colour selection: see fig. 6 on page 21.

You can break down the frequency by classes of your choice. But beware of Simpson's paradox. The information you think you see may be strongly affected by your choices - what you see are reflections of conditional distributions. These may be very different from the global picture.

If package wordcloud is installed, a different view is possible. This is added in the plots above.

## 2. A BETTER GRIP ON PROFILE INFORMATION

The basic information provided by all profilers in R is a protocol of sampled stacks. The conventional approach is to break the information down to nodes and edges. The stacks provide more information than this. One way to access it is to use linking to pass information. This has already been used on the node level in section 1.1.2 on page 10.

2.1. The internal details. For each recorded event, the protocol records one line with a text string showing the sampled stack (in reverse order: most recent first). The stack lines may be preceded by header lines with event specific information. The protocol may be interspersed with control information, such as information about the timing interval used.

We know that the structural information, static information as well as dynamic information, can be represented with the help of a graph. For a static analysis, the graph representation may be the first choice. For a dynamic analysis, the stack information is our first information. A stack is a connected path in the program graph. If we start with nodes and edges, we loose information which is readily available in record of stacks.

As we know that we are working with stacks, we know that they have their peculiarities. Stacks tend to grow and shrink. Subsequent events will have extensions and shrinkages of stacks (if the recording is on a fine scale), or stack sharing common stumps (if the recording is on a coarser scale). We could exploit this information, but it does not seem worth the effort.

ToDo: add attributes to stacks, and discuss scope
ToDo: sorting/arranging stacks

ToDo: re-think: sort stacks

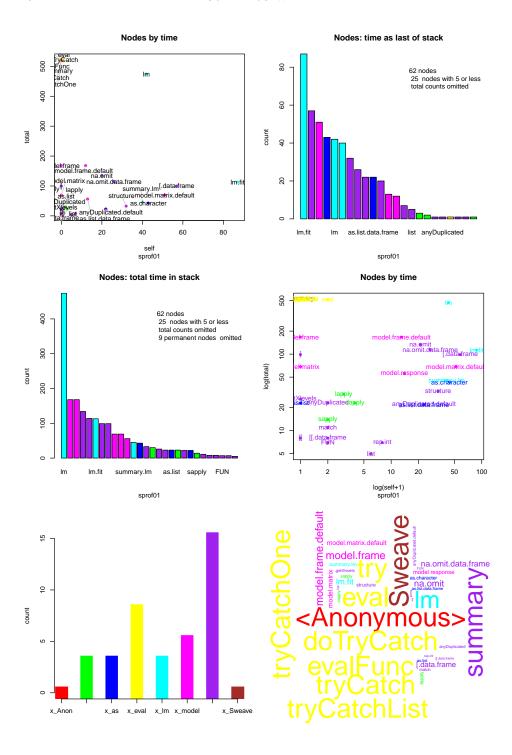


FIGURE 5. Nodes by class. User defined colors.

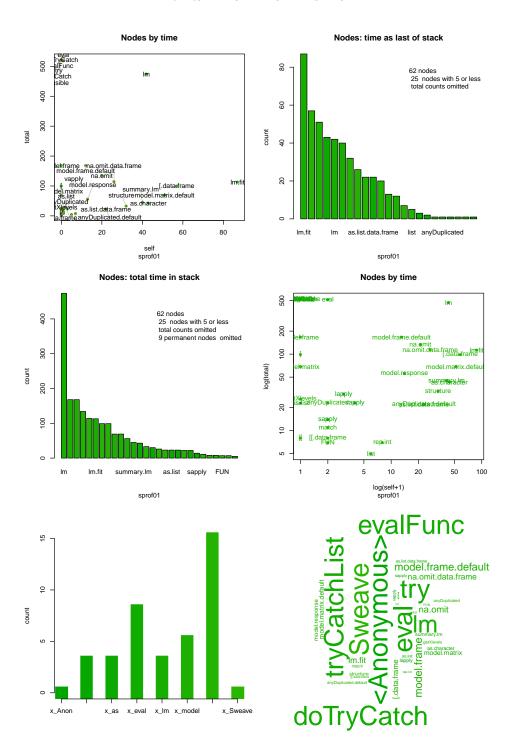


Figure 6. 8 12 Nodes by class: default colour selection

There have always been interrupts, and these show up in profiles. In R, there is a related problem: garbage collection (GC) may interfere and leave traces in the stack.

Stack information is first. The call graph is a second instance that is (re)constructed from the stack recording. The graph represents cumulated one-step information. Longer scale information contained in the stacks is lost in the graph.

Here is the way we represent the profile information:

The profile log file is sanitised:

- Control lines are extracted and recorded in a separate list.
- Head parts, if present, are extracted and recorded in a matrix that is kept line-aligned with the remainder
- Line content is standardised, for example by removing stray quotation marks etc.

After this, the sanitised lines are encoded as a vector of stacks, and references to this.

If necessary, these steps are done by chunks to reduce memory load.

From the vector of stacks, a vector of nodes (or rather node names) is derived.

The stacks are now encoded by references to the nodes table. For convenience, we keep the (sanitised) textual representation of the stacks.

So far, texts are in reverse order. For each stack, we record the trailing leaf, and then we reverse order. The top of stack is now on first position.

Several statistics can be accumulated easily as a side effect.

Conceptually, the data structure consist of three tables (the implementation may differ, and is subject to change).

The profiles table is the representation of the input file. Control lines are collected in a special table. With the control lines removed, the rest is a table, one row per input line. The body of the line, the stack, is encoded as a reference to a stacks table (obligatory) and header information (optional).

The stacks table contains the collected stacks, each stack encoded as a list of references to the node table. This is obligatory. This list is kept in reverse order (root at position 1). A source line representing the stack information may be kept (optional).

The nodes table keeps the names at the nodes.

Sometimes, it is more convenient to use a simple representation, such as a matrix. Several extraction routines are provided for this, and the display routines make heavy use of this. See table 5 on the next page.

We now can go beyond node level.

This is what we get for free from the node information on our three levels: node, stack, and profile.

<u>ToDo:</u> complete matrix conversion

ToDo: check and stabilise colour linking

<u>ToDo:</u> clean up colour handling

Table 5. Extraction and conversion routines

<pre>profiles_matrix()</pre>	incidence matrix: nodes by event
stacks_matrix()	incidence matrix: nodes by stack
list.as.matrix()	fill list to equal length and convert to matrix
stackstoadj()	stacks to (correspondence) adjacency matrix
adjacency()	sprof to (correspondence) adjacency matrix

#8 rainbow sprof01\$nodes\$icol <- freqrank01; freqrankcol <- rainbow(62) shownodes(sprof01, col=freqrankcol)

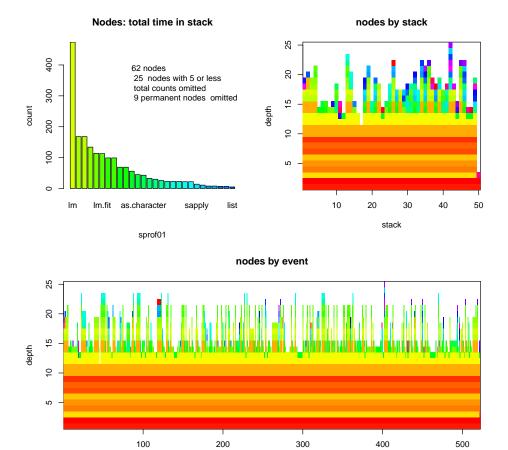


FIGURE 7. Nodes by stack and profile

event

See fig. 7 for a summary by nodes.

The obvious message is that if seen by stack level, there are different structures. Profiling usually takes place in a framework. So at the base of the stacks, we find entries that are (almost) persistent. Then usually we have some few steps where the algorithm splits, and then we have the finer details. These can be identified using information on the stack level, but of course they are not visible on the node or edge level in a graph representation. On the stack level, we see a socket. If we want a statistic, we can look at number of different nodes by level.

```
stacks_nodes <- list.as.matrix(sprof01$stacks$nodes)
 nrnodes <- apply(stacks_nodes,1,function(x) {length(unique(x))})</pre>
 cat("nr unique nodes per stack level\n")
                                     Output
nr unique nodes per stack level
                                     Input
 nrnodes
                                    Output ______ 2 2 4 11 12 10 10 16
 [1]
[23]
     3
         2
                                     Input
 plot(x=nrnodes, y= 1:length(nrnodes), xlab="nr of unique nodes", ylab="stack level")
 abline(h=2.5,col="green")
 abline(h=12.5,col="green")
```

<u>ToDo:</u> check and synchronise

Nr. of unique nodes by stack level: See fig. 8 on the next page.

We will come to finer tools in section 2.4 on page 34 but for the moment the rough information should suffice to take a decision. In our example, it is only a matter of taste whether we cut off 12 levels, or we want to work with five components after cutting 13 levels

Not so often, but a frequent phenomenon is to have some "burn in" or "fade out". To identify this, we need to look at the profile level. The indicator to check is to whether we have very low frequency stacks at the beginning or the end of our recording. The counts to be takes as reference can be seen from the summary.

<u>ToDo:</u> we could do smart smoothing of the stacks here

This summary has to be taken with caution. As the program runs, the stacks are build up und teared down, and we only take random samples. So in dynamic parts, we see images with some fluctuation, as one stack may be a snapshot of an other under construction. A better information is to cut off fluctuations and use this summary as a reference.

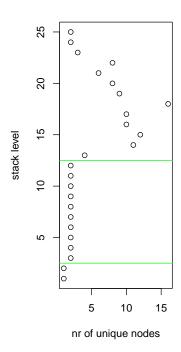


FIGURE 8. Nr of unique nodes by stack level.

```
summary(sprof01$stacks$refcount[sprof01$stacks$refcount>2])
                          _____ Output
Mean 3rd Qu.
  Min. 1st Qu.
                Median
                                           Max.
  3.00
          7.75
                 16.00
                          24.30
                                 40.50
                                  \_ Input \_
df <- data.frame(stack=sprof01$profiles$data,
        count=sprof01$stacks$refcount[sprof01$profiles$data])
prxt(df, caption="Stacks by event: burn in/fade out",
        label="tab:margin",
        digits=c(0,0,0) )
                 Table 6: Stacks by event: burn in/fade out
```

	stack	count
1	1	1
2	2	5

3	2	5
4	3	26
5	4	55
6	4	55
7	5	13
8	5	13
9	6	43
10	7	51
< cut >	:	:
513	3	26
514	3	26
515	3	26
516	3	26
517	4	55
518	3	26
519	36	2
520	16	42
521	44	2
522	50	1

Here at least one recording on either side is a candidate to be off. We may have a look at the next recordings and decide to go beyond and cut off events 1:3 and 519:522.

At a closer look, we may find stack patterns (maybe marked by specific nodes) that indicate administrative intervention and rather should be handled as separators between distinct profiles rather than as part of the general dynamics. Again we may use some indicator nodes to be used as marker for special stacks.

Stable framework effects sometimes are obvious and can be detected automatically. "burn in" or "fade out' may need a closer look, and special stacks need and individual inspection on low frequency stacks. Tools for trimming are in section 2.3.1 on page 30.

2.2. **The free lunch.** What you have seen so far is what you get for free when using package *sprof*.

If you want to wrap up the information and look at it from a graph point of view, here is just one example. More are in section 3 on page 39 and vrefsec:moregraph. But before changing to the graph perspective, we recommend to see the next sections, not to skip them.

The preview, at this point, taking package graph as an example<sup>1</sup>. graph on its side has an undocumented feature: it needs Rgraphviz to handle graph attributes<sup>2</sup>. We

**ToDo:** example

**ToDo:** colours. recolour. Propagate colour to graph.

<sup>&</sup>lt;sup>1</sup>Package 'graph' was removed from the CRAN repository.

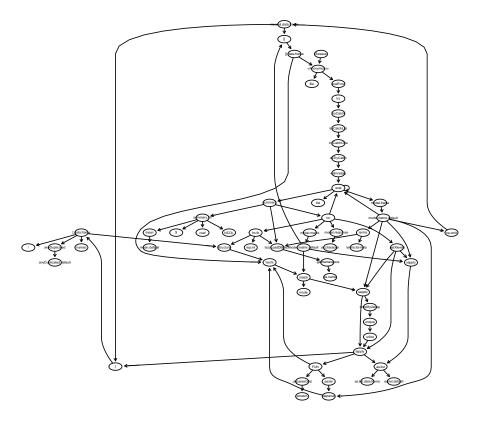
This package is now available from Bioconductor only, see http://www.bioconductor.org/packages/release/bioc/html/graph.html.

<sup>&</sup>lt;sup>2</sup>Package 'Rgraphviz' was removed from the CRAN repository.

This package is now available from Bioconductor only.

have to take two steps. We extract the graph information from *sprof*. Using an adjacency matrix is a simple solution here. This is then converted to the "graph-NEL" format which is shared by graph and Rgraphviz. Rgraphviz is hidden in the use of plot(). So here is a bare foot approach. A more sophisticated function implementation is in section 3 on page 39.

# sprof01: graph layout example



See http://www.bioconductor.org/packages/release/bioc/html/Rgraphviz.html.

ToDo: updateRprof needs careful checking. For now, we are including long listings here to provide the necessary information

2.3. Cheap thrills. Before starting additional inspection, the data better be trimmed. Trimming routines are in section 2.3.1 on page 30, but the data structure is robust enough to allow manual intervention as used here.

```
sprof02 <- sprof01; sprof02$info$id <- "sprof02: trimmed"</pre>
```

On the stack level, we take brute force to cut off the basic stacks.

```
basetrim <- 13
sprof02$stacks$nodes <- sapply(sprof02$stacks$nodes,
   function (x){if (length(x)> basetrim) x[-(1:basetrim)] })
```

We have noted burn in/fade out. This is on the profile level. Taking the big knife is not advisable, since time information and stack data must be synchronised. So we are more cautious.

```
sprof02$profiles$data[1:3] <- NA
sprof02$profiles$data[519:522] <- NA
```

At this point, it is a decision whether to adapt the timing information, or keep the original information. Since this decision does affect the structural information, it is not critical. But analysis is easier if unused nodes are eliminated. The <code>info</code> section is inconsistent at this point. Another reason to call <code>updateRprof()</code>.

```
_ Input _
strx(sprof02$info)
                                 _ Output _
##strx: sprof02$info
'data.frame': 1 obs. of 9 variables:
$ id : chr "sprof02: trimmed"
$ date : POSIXct, format: "2013-08-15 22:02:43"
$ nrnodes : int 62
$ nrstacks : int 50
$ nrrecords : int 522
$ sample.interval: num 0.001
$ sampling.time : num 0.522
$ ctllines : Factor w/ 1 level "sample.interval=1000": 1
$ ctllinenr : num 1
                                  _ Input _
prxt(sprof02$nodes,
         caption="sprof02, before update",
         label="tab:sprof02info1",
         digits=c(0,0,0,2,0,2,0),
         zero.print=" . "
         )
                       Table 7: sprof02, before update
```

ToDo: handle empty stacks and zero counts gracefully

<u>ToDo:</u> add a purge function

	name	self.time	self.pct	total.time	total.pct	icol
1	!	2	0.38	2	0.03	46
2	getNamespace		0.00	1	0.01	60
3	. departs = Opts	2	0.38	4	0.05	43
4	.get $X$ levels		0.00	26	0.34	28
5			0.00	99	1.29	20
6	data.frame	57	10.92	99	1.29	19
7	Ī		0.00	8	0.10	36
8	[[.data.frame	1	0.19	8	0.10	35
9	%in $%$	1	0.19	4	0.05	40
10	<anonymous></anonymous>	6	1.15	522	6.79	2
< cut >	$\setminus vdots$	:	:	:	:	:
53	terms		0.00	2	0.03	52
54	terms.formula	1	0.19	1	0.01	59
55	try		0.00	521	6.78	7
56	tryCatch		0.00	521	6.78	8
57	tryCatchList		0.00	521	6.78	10
58	tryCatchOne		0.00	521	6.78	5
59	unique	3	0.57	4	0.05	42
60	unlist		0.00	1	0.01	54
61	vapply	3	0.57	23	0.30	30
62	withVisible		0.00	521	6.78	4

\_\_\_\_\_ Input \_\_\_\_\_ sprof02 <- updateRprof(sprof02) sprof02\$info\$id <- "sprof02 updated"</pre> \_\_\_\_\_ Input \_\_\_\_\_ strx(sprof02\$info) \_\_\_\_\_ Output \_\_\_\_\_ ##strx: sprof02\$info 'data.frame': 1 obs. of 10 variables: \$ id : chr "sprof02 updated" \$ date : POSIXct, format: "2013-08-15 22:02:43" \$ nrnodes : int 62 \$ nrstacks : int 50 \$ nrecords : int 522\$ sample.interval: num 0.001 \$ sampling.time : num 0.522  $\$  ctllines : Factor w/ 1 level "sample.interval=1000": 1 \$ ctllinenr : num 1 \$ date\_updated : POSIXct, format: "2013-08-15 22:02:48"

\_\_\_\_\_\_ Input \_\_\_\_

Table 8: sprof02, after update

	name	self.time	self.pct	total.time	total.pct	icol
1	!	1	0.23	1	0.06	46
2	get Name space		0.00	1	0.06	60
3	. departs = Opts	2	0.46	4	0.25	43
4	.get $X$ levels		0.00	26	1.64	28
5	[		0.00	98	6.17	20
6	[.data.frame	57	13.16	98	6.17	19
7	[[		0.00	8	0.50	36
8	[[.data.frame	1	0.23	8	0.50	35
9	$\% \mathrm{in}\%$	1	0.23	4	0.25	40
10	<anonymous></anonymous>	6	1.39	6	0.38	2
< cut >	\vdots	:	:	:	:	:
53	terms		0.00	1	0.06	52
54	terms.formula	1	0.23	1	0.06	59
55	try		0.00		0.00	7
56	tryCatch		0.00		0.00	8
57	tryCatchList		0.00		0.00	10
58	tryCatchOne		0.00		0.00	5
59	unique	3	0.69	4	0.25	42
60	unlist		0.00	1	0.06	54
61	vapply	3	0.69	23	1.45	30
62	withVisible		0.00		0.00	4

#8 8
shownodes(sprof02)

Nodes by stack and profile: see fig. 9 on the next page.

ToDo: This section needs to be reworked ToDo: trimexample ToDo: add trim by keyword

2.3.1. Trimming. Note: trimming may be supported by the graph packages.

```
trimstacks <- function(sprof, level){
lapply(sprof$stacks$nodes, function(x) {x[-(1:level)]})
}</pre>
```

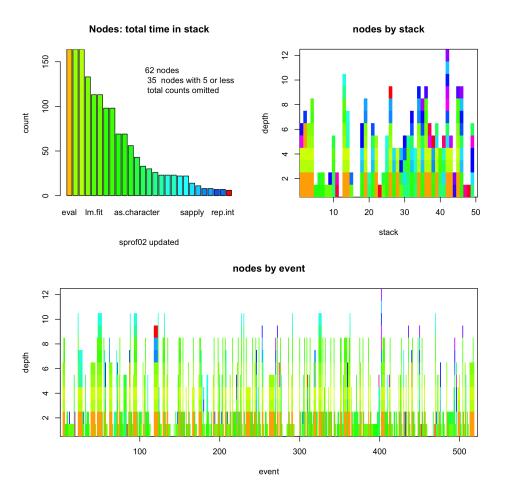


FIGURE 9. sprof02: Nodes by stack and profile

```
sprof01Tr <- trimstacks(sprof01, 11)
#profile_nodesTr <- profiles_matrix(sprof01Tr)
#image(x=1:ncol(profile_nodesTr),y=1:nrow(profile_nodesTr), t(profile_nodesTr),xlab="event", ylab="</pre>
```

There is no statistics on profiles. Profiles are our elementary data. However we can link to our derived data to get a more informative display. For example, going one step back we can encode stacks and use these colour codes in the display of a profile.

Or going two steps back, we can encode nodes in colour, giving coloured stacks, and use these in the display of profile data.

 $2.3.2.\ Surgery.$  Note: surgery may be supported by the graph packages. Use the implementation you are more familiar with.

Looking at nodes gives you a point-wise horizon. Looking at edges gives you a one step horizon. The stacks give a wider horizon, typically a step size of 10 or more. The stacks we get from R have peculiarities, and we can handle with this broader perspective. These are not relevant if we look point-wise, but may become dominating if we try to get a global picture. We take a look ahead (details to come in section 3 on page 399 and have a preview how our example is represented as a graph. Left is the original graph as recovered from the edge information, right the graph after we have cut off the scaffold effects.

**ToDo:** cut next level

Control structures may be represented in R as function, and these may lead to concentration points. Using information from the stacks, we can avoid these by introducing substitute nodes on the stack level. For example, lapply is appearing in various contexts and may be confusing any graph representation. We can avoid this by replacing a short sequence.

```
ToDo: Implement.
Currently best han-
dled on source=text
level
```

ToDo: function addnode using "call by reference" to be added

```
"[" "lapply" ".getXlevels" -> "<.getXlevels_[>"
```

If the node does not exist, we want to add it to our global variable. For now, we do it using expressions on the R basic level and avoid tricks like simulating "call by reference".

```
sprof03 <- sprof02; sprof03$info$id <- "sprof03: surgery"
node <- "<.getXlevels_[>"
#nodei <- function(sprofx, node, warn = TRUE)</pre>
        i <- match(node, sprof03$nodes$name, nomatch=0)</pre>
        if (i==0){
                 sprof03$nodes$name <- as.character(sprof03$nodes$name)</pre>
                 sprof03$nodes <- rbind(sprof03$nodes,NA)</pre>
                 i <- length(sprof03$nodes$name)</pre>
                 sprof03$nodes$name[i] <- node</pre>
        if (as.logical(options("warn")))
                 message("node added. An updateRprof() may be necessary.")
        nodei <- i
}
# sprof <- sprof01; nodei(sprof, "kiki"); sprof$nodes</pre>
```

Now we have to identify the stacks that may get a replacement. First find the candidates.

```
Input
 targeti <- match("lapply", sprof03$nodes$name, nomatch=0)
 found <- lapply(sprof03$stacks$nodes, function(X) match(targeti,X))</pre>
 found <- data.frame(stack=1:length(found), position=as.matrix(found))</pre>
 found[!is.na(found$position),]
                                    _ Output _
   stack position
10
      10
                2
21
      21
                 5
```

```
30
      30
                3
31
      31
                2
32
      32
                5
33
      33
                2
35
      35
                5
                6
39
     39
                2
40
     40
41
                3
      41
                6
42
      42
                5
46
      46
      49
                                    Input
 # as.factor(sprof03$stacks$nodes[!is.na(found)],
                  levels=1: length(sprof03$nodes$name), labels=sprof03$nodes$name)
For now, these are just candidates.
                                                                                 <u>ToDo:</u> implement
                                                                                replacement on the
Other candidates are:
                                                                                stack level.
"as.list" "vapply" "model.frame.default" -> "<model_as.list>"
or
"as.list" "vapply" "model.matrix.default" -> "<model_matrix_as.list>" ToDo: implement
newchopnode <- function(nodenames, chop) {
 tmpname <- paste("<",as.character(nodenames[chop]),">")
 # chec for existing.
 # add if necessary
 tmpname
 chopstack <- function(x , chop, replacement)</pre>
 # is chop in x`
 # y: cut x.
```

We use a prepared sanitised version of our data set.

# merge x <- head + replacement + tiail</pre>

**ToDo:** fix null name

```
sprof04 <- readRprof("RprofsRegressionExpl03.out", id="sprof04")
```

Applying our old classification:

return(x)

}

```
#8 12
oldpar <- par(mfrow=c(3,2))
plot_nodes(sprof04, which=1:6, col=classcol)
par(oldpar)
```

Sanitised nodes by class, user defined colours: see fig. 10 on the next page.

ToDo: needs serious
revision
ToDo: test na removal in rrle

2.4. Run length. For a visual inspection, runs of the same node and level in the profile are easily perceived. For an analytical inspection, we have to reconstruct the runs from the data. In stacks, runs are organised hierarchically. On the root level, runs are just ordinary runs. On the next levels, runs have to be defined given (within) the previous runs. So we need rrle(), a recursive version of rle, applied

level, runs are just ordinary runs. On the next levels, runs have to be defined given (within) the previous runs. So we need rrle(), a recursive version of rle, applied to the profile information. This gives a detailed information about the presence time of each node, by stack level.

ToDo: use sprof02 or sprof03?
ToDo: handle NA as special case

```
_ Output _
##strx: profile_nodes_rle
List of 12
$ :List of 2
..$ lengths: int [1:365] 1 1 1 3 3 1 7 1 ...
..$ values : int [1:365] NA NA NA 22 39 37 30 4 ...
..- attr(*, "class")= chr "rle"
$ :List of 2
..$ lengths: int [1:411] 1 1 1 3 1 1 1 1 ...
..$ values : int [1:411] NA NA NA 22 NA NA 14 38 ...
..- attr(*, "class")= chr "rle"
$ :List of 2
..$ lengths: int [1:431] 1 1 1 3 1 1 1 1 ...
..$ values : int [1:431] NA NA NA 35 NA NA NA NA ...
..- attr(*, "class")= chr "rle"
$ :List of 2
```

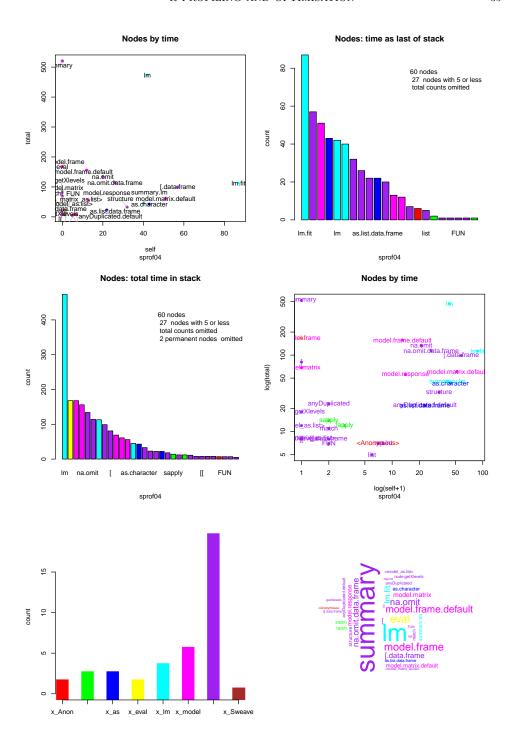


FIGURE 10. Sanitised. Nodes by class. User defined colours.

```
..$ lengths: int [1:431] 1 1 1 3 1 1 1 1 ...
..$ values : int [1:431] NA NA NA 36 NA NA NA NA ...
..- attr(*, "class")= chr "rle"
$ :List of 2
..$ lengths: int [1:452] 1 1 1 3 1 1 1 1 ...
..$ values : int [1:452] NA NA NA 40 NA NA NA NA ...
..- attr(*, "class")= chr "rle"
[list output truncated]
```

Input \_

On a given stack level, the run length is the best information on the time used per call, and the run count of a node is the best information on the number of calls. So this is a prime starting point for in-depth analysis.

If you need it, you can represent the run length information by level as a matrix. This is expanding a sparse matrix to full and should be avoided.

```
profile_nodes_rlearray <- nodesprofile(sprof02)
strx(profile_nodes_rlearray)
```

```
##strx: profile_nodes_rlearray
num [1:61, 1:12, 1:7] 0 1 0 17 0 0 0 0 ...
- attr(*, "dimnames")=List of 3
..$ node : chr [1:61] "!" "..getNamespace" ".deparseOpts" ...
..$ level : chr [1:12] "1" "2" "3" ...
..$ run_length: chr [1:7] "1" "2" "3" ...
```

This allows us to extract marginal from provlev[ node, level, run length].

```
nn <- profile_nodes_rlearray["model.frame", , ]
print.table(addmargins(nn), zero.print = ".")</pre>
```

```
Output
```

ToDo: keep as factor. This is a sparse cube with margins node, stack level, run length. Nodes are mostly concentrated on few levels.

<u>ToDo:</u> Warning: data structure still under discussion

<u>ToDo:</u> hack. replace by decent vector/array based implementation

<u>ToDo:</u> add names for node dimension

 $\begin{array}{ll} {\bf \underline{ToDo:}} & {\rm add} & {\rm summary \ for \ NA} \end{array}$ 

ToDo: add marginals and conditionals.

 $\begin{array}{ll} Provide & function \\ node\_summary. \end{array}$ 

<u>**ToDo:**</u> rescale to application scale

<u>**ToDo:**</u> replace sum by weighted sum

ToDo: allow sorting, e.g. by marginals

\_ Input \_

```
amt <- nodesrunlength(sprof02)</pre>
 # print.table(as.table(amt[,1:9]), zero.print = ".")
 ## ??? print.table changes format when col.10 is present.
But print.xtable has no zero.print ...
                                                                                      ToDo:
                                                                                               needs
                                                                                      to be hidden for
                                                                                      print.xtable
                                     \_ Input \_
 #special sanitising for xtable
                                                                                      <u>ToDo:</u> handle gaps
 xr <- rownames(amt)</pre>
                                                                                      in run length = NA
 \#for \ (i \ in \ (1:length(xr))) \ \{xr[i] \leftarrow sub(xr[i],"\backslash [","\$[\$", \ fixed=TRUE)\}
                                                                                      counts
 #xr <- paste("$",xr,"$")</pre>
 xr <- gsub("[","{[}",xr, fixed=TRUE)</pre>
 xr <- gsub("_","\\_",xr, fixed=TRUE)</pre>
 xr <- gsub("^","\\^",xr, fixed=TRUE)</pre>
 rownames(amt)<- xr</pre>
 prxt(amt,
  caption="Marginal statistics on nodes by run length, sorted by total time used, count > 1 only",
         label="tab:pramt1",
         digits=c(rep(0,dim(amt)[2]),2),
         zero.print=" . ") #dim(amt)[2]-1, +1 for rownames
```

Table 9: Marginal statistics on nodes by run length, sorted by total time used, count > 1 only

	1	2	3	4	5	6	7	nr_runs	total_time	avg_time
eval	86	34	14	8	4	12	2	160	334	2.09
model.frame	40	17	7	4	2	6	1	77	164	2.13
model.frame.default	40	17	7	4	2	6	1	77	164	2.13
na.omit	46	10	5	4	1	4	1	71	133	1.87
$_{ m lm.fit}$	46	18	3	1	1	1	1	71	113	1.59
na.omit.data.frame	43	7	4	5		4		63	113	1.79
{[}	59	4	3	4		1		71	98	1.38
$\{[]\}.data.frame$	59	4	3	4		1		71	98	1.38
model.matrix	55	4	2			•		61	69	1.13
model.matrix.default	55	4	2					61	69	1.13
model.response	35	3	3			1		42	56	1.33
as.character	34	3	1					38	43	1.13
structure	21	1		1		1		24	33	1.38
lapply	26			1				27	30	1.11
.get $X$ levels	17	1	1	1				20	26	1.30
anyDuplicated	10			2	1			13	23	1.77
as.list	16	1			1			18	23	1.28
vapply	13	1	1		1			16	23	1.44
anyDuplicated.default	9			2	1			12	22	1.83
as.list.data.frame	15	1			1			17	22	1.29
sapply	14							14	14	1.00
match	12							12	12	1.00
{[}{[}	3				1			4	8	2.00
$\{[\}\{[\}.data.frame$	3				1			4	8	2.00
FUN	7							7	7	1.00

rep.int	7				7	7	1.00
<anonymous></anonymous>	1		1		2	6	3.00
. departs = Opts	4				4	4	1.00
%in $%$	4				4	4	1.00
simplify2array	4				4	4	1.00
unique	4				4	4	1.00
list	3				3	3	1.00
deparse	2				2	2	1.00
$\operatorname{mode}$	2				2	2	1.00
names	2				2	2	1.00
pmatch	2				2	2	1.00
!	1				1	1	1.00
get Name space	1				1	1	1.00
\$	1				1	1	1.00
as.list.default	1				1	1	1.00
as.name	1				1	1	1.00
coef	1				1	1	1.00
lazyLoadDBfetch	1				1	1	1.00
mean	1				1	1	1.00
mean.default	1				1	1	1.00
NCOL	1				1	1	1.00
paste	1				1	1	1.00
terms	1				1	1	1.00
terms.formula	1				1	1	1.00
unlist	1				1	1	1.00

ToDo: add current level
ToDo: generate a coplot representation
ToDo: remove counts per level?
ToDo: hack. keep length in nodesrun-

length

See table 9: Marginal statistics on nodes by run length.

We can add this information to our base and use it to enhance graphical displays. See: nodes marked by run length and run count, fig. 15 on page 47.

eval is the base of all evaluation in R, so we should not be surprised to find it at the top of the list. The next two entries, model.frame and model.frame.default are seen 77 times. We know that in our example we had a loop with 100 repetitions, so a frequency of 100 would not be surprising. A closer look shows that both functions occur as isolated events 40 times, but frequency decreases with run length until we see another high at run length 6. We are sampling, and of course sampling can miss some function calls. But this pattern is the oopposite. This is what if occurs if we do not miss a function call, but we miss a gap. So several function calls are joined and appear as some longer runs, typically a multiple of the original run length.

We can improve the hit rate by increasing the sampling rate. But of course this is at the expense of using more space for the log files, and increased time for the overhead. In our case, 1 ms seems to be a good compromise, but 0.1 ms might be another feasible choice.

As we walk down the list, na.omit is next, followed closely by na.omit.data.frame. In our problem, there are no missing data. But R does not have something like a "vanilla mode". The overhead used for the handling of potentially missing values

would need a restructuring of the linear model algorithm, e.g. by introducing a "has.na" flag.

Walking further down, we see other candidates such as **as.character** that may be avoidable in this problem.

```
Input

nr <- nodesrunlength(sprof02, clean=FALSE)

#nr <- rbind(nr,0) ## hack. keep length in nodesrunlength

rownames(nr) <- sprof02$nodes$name

sprof02$nodes <- cbind(sprof02$nodes, nr)

rownames(sprof02$nodes) <- sprof02$nodes$name
```

3. Graph Package

#runs min median run length max

ToDo: table: node

What we have achieved so far can be seen from the graph representations.

```
_ Input _
 library(graph)
 search()
                                   Output .
 [1] ".GlobalEnv"
                             "package:Rgraphviz"
 [3] "package:graph"
                             "package:sna"
 [5] "package:grid"
                             "package:wordcloud"
 [7] "package:RColorBrewer" "package:Rcpp"
[9] "package:xtable"
                             "package:sprof"
[11] "package:stats"
                             "package:graphics"
[13] "package:grDevices"
                             "package:utils"
[15] "package:datasets"
                             "package:methods"
[17] "Autoloads"
                             "package:base"
                                   _{-} Input
#12 6
 oldpar <- par(mfrow=c(1,2))</pre>
 plotviz(as(adjacency(sprof01), "graphNEL"),
          sub=sprof01$info$id)
 plotviz(as(adjacency(sprof02), "graphNEL"),
          sub=sprof02$info$id)
 par(oldpar)
```

See fig. 11 on the following page for a comparison before and after trimming. The scaffold effect are removed from the picture on the right side. This cuts off the uninformative spine, and induces minor changes in the body of the graph. You can do additional trimming, if you want.

R is function based, and control structures in general are implemented as functions. In a graph representation, they appear as nodes, concentrating and seeding to unrelated paths. We can detect these on the stack level and replace them by surrogates, introducing new nodes. This is a case for surgery.

We use a prepared sanitised version of our data set.

**ToDo:** fix null name

#### plotviz( as(adjacency(sprof01), "graphNEL"), dot ) plotviz( as(adjacency(sprof02), "graphNEL"), dot )

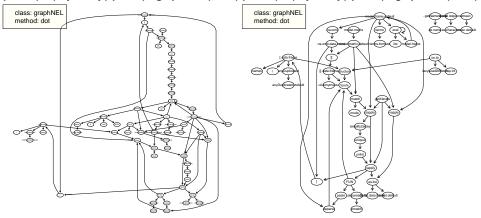


FIGURE 11. Sprof graph, before and after trimming.

```
sprof04 <- readRprof("RprofsRegressionExpl03.out", id="sprof04")</pre>
```

#### plotviz( as(adjacency(sprof02), "graphNEL"), dot ) plotviz( as(adjacency(sprof04), "graphNEL"), dot )

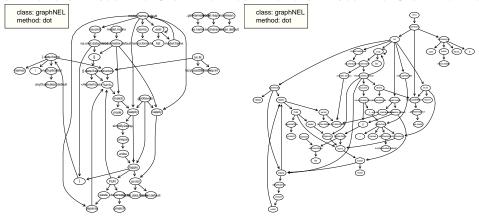


FIGURE 12. Sprof graph, after trimming, and after trimming and some sanitising.

**ToDo:** cut top levels

Now the structure becomes obvious. Cutting off may be taken two levels deeper. This would completely separate the lm() branch from the summary.lm() In the lm() branch, there are tree nodes (lapply(), %in% and [.data.frame) that are a cases for additional surgery to avoid focusing affects in the graph display. It is your choice to remove them, or live with them.

We know ho to create standard graph displays from this. The next step is to encode additional information we have from the profiles as attribute to the graph.

The derived edge frequency is the first bit of information. Implicitly, it can be used as weight in the graph placement routines. We make this explicit by giving a choice whether to use it or not. Irrespective of this choice, we encode reference counts as line width of the edges.

This function is not included in the package to avoid dependency of **sprof** on **graph** and other graph packages.

```
Input -
library(graph)
search()
                                   Output
[1] ".GlobalEnv"
                            "package:Rgraphviz"
 [3] "package:graph"
                            "package:sna"
[5] "package:grid"
                             "package:wordcloud"
[7] "package:RColorBrewer" "package:Rcpp"
[9] "package:xtable"
                            "package:sprof"
[11] "package:stats"
                             "package:graphics"
[13] "package:grDevices"
                             "package:utils"
[15] "package:datasets"
                             "package:methods"
[17] "Autoloads"
                             "package:base"
                                   _ Input _
library(Rgraphviz)
search()
                                   Output
[1] ".GlobalEnv"
                             "package:Rgraphviz"
[3] "package:graph"
                             "package:sna"
[5] "package:grid"
                             "package:wordcloud"
[7] "package:RColorBrewer" "package:Rcpp"
[9] "package:xtable"
                             "package:sprof"
[11] "package:stats"
                             "package:graphics"
[13] "package:grDevices"
                             "package:utils"
[15] "package:datasets"
                             "package:methods"
[17] "Autoloads"
                             "package:base"
                                  _ Input _
```

This is a common routine for the graph and Rgraphviz package.

```
as_graphNEL_sprof <- function(sprof, weight=TRUE){</pre>
a04<-adjacency(sprof)
rnames <- rownames(a04)</pre>
if (!weight) {
dimold \leftarrow dim(a04); a04 \leftarrow as.numeric(a04); dim(a04) \leftarrow dimold
rownames(a04)<- rnames; colnames(a04)<- rnames;</pre>
} #! define lwd first
el04 <- edgematrix(a04)</pre>
el04$lwd <- rkindex(el04$count, maxindex=6, ties.method="min")
a04NEL <- as(a04, "graphNEL")
nodeDataDefaults(a04NEL, "shape") <- "ellipse"</pre>
nodeDataDefaults(a04NEL, "cex") <- 0.6</pre>
nodeDataDefaults(a04NEL, "weigth") <- 1</pre>
nodeDataDefaults(a04NEL, "fill") <- "green"</pre>
nodeDataDefaults(a04NEL, "col") <- "yellow"</pre>
a04NEL <- layoutGraph(a04NEL)
nodeRenderInfo(a04NEL) <- list(shape="ellipse")</pre>
nodeRenderInfo(a04NEL) <- list(cex=0.6, shape="ellipse")</pre>
nodeRenderInfo(a04NEL) <- list(weight=1)</pre>
#nodeRenderInfo(a04NEL) <- list(color="yellow")</pre>
nodeRenderInfo(a04NEL) <- list(fill="yellow", col="blue")</pre>
edgeDataDefaults(a04NEL,"lwd") <- 1</pre>
edgeDataDefaults(a04NEL, "col") <- "grey"</pre>
#nodeRenderInfo(a04NEL) <- list(weight=1)</pre>
#edgeRenderInfo(a04NEL) <- list(lwd=e104$1wd)</pre>
#edgeRenderInfo(a04NEL)$1wd <- e104$1wd</pre>
for (i in 1:length(el04$lwd))
{edgeRenderInfo(a04NEL)$lwd[i] <- e104$lwd[i]}</pre>
a04NEL
plot_graphNEL_sprof <- function(graphNEL)</pre>
         main1 <- deparse(substitute(graphNEL))</pre>
         if (inherits(graphNEL, "sprof")) {
         main1 <- graphNEL$info$id
         graphNEL <-
                  as_graphNEL_sprof(graphNEL, weight=FALSE)
         renderGraph(graphNEL)
```

```
a04NEL <- as_graphNEL_sprof(sprof04)
plot_graphNEL_sprof(a04NEL)
```

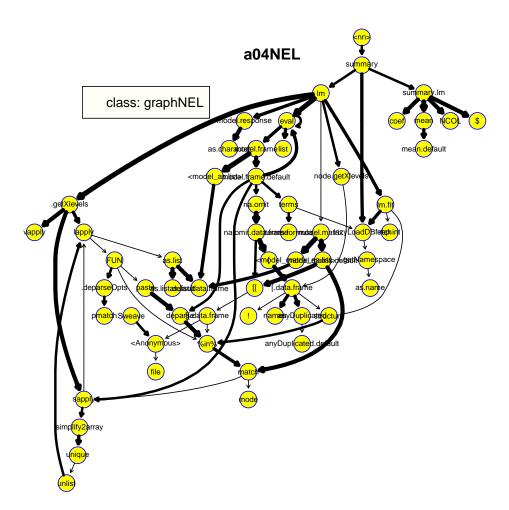


FIGURE 13. Rgraphviz/graph basic plot

Rgraphviz/graph basic plot: see fig. 13.

 $as\_graphNEL\_sprof$ 

**ToDo:** merge with To use attributes on nodes and edges, we need Rgraphviz.

```
_ Input _
library(graph)
# see help(nodeRenderInfo) for attributes.
a04NEL0 <- as_graphNEL_sprof(sprof04, weight=FALSE)
#nDD0 <- (nodeRenderInfo(a04NEL0))</pre>
#a04NEL0 <- layoutGraph(a04NEL0)</pre>
#nDD1 <- (nodeRenderInfo(a04NEL0))# fill & col corrupted</pre>
nodeDataDefaults(a04NEL0, "shape") <- "ellipse"</pre>
#nodeDataDefaults(a04NEL0, "cex") <- 1.0</pre>
#nodeDataDefaults(a04NEL0, "weigth") <- 1</pre>
#nodeDataDefaults(a04NEL0, "fill") <- "green"</pre>
#nodeDataDefaults(a04NELO, "col") <- "yellow"</pre>
nDD2 <- (nodeRenderInfo(a04NEL0))</pre>
a04NEL0 <- layoutGraph(a04NEL0)</pre>
# node attributes
#sprof04$nodes$self.time -> fill
fill_list <- heat.colors(12)[</pre>
        rkindex(-sprof04$nodes$self.time,
                pwr=0.5, maxindex=12, ties.method="min")]
names(fill_list) <- sprof04$nodes$name</pre>
#sprof04$nodes$total.time -> lwd
lwd_list <- rkindex(sprof04$nodes$total.time,</pre>
        pwr=0.5, maxindex=6, ties.method="min")
names(lwd_list) <- sprof04$nodes$name</pre>
#strx(nodeRenderInfo(a04NEL0))
nodeRenderInfo(a04NEL0) <- list(lwd=lwd_list,</pre>
        fill=fill_list,
        col="#0000FF80"
        shape="ellipse",
        weight=1)
#strx(nodeRenderInfo(a04NEL0))
# edge attributes
# strx(edgeRenderInfo(a04NEL0))
edgeRenderInfo(a04NEL0) <- list(col= "#80808080")</pre>
# strx(edgeRenderInfo(a04NEL0))
renderGraph(a04NEL0)
title(main="nodes (time)", sub=sprof04$info$id, col.sub=grey(0.5))
legend("bottomright",legend= c("node color: self.time, pwr=0.5",
        "node lwd: total.time, pwr=0.5",
        "edge lwd: frequency"),
        bty="n",
        text.font=3,
        bg="#00004040",
        seg.len=0)
        legend("topleft",
                 legend=paste0("class: ",class(a04NEL0)),
                 bg="#FFFE040",
```

seg.len=0)

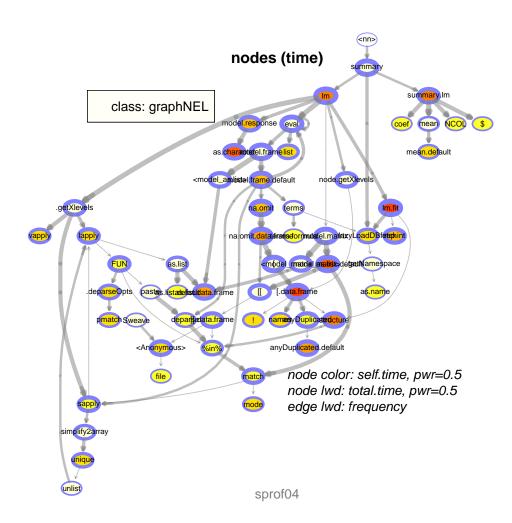


FIGURE 14. Rgraphviz/graph plot with attributes

Rgraphviz/graph plot with attributes: see fig. 14. This plot gives us the traditional view, highlighting nodes and edges with overall time presence.

Input

detach(package:Rgraphviz)

We should not overload the plot. We could use colour encoding for the edges, but this would conflict visually with the colour encoding of the nodes. We could use different colours for classes of nodes, but then we would need an additional display to explain the shape information. But within the choice of attributes used, we still can select the information shown. To close this round, instead of showing the node information from the rough summary, we can show the information from the run length discussed in section 2.4 on page 34.

```
_ Input _
amt04 <- nodesrunlength(sprof04)</pre>
library(graph)
library(Rgraphviz)
# see help(nodeRenderInfo) for attributes.
a04NELOrle <- as_graphNEL_sprof(sprof04, weight=FALSE)
#nDD0 <- (nodeRenderInfo(a04NELOrle))</pre>
#a04NELOrle <- layoutGraph(a04NELOrle)</pre>
#nDD1 <- (nodeRenderInfo(a04NELOrle))# fill & col corrupted</pre>
nodeDataDefaults(a04NELOrle, "shape") <- "ellipse"</pre>
#nodeDataDefaults(a04NELOrle, "cex") <- 1.0</pre>
#nodeDataDefaults(a04NELOrle, "weigth") <- 1</pre>
#nodeDataDefaults(a04NELOrle, "fill") <- "green"</pre>
#nodeDataDefaults(a04NELOrle, "col") <- "yellow"</pre>
nDD2 <- (nodeRenderInfo(a04NELOrle))</pre>
a04NELOrle <- layoutGraph(a04NELOrle)</pre>
# node attributes
#sprof04$nodes$self.time -> fill
fill_list <- heat.colors(12)[rkindex(-amt04[,"avg_time"], pwr=0.5, maxindex=12, ties.method="min")]
names(fill_list) <- sprof04$nodes$name</pre>
#sprof04$nodes$total.time -> lwd
lwd_list <- rkindex(amt04[,"nr_runs"], pwr=0.5,maxindex=6, ties.method="min")</pre>
names(lwd_list) <- sprof04$nodes$name</pre>
#strx(nodeRenderInfo(a04NELOrle))
nodeRenderInfo(a04NELOrle) <- list(lwd=lwd_list,</pre>
        fill=fill_list,
        col="#0000FF80"
        shape="ellipse",
        weight=1)
#strx(nodeRenderInfo(a04NELOrle))
# edge attributes
# strx(edgeRenderInfo(a04NELOrle))
edgeRenderInfo(a04NELOrle) <- list(col= "#80808080")</pre>
# strx(edgeRenderInfo(a04NELOrle))
renderGraph(a04NELOrle)
title(main="nodes (run length)", sub=sprof04$info$id, col.sub=grey(0.5))
legend("bottomright",
        legend= c("node lwd: nr runs, pwr=0.5",
         "node color: avg time, pwr=0.5",
         "edge lwd: frequency"),
                          bg="#FFFE040",
                seg.len=0,
        bty="n",
```

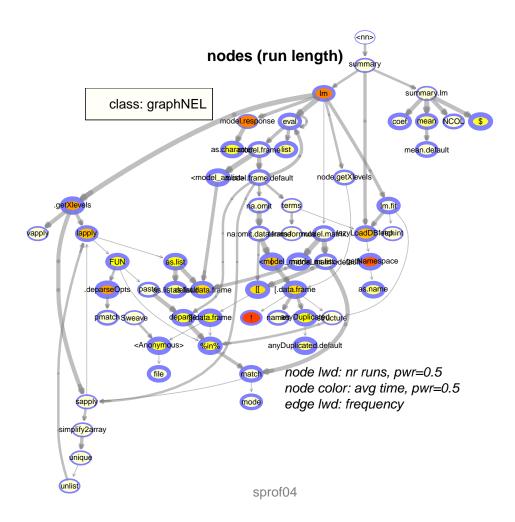


FIGURE 15. nodes (run length)

Nodes marked by run length and run count: see fig. 15. This plot gives us a more informative view, highlighting nodes by number of runs and average run length and edges with overall time presence. It puts the information in place, showing us the context where the resources are consumed. Some are consumed with good reason, since here work is done. Others are very doubtful and seem to be mere administration. These are candidates for improvement.

#### 4. Standard output

For a reference, here are the standard functions.

sprof <- sprof01	1	
4.1. <b>Print.</b> We omit the (lengt as a reference.	hy) print output here and just give th	e command
<pre>print_nodes(sprof)</pre>	Input	
print_stacks(sprof)	Input	
<pre>print_profiles(sprof)</pre>	Input	
The print() method for sprof	objects concatenates these three fund	tions.
4.2. Summary.		
summary_nodes(sprof)	Input	
summary_stacks(sprof)	Input	
summary profiles(sprof)	Input	

The summary() method for sprof objects concatenates these three functions.

ToDo: Clarify:"print prints its argument and returns it invisibly (via invisible(x))." Return the argument, or some print representation?

ToDo: is there a print=FALSE variant to postpone printing to e.g. xtable?

4.3. Plot. Examples are given in the reference manual for sprof.

plot_nodes(sprof)	Input	
plot_stacks(sprof)	Input	
plot_profiles(sprof)	Input	

The plot.sprof() method for sprof objects concatenates these three functions, see fig. 16. Using the plot functions above allows better control and will be preferred. shownodes() may be a sufficient summary, see fig. 7 on page 23.

```
#12 16
oldpar <- par(mfrow=c(3,4))
plot.sprof(sprof04)
par(oldpar)
```

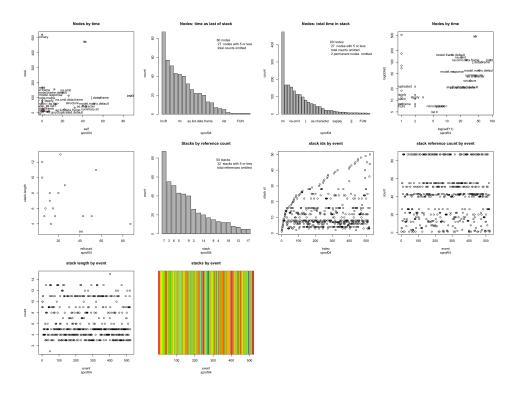


FIGURE 16. plot.sprof(sprof04)

#### 5. More Graphs

Note: This section is collecting experiments with various graph packages. It is only of interest for you, if you have a preference for some graphic package and want to look up a wrapper here. On the other side: contributions and suggestions are welcome.

Graph layout is a theme of its own. Proposals are readily available, as are their implementation. For some of them, there are R interfaces or re-implementations in R. Their usefulness in our context has to be explored, and the answers will vary with personal preferences.

For some graph layout packages we illustrate an interface here and show a sample result. We use the original profile data here. This is a nasty graph with some R stack peculiarities. The corresponding results for the trimmed profile data are shown in the next section 6 on page 77. This is a more realistic example of the kind of graphs you will have to work with.

5.1. **Example: regression.** In this section, we use the recent version of our example, *sprof02* for demonstration. You can re-run it, using your *sprof* data by modifying this instruction by replacing *sprof02* with your profile information.

\_ Input -

sprof <- sprof02

To interface **sprof** to a graph handling package, **adjacency()** can extract the adjacency matrix from the profile.

There are various packages for finding a graph layout, and the choice is open to your preferences. The R packages for most of these are just wrapper

sprofadi02 <- adjacency(sprof)

This is a format any graph package can handle (maybe). To be on the save side, we provide an (extended) edge list. The added component lwd is a proposal for the line width in the graph rendering.

Input \_\_\_\_\_ Input \_\_\_\_ sprofedge102 <- edgematrix(sprofadj02) sprofedge102\$lwd <- rkindex(sprofedge102\$count, maxindex=12, ties.method="min")

**ToDo:** by graph package: preferred input format?

ToDo: use attributes. Edge width should be easy.

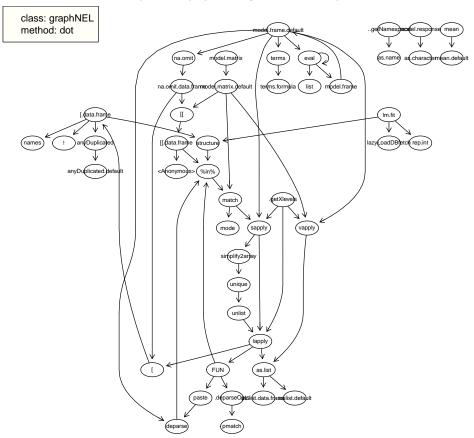
<u>ToDo</u>: include information from stack connectivity.

 $5.1.1.\ graph\ package.$  Package 'graph' was removed from the CRAN repository.

This package is now available from Bioconductor only, see http://www.bioconductor.org/packages/release/bioc/html/graph.html.

```
library(graph)
#search()
sprofadjNEL02 <- as(sprofadj02, "graphNEL")
plotviz(sprofadjNEL02, sub=sprof$info$id)
detach("package:graph")</pre>
```

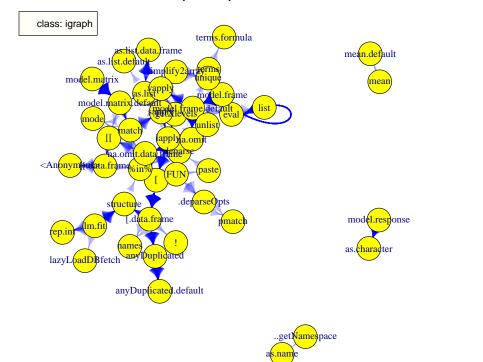
#### plotviz( sprofadjNEL02, dot )



```
5.1.2. igraph package. Attributes for igraph are documented in help(igraph.plotting).
Layouts available for igraph:
layout.auto(graph, dim=2, ...)
layout.random(graph, params, dim=2)
layout.circle(graph, params)
layout.sphere(graph, params)
layout.fruchterman.reingold(graph, ..., dim=2, params)
layout.kamada.kawai(graph, ..., dim=2, params)
layout.spring(graph, ..., params)
layout.reingold.tilford(graph, ..., params)
layout.fruchterman.reingold.grid(graph, ..., params)
layout.lgl(graph, ..., params)
layout.graphopt(graph, ..., params=list())
layout.svd(graph, d=shortest.paths(graph), ...)
layout.norm(layout, xmin = NULL, xmax = NULL, ymin = NULL, ymax = NULL,
zmin = NULL, zmax = NULL)
                                  \_ Input \_
 library(igraph)
 search()
                                  Output
 [1] ".GlobalEnv"
                            "package:igraph"
 [3] "package:sna"
                            "package:grid"
 [5] "package:wordcloud"
                            "package: RColorBrewer"
 [7] "package:Rcpp"
                            "package:xtable"
 [9] "package:sprof"
                            "package:stats"
[11] "package:graphics"
                            "package:grDevices"
[13] "package:utils"
                            "package:datasets"
[15] "package:methods"
                            "Autoloads"
[17] "package:base"
                                   Input
 as_igraph_sprof <- function(sprof, layoutfun, params=NULL,...){
         adj <- adjacency(sprof)</pre>
         adj[adj!=0] <-1
         sprof_igraph <- graph.adjacency(adj)</pre>
         sprof_igraph <- set.graph.attribute(sprof_igraph, "layout",</pre>
                 layoutfun(sprof_igraph,params=params,...),...)
         V(sprof_igraph)$color <- "yellow"</pre>
         E(sprof_igraph)$color <- "#0000FF20"</pre>
         E(sprof_igraph)$width <- c(1,2)</pre>
         return(sprof_igraph)
```

```
}
```

#### igraph kamada.kawai layout sprof02 updated



Experiments with various layouts for igraph follow.

```
_ Input _
#8 8
sprof_ig_auto <- as_igraph_sprof(sprof, layout.auto)</pre>
plot(sprof_ig_auto, main=paste0("igraph auto layout\n", sprof$info$id))
           legend("topleft",
                      legend= paste0("class: ",class(sprof_ig_auto)),
                      bg="#FFFE040",
                      seg.len=0
                                           igraph auto layout
                                            sprof02 updated
             class: igraph
                                                        <An<mark>onym</mark>ous>
                        as.l(st.default model.natrix).default .default .default [[.data.frame
                   as.list.data.frameas.list
                                    .getXlevels vapply
                                          apply
                             unlist
                                             sappl
                                            na.om(t.dan) feln rame default
                         uniqueimplify2array
                                                                             terms.formula
                                                        n<mark>a.om)</mark>t mo<mark>del.fra</mark>me
            model.response
                                       [.data.frame
                              any Duplicated
                     anyDuplicated.default
                                                                              .nan
```

<mark>name</mark>

rep.in

lazyLoadDBfetch

```
#6 6

sprof_ig_random <- as_igraph_sprof(sprof, layout.random)

plot(sprof_ig_random, main=paste0("igraph random layout\n", sprof$info$id))

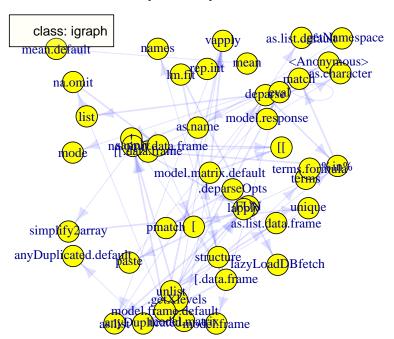
legend("topleft",

legend= paste0("class: ",class(sprof_ig_random)),

bg="#FFFFE040",

seg.len=0
)
```

## igraph random layout sprof02 updated



```
#6 6

sprof_ig_circle <- as_igraph_sprof(sprof, layout.circle)

plot(sprof_ig_circle, main=paste0("igraph circle layout\n", sprof$info$id))

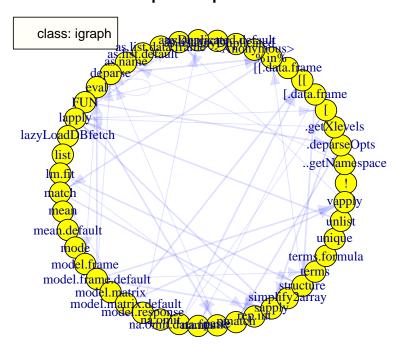
legend("topleft",

legend= paste0("class: ",class(sprof_ig_circle)),

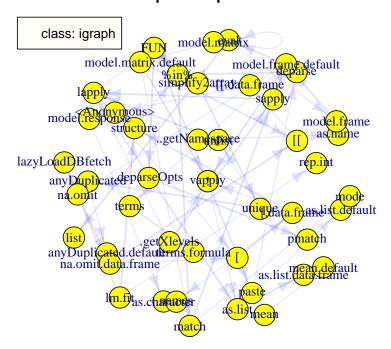
bg="#FFFFE040",

seg.len=0
)
```

## igraph circle layout sprof02 updated



#### igraph sphere layout sprof02 updated



```
#6 6

sprof_ig_auto <- as_igraph_sprof(sprof, layout.fruchterman.reingold)

plot(sprof_ig_auto, main=paste0("igraph fruchterman.reingold layout\n", sprof$info$id))

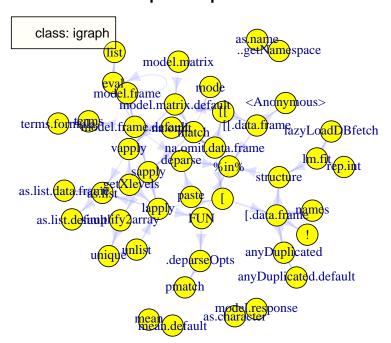
legend("topleft",

legend= paste0("class: ",class(sprof_ig_auto)),

bg="#FFFFE040",

seg.len=0
)
```

## igraph fruchterman.reingold layout sprof02 updated



```
#6 6

sprof_ig_auto <- as_igraph_sprof(sprof, layout.kamada.kawai)

plot(sprof_ig_auto, main=paste0("igraph kamada.kawai layout\n", sprof$info$id))

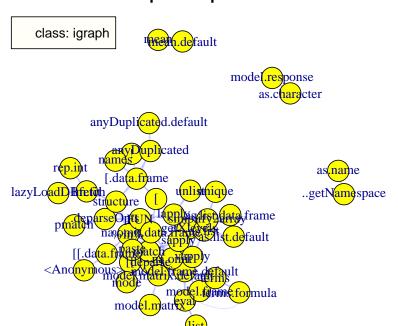
legend("topleft",

legend= paste0("class: ",class(sprof_ig_auto)),

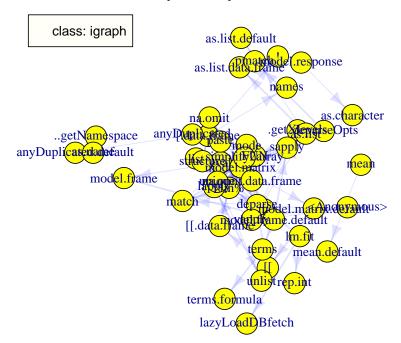
bg="#FFFFE040",

seg.len=0
)
```

#### igraph kamada.kawai layout sprof02 updated



## igraph spring layout sprof02 updated



```
#6 6

sprof_ig_auto <- as_igraph_sprof(sprof, layout.fruchterman.reingold.grid)

plot(sprof_ig_auto, main=paste0("igraph fruchterman.reingold.grid layout\n", sprof$info$id))

legend("topleft",

legend= paste0("class: ",class(sprof_ig_auto)),

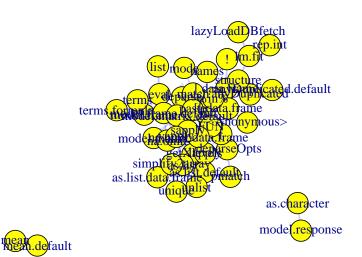
bg="#FFFFE040",

seg.len=0
)
```

## igraph fruchterman.reingold.grid layout sprof02 updated

class: igraph

..getNamespace



```
#6 6

sprof_ig_auto <- as_igraph_sprof(sprof, layout.lgl)

plot(sprof_ig_auto, main=paste0("igraph lgl layout\n", sprof$info$id))

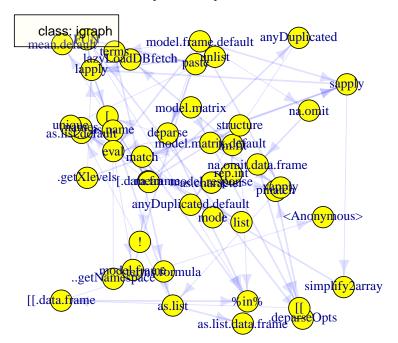
legend("topleft",

legend= paste0("class: ",class(sprof_ig_auto)),

bg="#FFFFE040",

seg.len=0
)
```

## igraph Igl layout sprof02 updated



```
#6 6

sprof_ig_auto <- as_igraph_sprof(sprof, layout.graphopt)

plot(sprof_ig_auto, main=paste0("igraph graphopt layout\n", sprof$info$id))

legend("topleft",

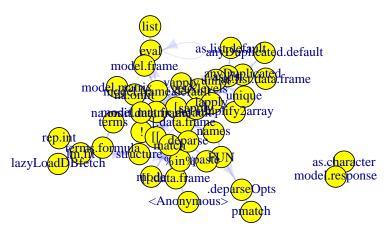
legend= paste0("class: ",class(sprof_ig_auto)),

bg="#FFFFE040",

seg.len=0
)
```

## igraph graphopt layout sprof02 updated







```
#6 6

sprof_ig_auto <- as_igraph_sprof(sprof, layout.svd)

plot(sprof_ig_auto, main=paste0("igraph svd layout\n", sprof$info$id))

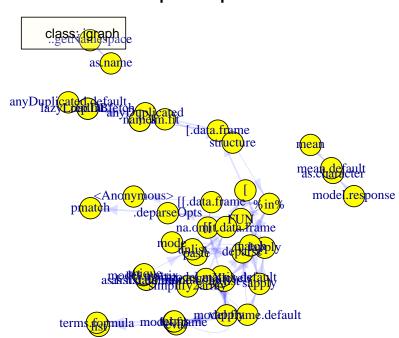
legend("topleft",

legend= paste0("class: ",class(sprof_ig_auto)),

bg="#FFFFE040",

seg.len=0
)
```

## igraph svd layout sprof02 updated



**ToDo:** propagate

5.1.3. network package.

```
detach("package:igraph")
 #detach("package:sna")
 search()
                                   _ Output
[1] ".GlobalEnv"
                             "package:sna'
 [3] "package:grid"
                              "package:wordcloud"
 [5] "package: RColorBrewer"
                             "package:Rcpp"
 [7] "package:xtable"
                             "package:sprof"
[9] "package:stats"
                              "package:graphics"
[11] "package:grDevices"
                              "package:utils"
[13] "package:datasets"
                              "package:methods"
[15] "Autoloads"
                             "package:base"
                                    _ Input _
 library(network)
 search()
                                    Output
[1] ".GlobalEnv"
                             "package:network"
 [3] "package:sna"
                             "package:grid"
 [5] "package:wordcloud"
                             "package: RColorBrewer"
 [7] "package:Rcpp"
                             "package:xtable"
 [9] "package:sprof"
                             "package:stats"
[11] "package:graphics"
                             "package:grDevices"
[13] "package:utils"
                             "package:datasets"
[15] "package:methods"
                             "Autoloads"
[17] "package:base"
as_network_sprof <- function(sprof) {
         sprofadj02 <- adjacency(sprof)</pre>
         nwsprof02 <- as.network(sprofadj02)</pre>
         network.vertex.names(nwsprof02) <-</pre>
                 rownames(sprofadj02) # not honoured by plot
         return(nwsprof02)
 }
                                     Input
plot_network_sprof <- function( nwsprof,</pre>
         mode = "fruchtermanreingold", main=NULL, label=NULL,...) {
         classnwsprof <- class(nwsprof)</pre>
         main1 <- deparse(substitute(nwsprof))</pre>
         if (!is.null(main)) main1 <- paste0(main,"\n",main1)</pre>
         if (inherits(nwsprof, "sprof")) nwsprof <-</pre>
                 as_network_sprof(nwsprof)
```

if (!is.null(label)) warning("explicit label supplied, but will use vertex names")

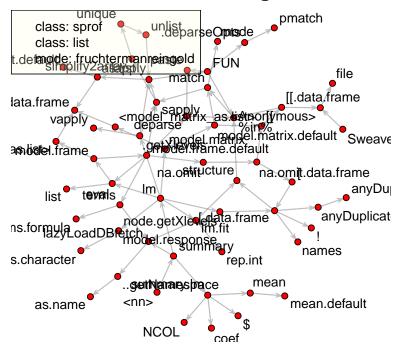
}

\_ Input -

#6 6

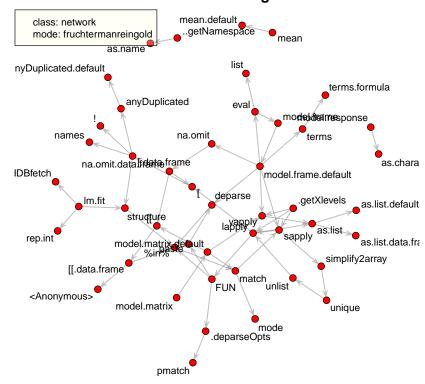
plot\_network\_sprof(sprof04)

# sprof04: network layout fruchtermanreingold



sprof04

#### nwsprof02: network layout fruchtermanreingold



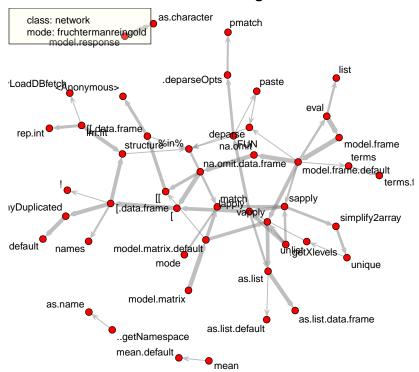
nwsprof02

Experiments to include weight.

#### ToDo: maximum edge.lwd?

```
edge.lwd<- sprofadj02
edge.lwd[edge.lwd>0]<- rkindex(edge.lwd[edge.lwd>0], maxindex=12, ties.method="min")
plot_network_sprof(nwsprof02, label=rownames(sprofadj02),
edge.lwd=edge.lwd)
detach("package:network")
```

#### nwsprof02: network layout fruchtermanreingold



nwsprof02

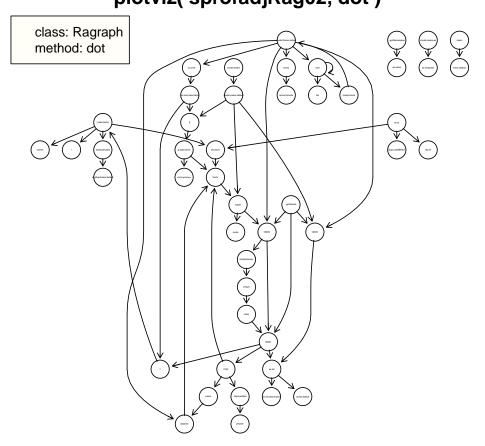
 $5.1.4.\ Rgraphviz$  package 'Rgraphviz' was removed from the CRAN repository.

This package is now available from Bioconductor only, see http://www.bioconductor.org/packages/release/bioc/html/Rgraphviz.html.

```
_ Input _
library(Rgraphviz)
search()
                                   Output
[1] ".GlobalEnv"
                            "package:Rgraphviz"
[3] "package:graph"
                            "package:sna"
[5] "package:grid"
                            "package:wordcloud"
[7] "package:RColorBrewer" "package:Rcpp"
[9] "package:xtable"
                            "package:sprof"
[11] "package:stats"
                            "package:graphics"
                            "package:utils"
[13] "package:grDevices"
                            "package:methods"
[15] "package:datasets"
[17] "Autoloads"
                            "package:base"
sprofadjRag02 <- agopen(sprofadjNEL02, name="Rprof Example")</pre>
```

# 6 6
plotviz(sprofadjRag02, sub=sprof\$info\$id)

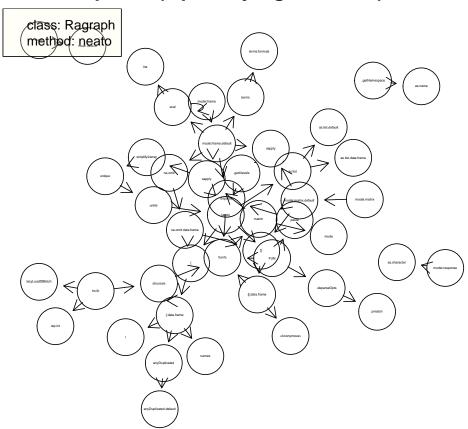
#### plotviz( sprofadjRag02, dot )



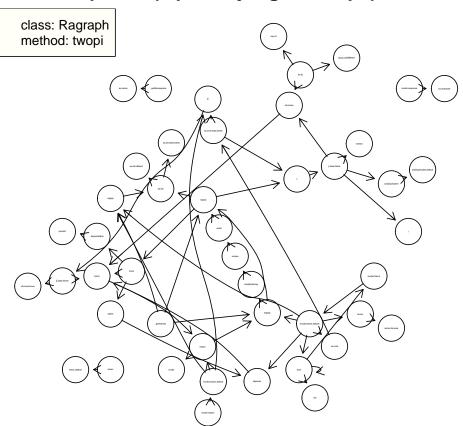
\_\_\_\_ Input \_

#6 6

### plotviz( sprofadjRag02, neato )

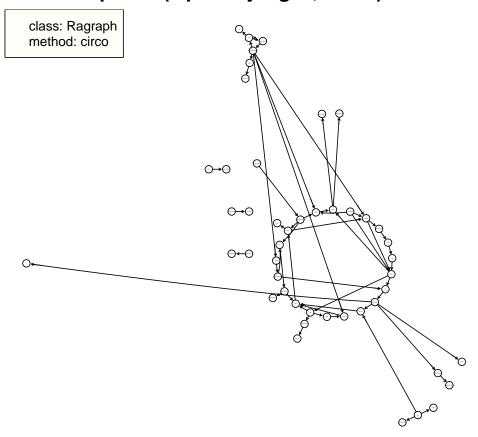


### plotviz( sprofadjRag02, twopi )

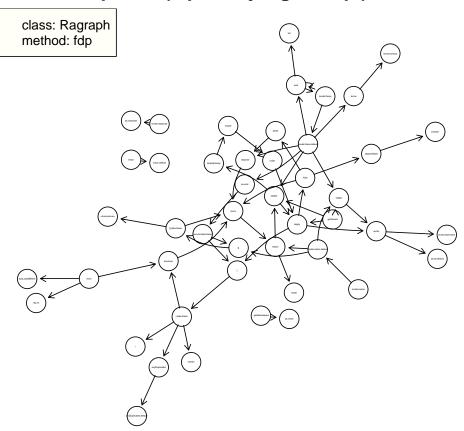


#6 6

### plotviz( sprofadjRag02, circo )



### plotviz( sprofadjRag02, fdp )



#### 6. Trimmed example: regression

In this section, we use the reduced version of our example, *sprof04* for demonstration. Except for the change of the data set, this is just a copy of the previous chapter, collecting the various layouts for easy reference.

Note: This section will be replaced by the template section. For internal use only. Skip it.

Some experiments may have found their way to this chapter. They will be expelled.

You can re-run it, using your **sprof** data by modifying this instruction:

```
sprof <- sprof04
```

To interface sprof to a graph handling package, until() can extract the adjacency matrix from the profile.

```
sprofadj02 <- adjacency(sprof)
adjname <- colnames(sprofadj02)
adjname[adjname==""] <- "<NULL>"
colnames(sprofadj02) <- adjname
rownames(sprofadj02) <- adjname
```

This is a format any graph package can handle (maybe).

6.0.5. graph package.

```
library(graph)
search()
```

```
- Output
[1] ".GlobalEnv"
                            "package:Rgraphviz"
[3] "package:graph"
                            "package:sna"
[5] "package:grid"
                            "package:wordcloud"
[7] "package:RColorBrewer" "package:Rcpp"
[9] "package:xtable"
                            "package:sprof"
[11] "package:stats"
                            "package:graphics"
[13] "package:grDevices"
                            "package:utils"
[15] "package:datasets"
                            "package:methods"
[17] "Autoloads"
                            "package:base"
```

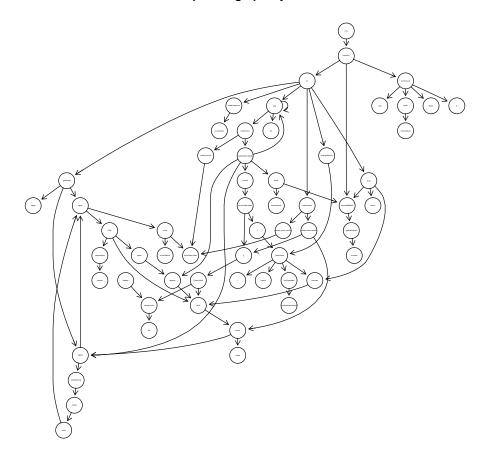
Some tests for scaling  $\dots$ 

```
sprofadjNEL02 <- as(sprofadj02,"graphNEL")
```

```
____ Input _____
```

plot(sprofadjNEL02, main="sprof04: graph layout",)
#detach("package:graph")

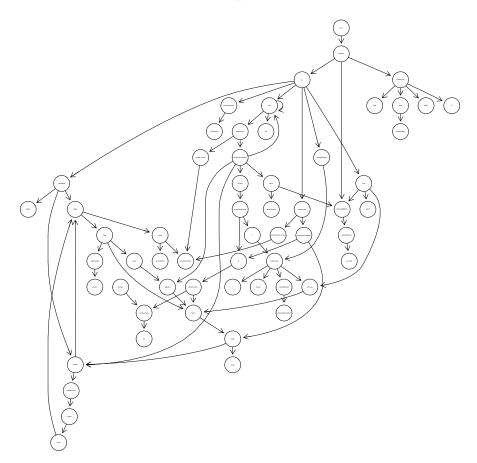
sprof04: graph layout



#18

plot(sprofadjNEL02, main="sprof04: graph layout",)
#detach("package:graph")

sprof04: graph layout

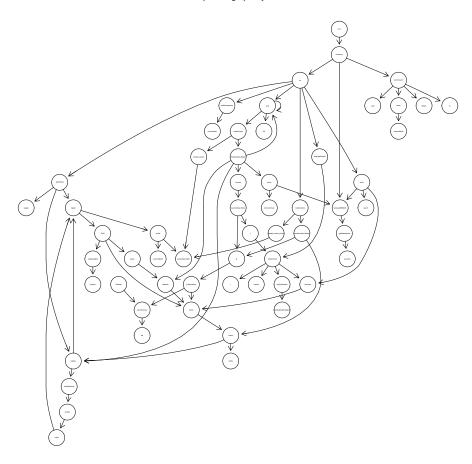


\_\_ Input \_\_\_\_\_

#12

plot(sprofadjNEL02, main="sprof04: graph layout")
#detach("package:graph")

#### sprof04: graph layout



6.0.6. igraph package.

```
_ Input _
library(igraph)
search()
```

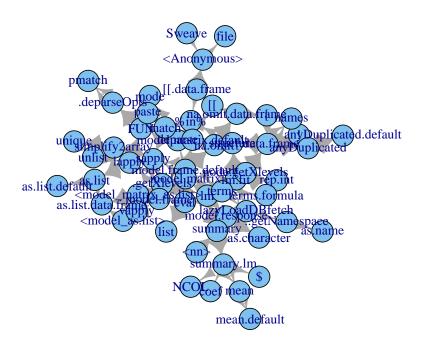
```
_____Output ___
"package:igraph"
[1] ".GlobalEnv"
 [3] "package:Rgraphviz"
                             "package:graph"
[5] "package:sna"
                             "package:grid"
[7] "package:wordcloud"
                             "package:RColorBrewer"
[9] "package:Rcpp"
                             "package:xtable"
[11] "package:sprof"
                             "package:stats"
[13] "package:graphics"
                             "package:grDevices"
[15] "package:utils"
                             "package:datasets"
[17] "package:methods"
                             "Autoloads"
[19] "package:base"
```

```
sprof_igraph <- graph.adjacency(sprofadj02)</pre>
```

```
_{-} Input _{--}
#plot(sprof_igraph, main="sprof04: igraph layout", cex.main=2)
plot(sprof_igraph, main="sprof04: igraph layout: trimmed data")
```

detach("package:igraph")

### sprof04: igraph layout: trimmed data



 $6.0.7.\ network\ package.$ 

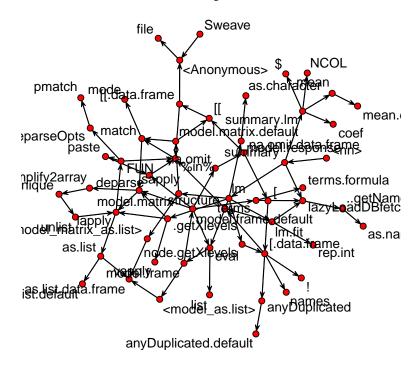
library(network)
search()

\_\_\_\_\_Output \_\_\_ "package:network" [1] ".GlobalEnv" [3] "package:Rgraphviz" "package:graph" [5] "package:sna" "package:grid" [7] "package:wordcloud" "package: RColorBrewer" [9] "package:Rcpp" "package:xtable" [11] "package:sprof"
[13] "package:graphics" "package:stats" "package:grDevices" [15] "package:utils" "package:datasets" [17] "package:methods" "Autoloads" [19] "package:base"

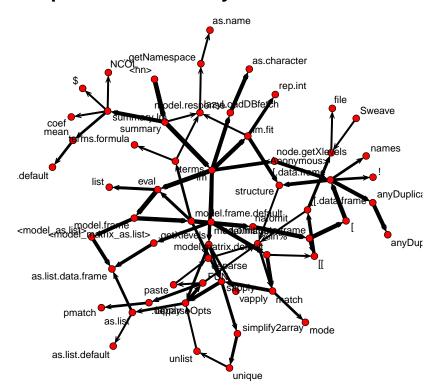
\_\_\_\_\_ Input \_\_\_\_\_

#6

# sprof04: network layout: trimmed dat

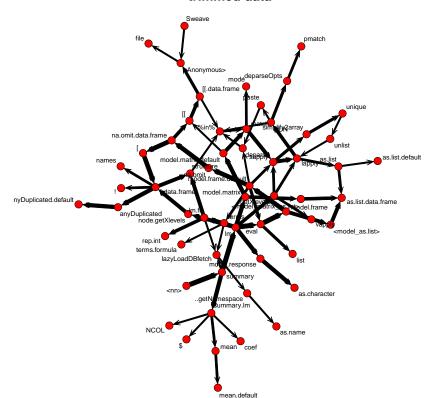


### sprof04: network layout: trimmed data



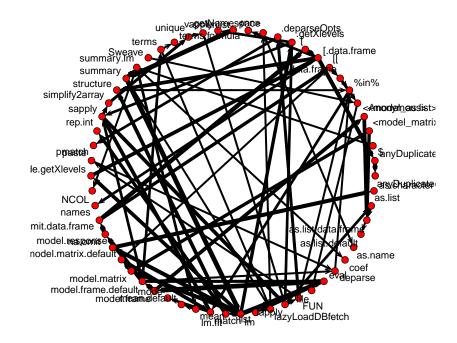
```
#12
plot(nwsprof02, label=rownames(sprofadj02),
main="sprof04: network kamadakawai layout: \n trimmed data",
mode="kamadakawai",
cex.main=2, edge.lwd=edge.lwd)
```

# sprof04: network kamadakawai layout: trimmed data



```
plot(nwsprof02, label=rownames(sprofadj02),
    main="sprof04: network circle layout: \n trimmed data",
    mode="circle",
    cex.main=2, edge.lwd=edge.lwd)
```

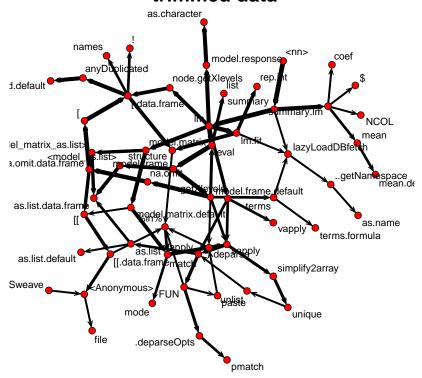
# sprof04: network circle layout: trimmed data



```
plot(nwsprof02, label=rownames(sprofadj02),
    main="sprof04: network fruchtermanreingold layout: \n trimmed data",
    mode="fruchtermanreingold",
    cex.main=2, edge.lwd=edge.lwd)

detach("package:network")
```

# sprof04: network fruchtermanreingold layout: trimmed data



6.0.8. Rgraphviz package.

library(Rgraphviz)
search()

Output

[1] ".GlobalEnv" "package:Rgraphviz"

[3] "package:graph" "package:sna"

[5] "package:grid" "package:wordcloud"

[7] "package:RColorBrewer" "package:Rcpp"

[9] "package:xtable" "package:sprof"

[11] "package:stats" "package:graphics"

[13] "package:grDevices" "package:utils"

[15] "package:datasets" "package:methods"

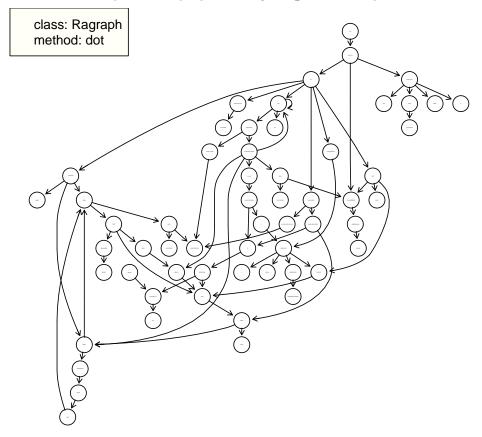
[17] "Autoloads" "package:base"

\_\_ Input \_\_\_\_\_

#6 6

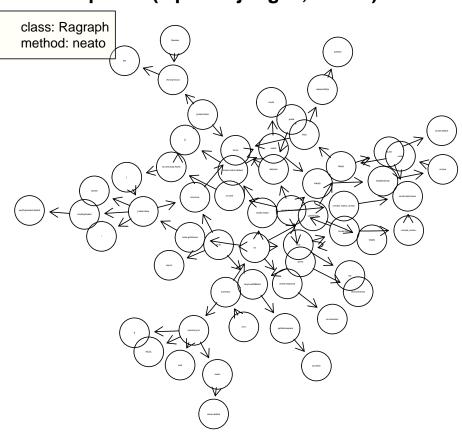
plotviz(sprofadjRag02, sub=as.character(sprof\$info\$id))

# plotviz( sprofadjRag02, dot )



#6
plotviz(x=sprofadjRag02,y="neato")

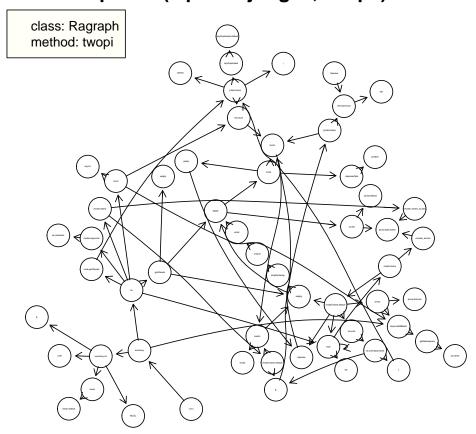
## plotviz( sprofadjRag02, neato )



#6

plotviz(sprofadjRag02,y="twopi")

### plotviz( sprofadjRag02, twopi )

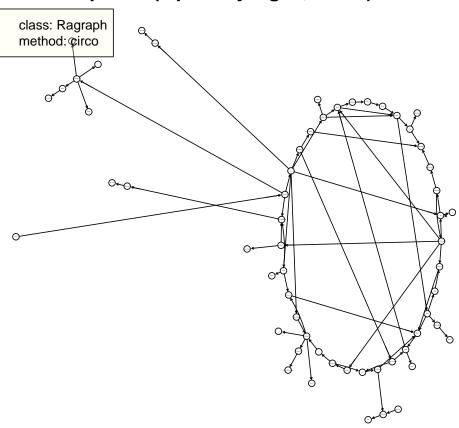


#6

\_ Input \_

plotviz(sprofadjRag02,y="circo")

## plotviz( sprofadjRag02, circo )

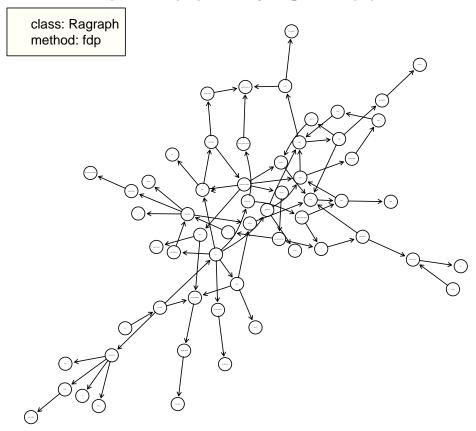


\_ Input -

#6

plotviz(sprofadjRag02, y="fdp")

### plotviz( sprofadjRag02, fdp )



#### 7. Template

- Run a profiling routine to profile your functions. You can do it on the fly
- $\bullet\,$  Read in the profile
- Get a survey
- Trim base level and burn-in/fade-out
- $\bullet$  Get a revised survey
- Use a graph display
- $\bullet$  Think!

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```
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#### R session info:

- R version 3.0.1 (2013-05-16), x86\_64-apple-darwin10.8.0
  Locale: en\_GB.UTF-8/en\_GB.UTF-8/en\_GB.UTF-8/en\_GB.UTF-8
  Base packages: base, datasets, graphics, grDevices, grid, methods, stats, utils
  Other packages: graph 1.38.3, RColorBrewer 1.0-5, Rcpp 0.10.3, Rgraphviz 2.4.0, sna 2.3-1, sprof 0.0-6, wordcloud 2.4, xtable 1.7-1
  Loaded via a namespace (and not attached): BiocGenerics 0.6.0, igraph 0.6.5-2, network 1.7.2, parallel 3.0.1, slam 0.1-28, stats4 3.0.1, tools 3.0.1

LATEX information: textwidth: 4.9823in linewidth:4.9823in

textheight: 8.0824in

#### Svn repository information:

\$Source: /u/math/j40/cvsroot/lectures/src/insider/profile/Rnw/profile.Rnw,v \$

\$Id: sprofiling.Rnw 210 2013-08-14 20:24:37Z gsawitzki \$

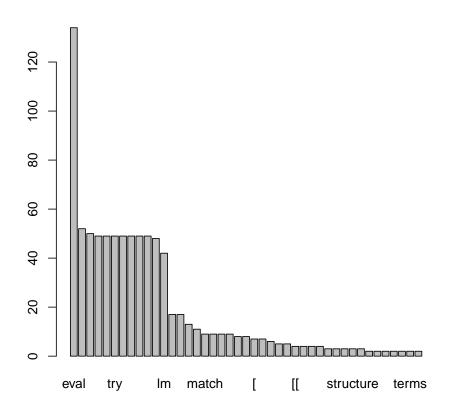
\$Revision: 210 \$

\$Date: 2013-08-14 22:24:37 0200(Wed, 14Aug2013) +

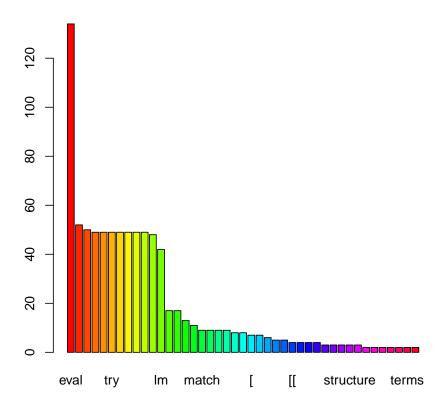
\$name: \$

\$Author: gsawitzki \$

#### 8. XXX - LOST & FOUND

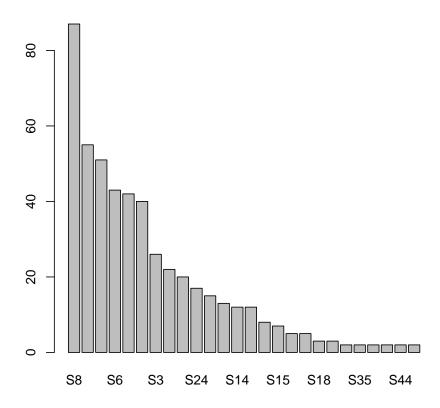


barplot(ndf[ondf], col=rainbow(length(ondf)))



Top frequent stacks.

```
x <- sprof01
xsrc <- as.matrix(x$stacks$refcount)
rownames(xsrc) <- rownames(xsrc, do.NULL=FALSE, prefix="S")
#stf <- x$stacks$refcount[x$stacks$refcount>1]
#names(stf) <- x$stacks$shortname[x$stacks$refcount>1]
stf <- xsrc[xsrc>1]
names(stf) <- rownames(xsrc)[xsrc>1]
ostf <- order(stf,decreasing=TRUE)
barplot(stf[ostf])</pre>
```



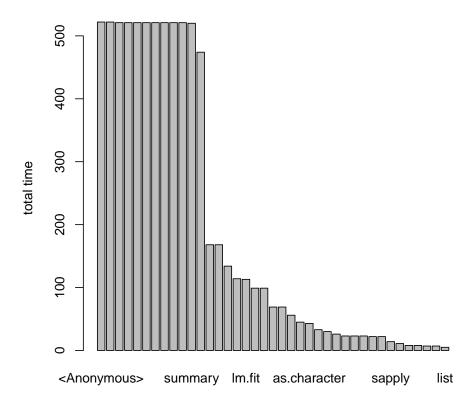
```
# xtable cannot handle posix, so we use print output here
# str(profile_nodes_rle, max.level=2, vec.len=3, nchar.max=40, list.len=6)
strx(sprof01$info)
```

```
##strx: sprof01$info
'data.frame': 1 obs. of 9 variables:
$ id : chr "sprof01"
$ date : POSIXct, format: "2013-08-15 22:02:43"
$ nrnodes : int 62
$ nrstacks : int 50
$ nrrecords : int 522
$ sample.interval: num 0.001
$ sampling.time : num 0.522
$ ctllines : Factor w/ 1 level "sample.interval=1000": 1
$ ctllinenr : num 1
```

Selections are recorded as selection vectors, with reference to the original order. This needs some caution to align them with the order choices.

```
rownames(sprof01$nodes) <- sprof01$nodes$names
nodesperm <- order(sprof01$nodes$total.time,decreasing=TRUE)
nodesnrobsok <- sprof01$nodes$total.time > 4
sp <- sprof01$nodes$total.time[nodesperm][nodesnrobsok[nodesperm]]
names(sp) <- sprof01$nodes$name[nodesperm][nodesnrobsok[nodesperm]]
barplot(sp,
main="Nodes, by total time", ylab="total time")
```

#### Nodes, by total time



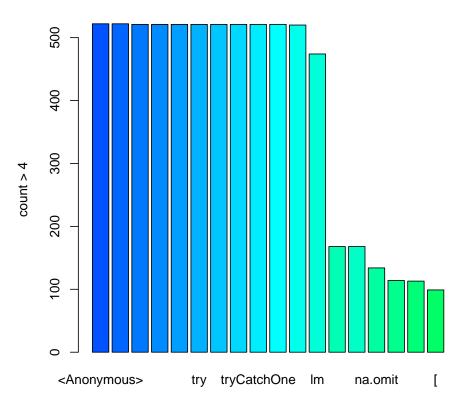
On the first look, information on the profile level is not informative. Profile records are just recordings of some step, taken at regular intervals. We get a minimal information, if we encode the stacks in colour.

We now do a step down analysis. Aggregating the information from the profiling events, we have the frequency of stack references. On the stack level, we encode the frequency in colour, and linking propagates this to the profile level.

<u>ToDo:</u> use stack colours

ToDo: use as\_rkindex

#### Stacks, by reference count (4 obs. minimum)



sprof01

```
prxt(sprof01$nodes,
caption="nodes",
label="tab:prSREnodes",
max.level=2, vec.len=3,nchar.max=40,
digits=c(0,0,0,2,0,2,0))
```

Table 10: nodes

	name	self.time	self.pct	total.time	total.pct	icol
1	!	2	0.38	2	0.03	46
2	get Name space	0	0.00	1	0.01	60
3	. departs = Opts	2	0.38	4	0.05	43
4	.get $X$ levels	0	0.00	26	0.34	28
5	[	0	0.00	99	1.29	20
6	[.data.frame	57	10.92	99	1.29	19
7	[[	0	0.00	8	0.10	36
8	[[.data.frame	1	0.19	8	0.10	35
9	$\% \mathrm{in}\%$	1	0.19	4	0.05	40
10	<anonymous></anonymous>	6	1.15	522	6.79	2
< cut >	\vdots	:	:	:	:	:
53	terms	0	0.00	2	0.03	52
54	terms.formula	1	0.19	1	0.01	59
55	try	0	0.00	521	6.78	7
56	tryCatch	0	0.00	521	6.78	8
57	tryCatchList	0	0.00	521	6.78	10
58	tryCatchOne	0	0.00	521	6.78	5
59	unique	3	0.57	4	0.05	42
60	unlist	0	0.00	1	0.01	54
61	vapply	3	0.57	23	0.30	30
62	withVisible	0	0.00	521	6.78	4

```
#str(sprof01$stacks, max.level=2, vec.len=6,
 # nchar.max=40, list.len=20,width=70, strict.width="wrap"
 strx(sprof01$stacks)
                              _____ Output _
##strx: sprof01$stacks
'data.frame': 50 obs. of 7 variables:
$ nodes :List of 50
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
..$ : int 52 10 23 55 56 57 58 21 ...
.. [list output truncated]
$ shortname : Factor w/ 50 levels
   "S<A>eFttCtCLtCOdTCwVeesleem.m..n.n...[["| __truncated__,..: 27 17
   19 1 35 36 37 30 ...
$ refcount : num 1 5 26 55 13 43 51 87 ...
$ stacklength : int 19 20 19 21 14 15 15 14 ...
$ stackheadnodes: int 52 52 52 52 52 52 52 ...
$ stackleafnodes: int 27 28 41 6 39 14 38 30 ...
$ stackssrc : Factor w/ 50 levels "! [.data.frame [
  na.omit.data.frame na."| __truncated__,..: 27 28 39 5 37 13 36 30
A summary is provided on request.
                                    Input
 sumsprof01 <- summary.sprof(sprof01)</pre>
                                 Output -
$id
[1] "Profile Summary Thu Aug 15 22:03:15 2013"
$len
[1] 522
$uniquestacks
[1] 50
$nr_runs
[1] 396
$nrstacks
Γ17 50
```

#### \$stacklength

[1] 3 25

\$nrnodesperlevel

[1] 1 1 2 1 1 1 1 1 1 1 1 1 3 10 11 9 9 15 8 7 5 7 [23] 2 1 1

shortname root leaf self.time self.pct ! - LEAF 2 0.383142 0 0.000000 ..gN ..getNamespace - LEAF 2 0.383142 .deparseOpts .dpO 0 0.000000 - -.getXlevels .gtX 0 0.000000 [ 57 10.919540 [.data.frame ſ.d. - LEAF ]] 11 0 0.000000 - LEAF [[.data.frame [[.. 1 0.191571 %in% %in% - LEAF 1 0.191571 6 1.149425 <Anonymous> <An>- LEAF 1 0.191571 - LEAF \$ - LEAF 1 0.191571 anyDuplicated anyD22 4.214559 - LEAF anyDuplicated.default anD. - LEAF 43 8.237548 as.character as.c as.list - -0.000000 as.l - LEAF 22 4.214559 as.list.data.frame a... - LEAF 1 0.191571 as.list.default as.. as.n - LEAF 1 0.191571 as.name coef coef - LEAF 1 0.191571 deparse dprs - LEAF 1 0.191571 0 0.000000 doTryCatch dTrC - -- LEAF 1 0.191571 eval eval 0 0.000000 evlF - evalFunc 1 0.191571 file - LEAF file 1 0.191571 FUN FUN - LEAF - LEAF 2 0.383142 lapply lppl lazyLoadDBfetch - LEAF 2 0.383142 1LDB list - LEAF 5 0.957854 list 42 8.045977 lm lm - LEAF lm.fit lm.f - LEAF 87 16.666667 1 0.191571 match mtch- LEAF 0 0.000000 mean mean - -- LEAF mean.default 2 0.383142 mn.d- LEAF 2 0.383142 mode mode 0 0.000000 - model.frame mdl.f model.frame.default mdl.f. - LEAF 12 2.298851 model.matrix mdl.m - -0 0.000000 - LEAF 51 9.770115 model.matrix.default mdl.m. mdl.r - LEAF 13 2.490421 model.response n.mt - LEAF 20 3.831418 na.omit na.omit.data.frame n... - LEAF 26 4.980843 nams - LEAF 2 0.383142 names NCOL - LEAF NCOL 1 0.191571 0 0.000000 paste past - LEAF 2 0.383142 pmatch pmtc

rep.int	rp.n	- LEAF	7	1.340996
sapply	sppl	- LEAF	1	0.191571
simplify2array	smp2		0	0.000000
structure	strc	- LEAF	32	6.130268
summary	smmr		0	0.000000
summary.lm	smm.	- LEAF	40	7.662835
Sweave	Swev I	ROOT -	0	0.000000
terms	trms		0	0.000000
terms.formula	trm.	- LEAF	1	0.191571
try	try		0	0.000000
tryCatch	tryC		0	0.000000
tryCatchList	trCL		0	0.000000
tryCatchOne	trCO		0	0.000000
unique	uniq	- LEAF	3	0.574713
unlist	unls		0	0.000000
vapply	vppl	- LEAF	3	0.574713
withVisible	wthV		0	0.000000
	total.time	total.pct		
!	2			
getNamespace	1	0.191571		
.deparseOpts	4			
.getXlevels	_	4.980843		
	99			
[.data.frame	99			
[[	8	1.532567		
[[.data.frame	8	1.532567		
%in%	4	0.766284		
<anonymous></anonymous>	_	100.000000		
\$	1	0.191571		
	23			
anyDuplicated		4.214559		
anyDuplicated.default as.character		8.237548		
as.list		4.406130		
	23			
as.list.data.frame				
as.list.default	1	0.191571		
as.name	1	0.191571		
coef	1	0.191571		
deparse	2			
doTryCatch	521			
eval	521	99.808429		
evalFunc		99.808429		
file	1			
FUN		1.340996		
lapply		5.747126		
${\tt lazyLoadDBfetch}$	3			
list	5			
lm	474			
lm.fit	113			
match	11			
mean		0.383142		
mean.default		0.383142		
mode	2			
model.frame	168	32.183908		

```
model.frame.default
                        168 32.183908
model.matrix
                         69 13.218391
                         69 13.218391
model.matrix.default
                         56 10.727969
model.response
                        134 25.670498
na.omit
na.omit.data.frame
                        114 21.839080
names
                          2 0.383142
NCOL
                          1 0.191571
                          1 0.191571
paste
                          2 0.383142
pmatch
                          7 1.340996
rep.int
                         14 2.681992
sapply
                          4 0.766284
simplify2array
structure
                          33
                              6.321839
summary
                        520 99.616858
summary.lm
                          45
                              8.620690
Sweave
                         522 100.000000
                         2 0.383142
terms
                          1 0.191571
terms.formula
                        521 99.808429
try
                        521 99.808429
tryCatch
                        521 99.808429
tryCatchList
tryCatchOne
                        521 99.808429
unique
                         4 0.766284
unlist
                          1 0.191571
                         23 4.406130
vapply
withVisible
                        521 99.808429
```

#str(profile\_nodes\_rle, max.level=2, vec.len=3, nchar.max=40, list.len=6)
strx(sumsprof01)

```
##strx: sumsprof01
'data.frame': 62 obs. of 7 variables:
$ shortname : Factor w/ 62 levels "!","..gN",".dp0",..: 1 2 3 4 5 6 7
8 ...
$ root : Factor w/ 2 levels "-","ROOT": 1 1 1 1 1 1 1 1 1 ...
$ leaf : Factor w/ 2 levels "-","LEAF": 2 1 2 1 1 2 1 2 ...
$ self.time : num 2 0 2 0 0 57 0 1 ...
$ self.pct : num 0.383 0 0.383 0 ...
$ total.time: num 2 1 4 26 99 99 8 8 ...
$ total.pct : num 0.383 0.192 0.766 4.981 ...
```

```
#str(sumsprof01, max.level=2, vec.len=3,
# nchar.max=40, list.len=6,
# width=70, strict.width="wrap")
```

The classical approach hides the work that has been done. Actually it breaks down the data to record items. This figure is not reported anywhere. In our case, it can be reconstructed. The profile data have 8456 words in 524 lines.

In our approach, we break down the information. Two lines of control information are split off. We have 522 lines of profile with 50 unique stacks, referencing 62 nodes. Instead of reducing it to a summary, we keep the full information. Information is always kept on its original level.

On the profiles level, we know the sample interval length, and the id of the stack recorded. On the stack level, for each stack we have a reference count, with the sample interval lengths used as weights. This reference count is added up for each node in the stack to give the node timings.

Cheap statistics are collected as the come by. For example, from the stacks table it is cheap to identify root and leaf nodes, and this mark is propagated to the nodes table. These are some attempts to recover the factor structures.

```
Input -
xfi <- levels(sprof02$nodes$name)
 profile_nodes_rlefac <- lapply(profile_nodes_rle,</pre>
         function(xl) {xl$values <- factor(xl$values,</pre>
                levels=1:62.
                labels=xfi); xl}) # seems ok
 profile_nodes_rletfac <- lapply(profile_nodes_rle,</pre>
         function(x) table(x,dnn=c("run length","node")) ) #factors lost again
         colnames(profile_nodes_rletfac[[1]]) <-</pre>
                sprof02$nodes$name[ as.integer(colnames(profile_nodes_rletfac[[1]]))]
 profile_nodes_rletfac1 <- lapply(profile_nodes_rletfac,</pre>
         function(x1) {colnames(x1) <-</pre>
                sprof02$nodes$name[ as.integer(colnames(x1))];
        x1} )
 invisible(lapply(profile_nodes_rletfac1,
 function(x) print.table(t(x),zero.print = ".") ))
                                \_ Output \_
     run length
      1 2 3 4 5 6 7
node
  <NA> 17 1 1 1 .
  <NA> 1 . . . .
  <NA> 1
  <NA> 40 17 7 4 2 6 1
  <NA> 46 18 3 1 1 1 1
  <NA> 1 .
             2 . .
  <NA> 55 4
  <NA> 35 3 3 . . 1
  <NA> 1
                     run length
                      1 2 3 4 5 6 7
node
  as.character
                      34
                          3 1
  as.name
                      1
                      40 17 7 4 2 6 1
  eval
                      16 .
  lapply
                                1
  lazyLoadDBfetch
                      1
  mean.default
                      1
  model.matrix.default 55 4 2
```

```
rep.int
                  7 . . . . . .
 sapply
                  3 .
                 10 1 . . . 1 .
 structure
 vapply
                 . . 1
           run length
node
            1 2 3 4 5 6 7
 [
            14 . . 1 . .
 ]]
             1 . . . .
            1 . . . .
 %in%
 as.list
             2 . . . . .
             2 . . . .
 lapply
 match
             9.
 model.frame 40 17 7 4 2 6 1
 simplify2array 1 . . . . .
 vapply 6 1 . .
                run length
                 1 2 3 4 5 6 7
node
                      . 1
 [.data.frame
                14
 [[.data.frame
                 1
                 7 1 .
 as.list
 as.list.data.frame 2
 FUN
                 1
 match
                 1
 model.frame.default 40 17 7 4 2 6 1
         9 . . . .
 sapply
                 1 . .
 unique
               run length
node
                1 2 3 4 5 6 7
 .deparseOpts
                1 . . . . .
 <Anonymous>
                1 . . . . .
 anyDuplicated
                1 . . . .
 as.list.data.frame 6 1 . .
 as.list.default 1 . . .
 deparse
                1 . . .
 eval
                 3
 lapply
                5
               46 10 5 4 1
 na.omit
 sapply
                2 .
 simplify2array
                3
                11 .
                     . 1
 structure
 terms
                1 .
 unlist
                1
                7 . . .
 vapply
               run length
                1 2 3 4 5 6
node
 as.list
 eval
 FUN
                5 . . .
                 3 .
 na.omit.data.frame 43 7 4 5 . 4
 terms.formula 1 . . . . .
                 3 . . . . .
 unique
               run length
```

```
node
                   1 2 3 4 5 6
                  3 .
  .deparseOpts
                  45 4 3 3 .
 [
 ]]
                  2 . .
                           . 1
 %in%
                  1 .
 as.list.data.frame 7 . . . 1
 FUN
                   1 . . . .
                   3 .
 list
           run length
           1 2 3 4 5 6
node
 [.data.frame 45 4 3 3 . 1
 [[.data.frame 2 . . . 1 .
 match 1 . . . .
 paste
              1 . . . . .
              2 . .
 pmatch
           run length
node
             1 4 5
             1 . .
 %in%
             1 . .
 <Anonymous> . . 1
 anyDuplicated 9 2 1
 deparse 1 . .
 mode
             1 . .
             2 . .
 names
                   run length
node
                    1 4 5
 %in%
                    1 . .
 anyDuplicated.default 9 2 1
    run length
node
     1
 match 1
     run length
node 1
 mode 1
GÜNTHER SAWITZKI
STATLAB HEIDELBERG
IM NEUENHEIMER FELD 294
D 69120 Heidelberg
E	ext{-}mail\ address: gs@statlab.uni-heidelberg.de}
```

URL: http://sintro.r-forge.r-project.org/