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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Private Version

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Profiling facilities in R

For the impatient: table 1 on page 9 and table 2 on page 11 give you the conventional information from profiling 100 *runs of a simple linear regression. You get a different view of the same information by fig. 18 on page 53. The additional information we derive is summarised in table 9 on page 45 and illustrated in fig. 19 on page 55. 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")
```

LaTeX 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.

To make sure we do not depend on packages collected along the way, we clean up packages. Calls to this funtion will be hidden.

Input

```
cleanpackages <- function()
{
        # make sure we do not get inherited methods or garbage here.
        try(detach("package:sna"), silent=TRUE)
        try(detach("package:igraph"), silent=TRUE)
        try(detach("package:network"), silent=TRUE)
        try(detach("package:graph"), silent=TRUE)
        try(detach("package:Rgraphviz"), silent=TRUE)
        try(detach("package:graph"), silent=TRUE)
}
```

The main library we are going to use is ***sprof***.

Input

```
library(sprof)
```

We want immediate warnings, if necessary. So we set **warn** to level 1. Set **warn** to level 2 if you want to handle warnings as error.

Input

```
message("switching  options(warn=1) -- immediate warning on")
options(warn=1)
```

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

Input

```
#options(error = recover)
```

Print parameters used here:

Input

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

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

Input

```
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"))
```

```
        str(x, max.level=max.level,
                vec.len=vec.len,
                nchar.max=nchar.max,
                list.len=list.len,
                width=width, strict.width=strict.width,...)
}
```

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

Input

```
xcutrows <- function(df,  cut, margin=5)
{
        if (!is.data.frame(df)) return(df)
        nrow <- nrow(df)
        # cut a range if it is not empty.
        # Quiet noop else.
        # Does not cut single lines.
         cutrng <- function(cutfrom,cutto, cutname="<cut>"){
                 if (cutfrom<cutto){
                df[cutfrom,] <- NA
                if (!is.null(rownames(df))) rownames(df)[cutfrom] <- cutname
                if (!is.null(df$name)) df$name[cutfrom] <- ""

                cutfrom <- cutfrom+1
                df[-(cutfrom:cutto),]
                }#if
        }

        if (!missing(cut))  {
                if (length(cut) ==2)
                {df <- cutrng(cut[1],cut[2])
                        return(df)
                } else {
                warning("cut ignored. Cut range must be a vector of length 2")
                        return(df)}
        }
        if (length(margin)==0) return(df)
        if (length(margin)==1) margin <-  c(margin, margin)
        if (length(margin)==2) {
                cut <- c(margin[1]+1,nrow-margin[2])
                df <-cutrng(cutfrom=cut[1],cutto=cut[2]);
                return(df)
        }

        #len=3: margin low,  center, margin up
        delta <- (nrow-sum(margin))  %/% 2
        if (delta<2) return(df)
        c1 <- margin[1]+1
        c2 <- c1 + delta -1
        c4 <- nrow-  margin[3]
        c3 <- c4 - delta
        df <- cutrng(c3,c4, cutname="<cut hi>")
        df <- cutrng(c1,c2, cutname="<cut low>")
```

```
        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,
        label=NULL, zero.print=NULL, print.results=TRUE,...)  {

        if (cut) {margin <- 10
        if (nrow(x)> 2*margin+3) x <-xcutrows(x, margin=margin)}

#special sanitising for xtable
xr <- rownames(x)
#for (i in (1:length(xr))) {xr[i] <- sub(xr[i],"\\[","$[$", fixed=TRUE)}
 #xr <- paste("$",xr,"$")
xr <- gsub("[","{[}",xr, fixed=TRUE)
xr <- gsub("_","\\_",xr, fixed=TRUE)
xr <- gsub("^","\\^",xr, fixed=TRUE)
rownames(x)<- xr

        pr <- print(
                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)

                if(!is.null(zero.print))
                        pr <- gsub( " 0 ",zero.print, x=pr, fixed=TRUE)


                if (print.results) cat(pr)
                invisible(pr)
        }
```

This is to be used with `<<print=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.

Input

```
plotviz <- function(x, layout="dot", main=NULL, sub=NULL,...)
{
        xid <- deparse(substitute(x))
        xsubid <- NULL
        class1 <- NULL
        if (inherits(x,"sprof")) {
                class1 <- paste0("class orig: ", paste(class(x), collapse=" "))
                xsubid <- x$info$id
                sub <- paste0(sub, " ", x$info$id)
                x <- as(adjacency(x), "graphNEL")
        }

        if (!is.null(sub)) sub <- as.character(sub)
        if (is.null(main))
                main <-paste0("plotviz( ", xid ,", ", layout, " )\n", xsubid) else
                main=paste0(main, "\n plotviz( ", xid ,", ", layout, " )\n", xsubid)

        if (inherits(x,"Ragraph")) {
        plot(x=x,  y=layout,
                        cex.main=1.2,
                main=main,sub=sub,...)
        } else {
                plot(x=x,  y=layout,
                attrs=list(
                        node=list(cex=4, fontsize=40,  shape="ellipse")),
                cex.main=1.2,
                main=main,sub=sub,...)
        }

#        title(sub = paste0(sub, " ", as.character(class(x)))        )
                title(sub = "xx try sub xx")
        legend("topleft",


                legend=c(class1,
                        paste0("class: ", paste(class(x), collapse=" ")),
                        paste0("layout: ",layout)),
                bg="#FFFFE040",
                seg.len=0
                )#"lightyellow" =#FFFFE0
}
```

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 *gprof* (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 *apply()* 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 *summaryRprof()* 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: rearrange stacks? detect order?

In this presentation, we will use a small list of examples. Since *Rprof* 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.

Input

```
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)
```

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.

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

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 12 how to use *sampleRprof* in package *sprof* for an easier solution.

Input

```
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.

Input

```
sumRprofRegressionExpl <- summaryRprof("RprofsRegressionExpl01.out")
#str(profile_nodes_rle, max.level=2, vec.len=3, nchar.max=40, list.len=6)
strx(sumRprofRegressionExpl)
```

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
   0.026 ...
..$ 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.

Input

```
prxt(sumRprofRegressionExpl$by.self,
        caption="summaryRprof result: by.self as final stack entry, all records",
        label="tab:prSRREbs")
```

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
”$”	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

Input

ToDo: improve barplot_s. Allow vars from matrix or data frame, keep names. Use horizontal names for horizontal layout.

ToDo: fix colour. Select colours >after< sorting. Explicit override.

```
s  <- sumRprofRegressionExpl$by.self$self.time
names(s) <-  rownames(sumRprofRegressionExpl$by.self)
barplot_s(s, horiz=TRUE, col=rainbow(length(s)), las=1, main="self.time, by self")
```

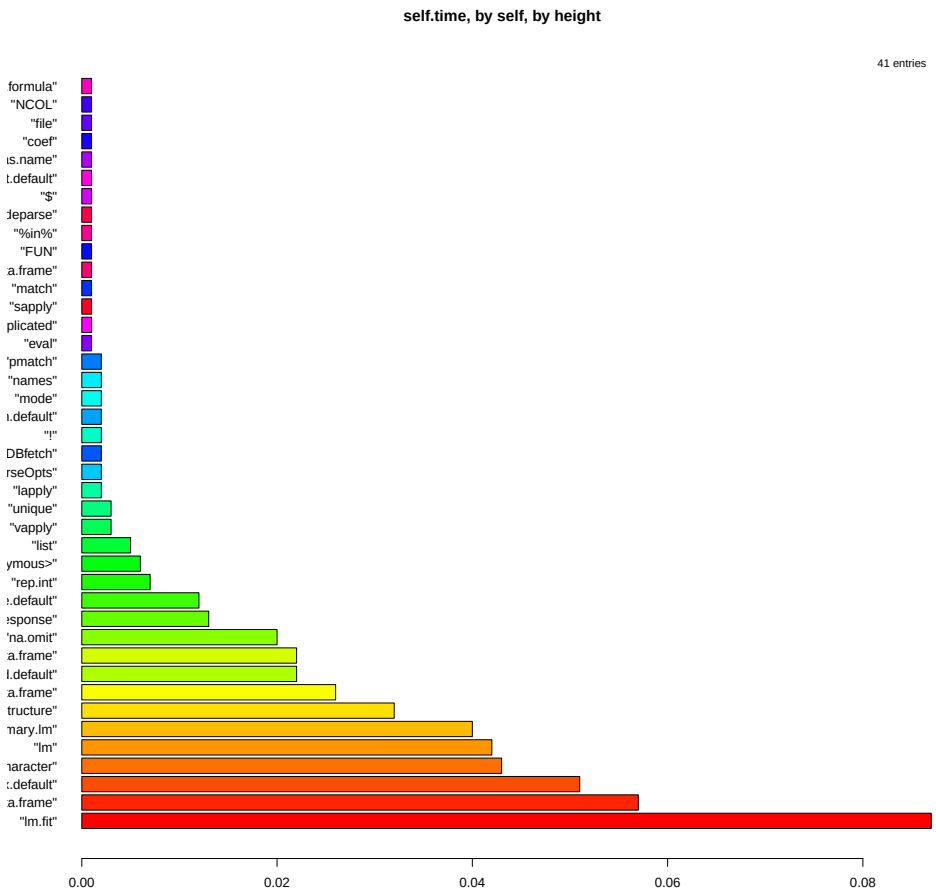


FIGURE 1. Nodes by by self.

Input

```
prxt(
        sumRprofRegressionExpl$by.total[
                sumRprofRegressionExpl$by.total$total.time>0.004,],
        caption="summaryRprof result: by.total, total time > 4ms",
        label="tab:prSRREbt")
```

Table 2: summaryRprof result: by.total, total time > 4ms

	total.time	total.pct	self.time	self.pct

"<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

Input

```
s  <- sumRprofRegressionExpl$by.total$total.time
names(s) <-  rownames(sumRprofRegressionExpl$by.total)
barplot_s(s, horiz=TRUE,
        col=rainbow(length(s)), las=1, main="total.time, by total")
```

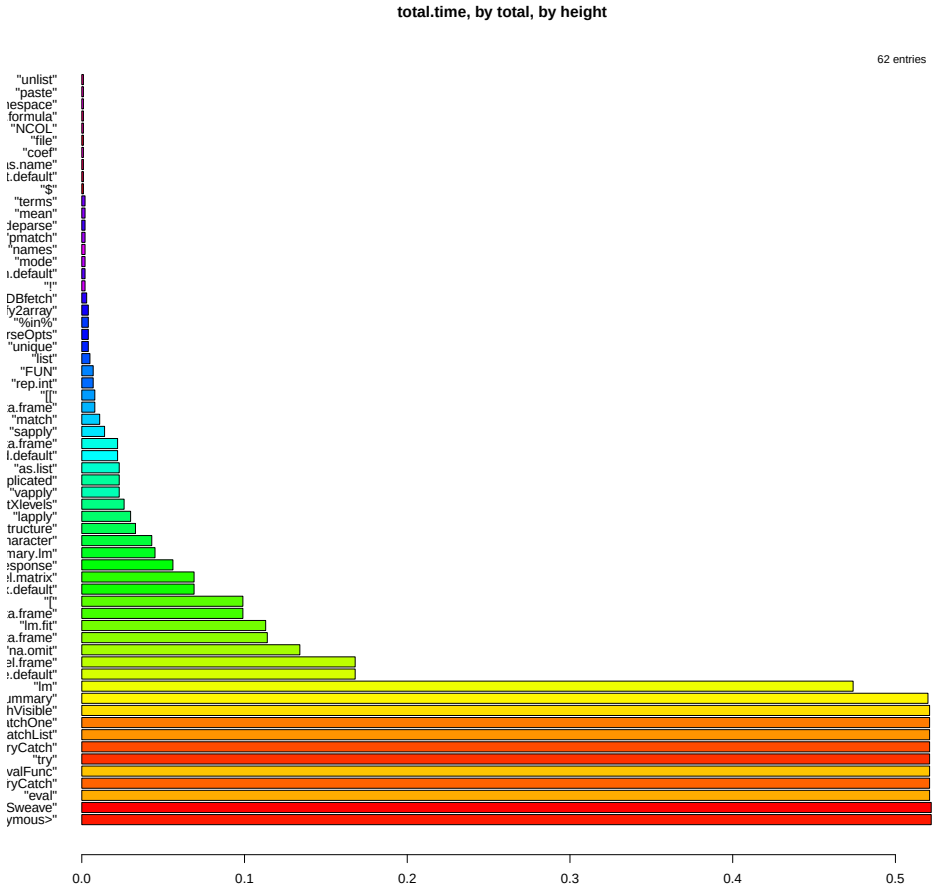


FIGURE 2. Nodes by by total.

1.1.2. *Package sprof.* In contrast to the common R profiling 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:
 $ id              : chr "RprofsRegressionExpl01.out 2013-06-13 23:46:04"
```

```
$ date           : POSIXct, format: "2013-09-10 18:00:07"
$ nrnodes        : int 62
$ nrstacks       : int 50
$ nrrecords      : int 522
$ sample.interval: num 0.001
$ sampling.time  : num 0.522
$ ctllines       : chr "sample.interval=1000"
$ ctllinenr      : num 1
```

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

Input

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

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

Input

```
save(sprof01, file="sprof01lm.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

Input

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

The basic data structure consists of four data frames. The *info* section collects global information from the input file, such as an identification strings and various global matrix. The *nodes* section initially gives the same information marginal information as *summaryRprof*. The *stacks* section puts the node information into their calling context as found in the input profile file. The *profiles* 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: variable sampling intervals will not be supported in future versions

Input

```
strx(sprof01)
```

Output

```
##strx: sprof01
List of 4
$ info :'data.frame': 1 obs. of 9 variables:
..$ id : chr "sprof01"
..$ date : POSIXct[1:1], format: "2013-09-10 18:00:07"
..$ nrnodes : int 62
..$ nrstacks : int 50
..$ nrrecords : int 522
..$ sample.interval: num 0.001
..$ sampling.time : num 0.522
..$ ctllines : chr "sample.interval=1000"
..$ ctllinenr : num 1
```

```
$ nodes :'data.frame': 62 obs. of 7 variables:
..$ name : chr [1:62] "!" "..getNamespace" ".deparseOpts" ...
..$ 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 ...
..$ nr_runs : num [1:62] 2 1 4 20 72 72 4 4 ...
..$ avg_time : num [1:62] 1 1 1 1.3 ...
$ stacks :'data.frame': 50 obs. of 6 variables:
..$ nodes :List of 50
.. .. [list output truncated]
..$ 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 52 ...
..$ stackleafnodes: int [1:50] 27 28 41 6 39 14 38 30 ...
..$ stackssrc : chr [1:50] "lazyLoadDBfetch terms model.frame.defau"|
   __truncated__ "list eval eval model.frame.default mode"|
   __truncated__ "na.omit.data.frame na.omit model.frame."|
   __truncated__ ...
$ profiles:List of 5
..$ data : int [1:522] 1 2 2 3 4 4 5 5 ...
..$ mem : NULL
..$ malloc : NULL
..$ timesRLE :List of 2
.. ..- attr(*, "class")= chr "rle"
..$ stringsAsFactors: logi FALSE
- attr(*, "class")= chr [1:2] "sprof" "list"
```

Input

```
str(sprof01$nodes)
```

Output

```
'data.frame':        62 obs. of  7 variables:
 $ name      : chr  "!" "..getNamespace" ".deparseOpts" ".getXlevels" ...
 $ self.time : num  2 0 2 0 0 57 0 1 1 6 ...
 $ self.pct  : num  0.38 0 0.38 0 0 ...
 $ total.time: num  2 1 4 26 99 99 8 8 4 522 ...
 $ total.pct : num  0.03 0.01 0.05 0.34 1.29 1.29 0.1 0.1 0.05 6.79 ...
 $ nr_runs   : num  2 1 4 20 72 72 4 4 4 3 ...
 $ avg_time  : num  1 1 1 1.3 1.38 ...
```

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 16 uses a permutation by total time, and a selection (compare to table 2 on page 11). One difference is that *sprof* uses an event count as base, usually sampled by milliseconds (ms), whereas R in general uses seconds as a base. Another difference is in two additional variables provided by *sprof*, the number of runs *nr_runs* and the average run length *avg_time*. These are discussed in section 2.4 on page 41.

Input

```
nodes <- sprof01$nodes[order(sprof01$nodes$self.time,
decreasing=TRUE),]
rownames(nodes) <- NULL
prxt(nodes[nodes$self.time>4,1:7],
digits=c(0,0,0,2,0,2,0,2),
caption="sprof result: by.self, self time > 4ms",
        label="tab:prspbtself")
```

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

	name	self.time	self.pct	total.time	total.pct	nr_runs	avg_time
1	lm.fit	87	16.67	113	1.47	71	1.59
2	[.data.frame	57	10.92	99	1.29	72	1.38
3	model.matrix.default	51	9.77	69	0.90	61	1.13
4	as.character	43	8.24	43	0.56	38	1.13
5	lm	42	8.05	474	6.16	30	15.80
6	summary.lm	40	7.66	45	0.59	29	1.55
7	structure	32	6.13	33	0.43	24	1.38
8	na.omit.data.frame	26	4.98	114	1.48	63	1.81
9	anyDuplicated.default	22	4.21	22	0.29	12	1.83
10	as.list.data.frame	22	4.21	22	0.29	17	1.29
11	na.omit	20	3.83	134	1.74	71	1.89
12	model.response	13	2.49	56	0.73	42	1.33
13	model.frame.default	12	2.30	168	2.18	77	2.18
14	rep.int	7	1.34	7	0.09	7	1.00
15	<Anonymous>	6	1.15	522	6.79	3	176.00
16	list	5	0.96	5	0.07	4	1.25

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.

Input

```
nodes <- sprof01$nodes[order(sprof01$nodes$total.time,
        decreasing=TRUE),]
rownames(nodes) <- NULL
prxt(nodes[nodes$total.time>4,1:7],
        digits=c(0,0,0,2,0,2,0,2),
        caption="sprof result: by.total, total time > 4ms",
        label="tab:prspbt")
```

Table 4: sprof result: by.total, total time > 4ms

	name	self.time	self.pct	total.time	total.pct	nr_runs	avg_time
1	<Anonymous>	6	1.15	522	6.79	3	176.00
2	Sweave	0	0.00	522	6.79	1	522.00

3	doTryCatch	0	0.00	521	6.78	1	521.00
4	eval	1	0.19	521	6.78	164	8.46
5	evalFunc	0	0.00	521	6.78	1	521.00
6	try	0	0.00	521	6.78	1	521.00
7	tryCatch	0	0.00	521	6.78	1	521.00
8	tryCatchList	0	0.00	521	6.78	1	521.00
9	tryCatchOne	0	0.00	521	6.78	1	521.00
10	withVisible	0	0.00	521	6.78	1	521.00
<cut>		⋮	⋮	⋮	⋮	⋮	⋮
30	vapply	3	0.57	23	0.30	16	1.44
31	anyDuplicated.default	22	4.21	22	0.29	12	1.83
32	as.list.data.frame	22	4.21	22	0.29	17	1.29
33	sapply	1	0.19	14	0.18	14	1.00
34	match	1	0.19	11	0.14	12	1.00
35	[[	0	0.00	8	0.10	4	2.00
36	[[.data.frame	1	0.19	8	0.10	4	2.00
37	FUN	1	0.19	7	0.09	7	1.00
38	rep.int	7	1.34	7	0.09	7	1.00
39	list	5	0.96	5	0.07	4	1.25

ToDo: remove text vdots from string/-name columns

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.

Input

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

Basic information on node level: see fig. 3 on the next 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 (See Basic information on node level - horizontal bars: fig. 4 on page 18).

Input

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

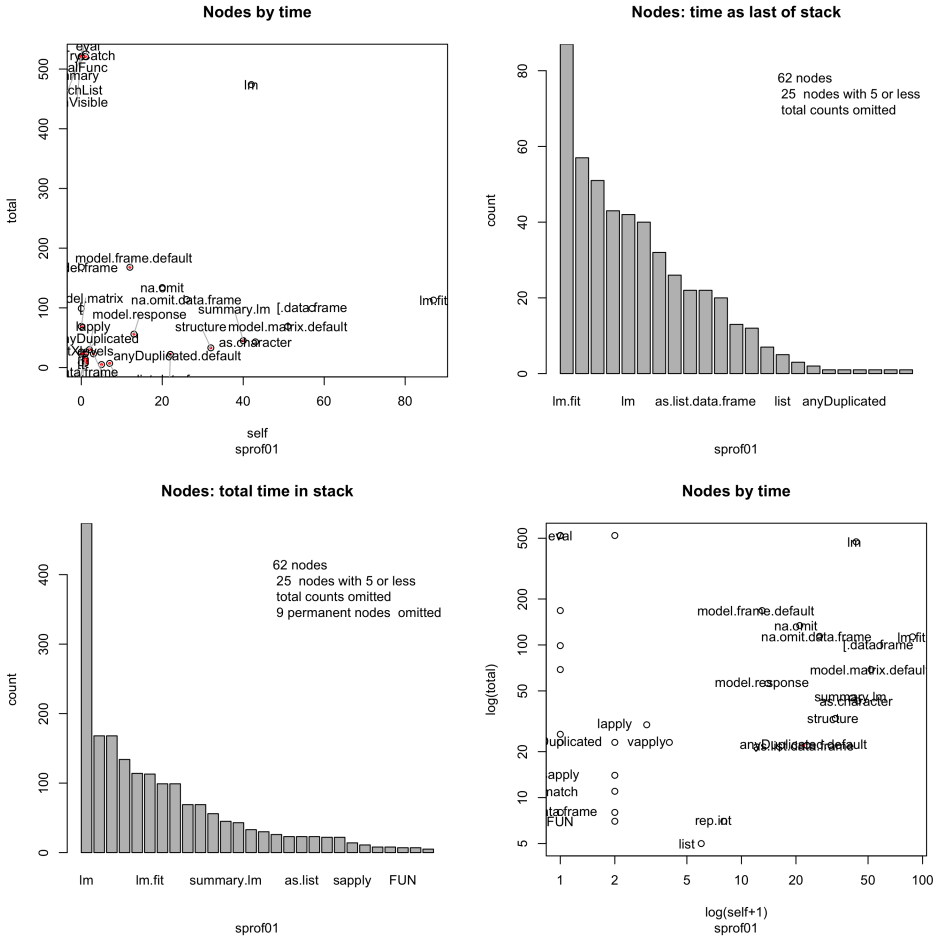


FIGURE 3. Basic information on node level

```
plot_nodes(sprof01, horiz=TRUE)
par(oldpar)
```

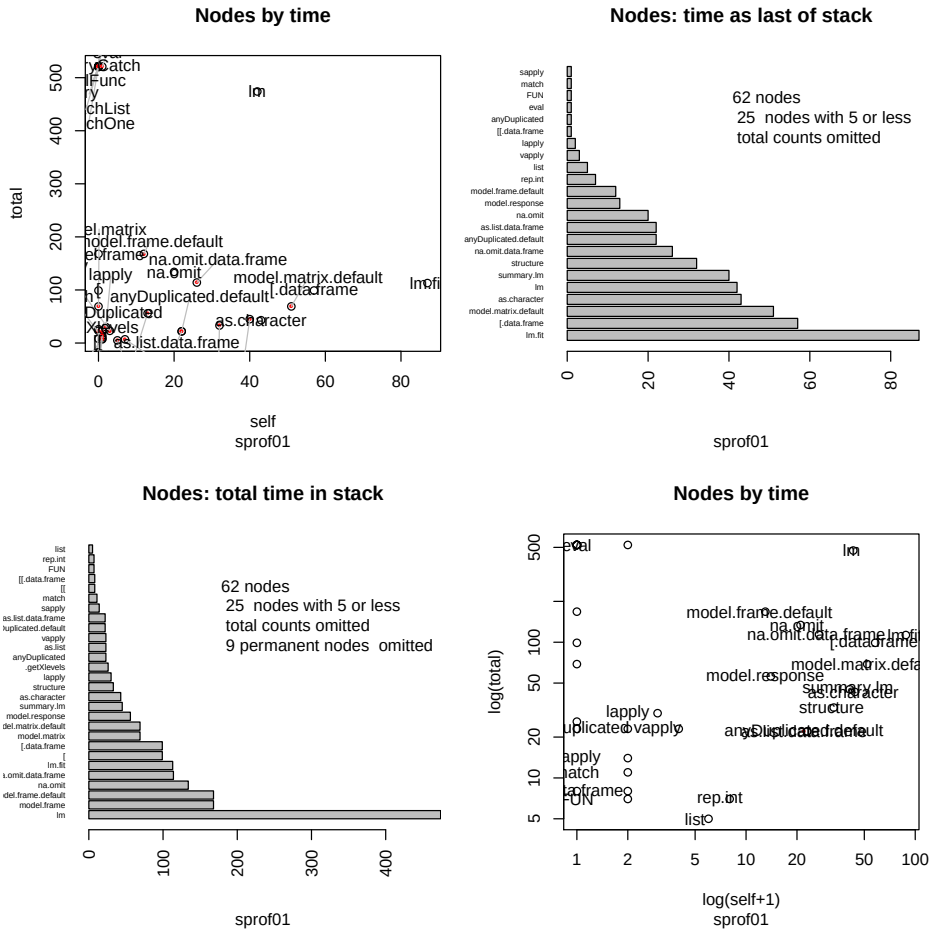


FIGURE 4. Basic information on node level - horizontal bars

1.2. **Node attributes.** The node information is kept as a *`data.frame`*, and this can be extended as long as the overall structure is preserved. Two variables, *`icol`* and *`ilwd`* are by convention used to hold graphical attributes. The semantics is not fixed, but up to your choice. *`rkindex()`* is provided to derive an index from some variable. If a display may have colour information, *`icol`* is to be used as an index into a colour table. If a display has a linewidth attribute, *`ilwd`* should be used to set the line width where feasible. The details are left to the plots.

Keeping in mind the limits of perception, *`icol`* and *`ilwd`* should be kept small. Using 4 colours or hues seems a save choice; aboout 16 colours seem to be the maximum useful in many diplays. For line width, about 3 is a save number, and about 7 seems to be the limit.

For two common models, attributes by time and attributes by run length information, a helper routine is provided to calculate indices.

ToDo: make call by reference function

Input

```
nodeattrs <- function(sprof,
        nodeattrs, # = c("default", "time", "runs"),
        maxcol=16, colpwr=0.5,
        maxlwd=7, lwdpow=0.5)
{
        #default
        if (missing(nodeattrs)) nodeattrs <- "default"
        icol <- rep(1,length(sprof$nodes$name))
        ilwd <- icol

        if (nodeattrs == "runs"){
                # may be skipped. if (is.null(sprof$nodes$avg_time)){
                nrl <- nodesrunlength(sprof, clean=FALSE)
                icol <-  rkindex(-nrl[,"avg_time"],
                                pwr=colpwr, maxindex=maxcol, ties.method="min")
                ilwd <-  rkindex(nrl[,"nr_runs"],
                        pwr=lwdpow, maxindex=maxlwd, ties.method="min")
        } else if (nodeattrs == "time"){
                icol <- rkindex(-sprof$nodes$self.time,
                                pwr=colpwr, maxindex=maxcol, ties.method="min")
                ilwd <-  rkindex(sprof$nodes$total.time,
                        pwr=lwdpow, maxindex=maxlwd, ties.method="min")
        }
        data.frame(icol=icol, ilwd=ilwd)
}
```

ToDo: propagate nodeattrs

ToDo: apply colour to selection?

So 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.

Note: we chose a colour palette for the full data set. If we restrict a display to a selection, only a fraction of the palette may be visible. As an alternative, we could adjust the palette to focus on your selection. However the recommendation is not to spend too much energy here. Instead of modifying the display, the recommended way is to extract the core information by trimming the data set (see section 2.3.1 on page 36) and surgery (see section 2.3.2 on page 38) if necessary, and tune the displays only after these steps.

Function *plot_nodes()*, and most plotting functions of *sprof*, allow various modes to control the use of colours. The recommended mode is to add a index variable *icol* giving an index to a colour table. See *rkindex()* for a choice of modifications. By default, the colour table is the current palette. The plot functions allow to chose a different colour function plot by plot.

ToDo: add general discussion of colours in *sprof*

Input

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

Here is the node view, using these choices (Basic information on node level, colour by total time, see fig. 5 on the next page).

Input

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

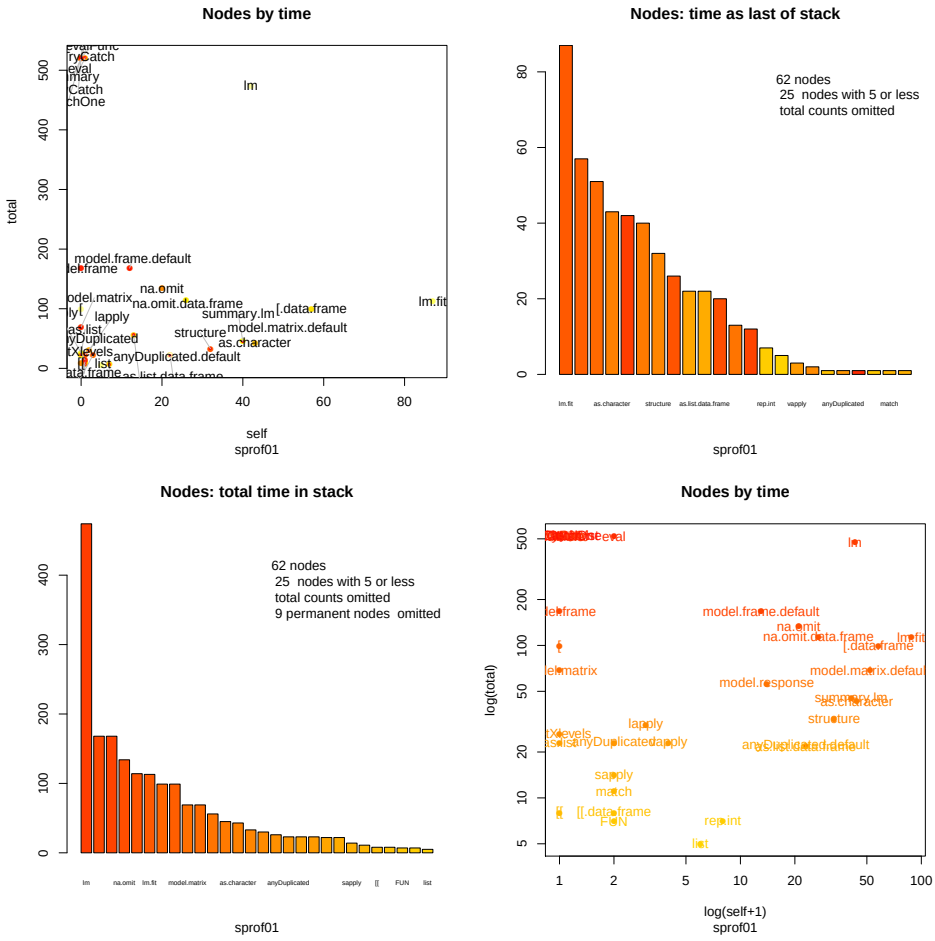


FIGURE 5. Basic information on node level, colour by total time.

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 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 `icol` is reserved for this) and use this as an index to a choice of colour palettes. So `icol` becomes a part of the data structure, and the colour palette to be used is passed as a parameter. Package `RColorBrewer` may be a helpful source for colour palettes. You find more information and references on this in the vignette "Bertin matrices" of package `bertin`, `http://bertin.r-forge.r-project.org`.

ToDo: improve colour: support colour in a structure

On the node level, the additional information provided by *sprof* is the information on runs. The total time is broken down to runs. Detailed information is accessible if requested. The nuber of runs, and the average run length are reported by default, giving an inside look into the total time reported conventionally. Making use of this information requires careful reading.

By itself, graphical inspection is not very helpful. The profiling process collapsed the real information to very few run numbers, and many ties in the average run lenght (fig. 6: Run length information.).

ToDo: add as additional information to classical basics

Input

```
plot(sprof01$nodes$avg_time,sprof01$nodes$nr_runs)
```

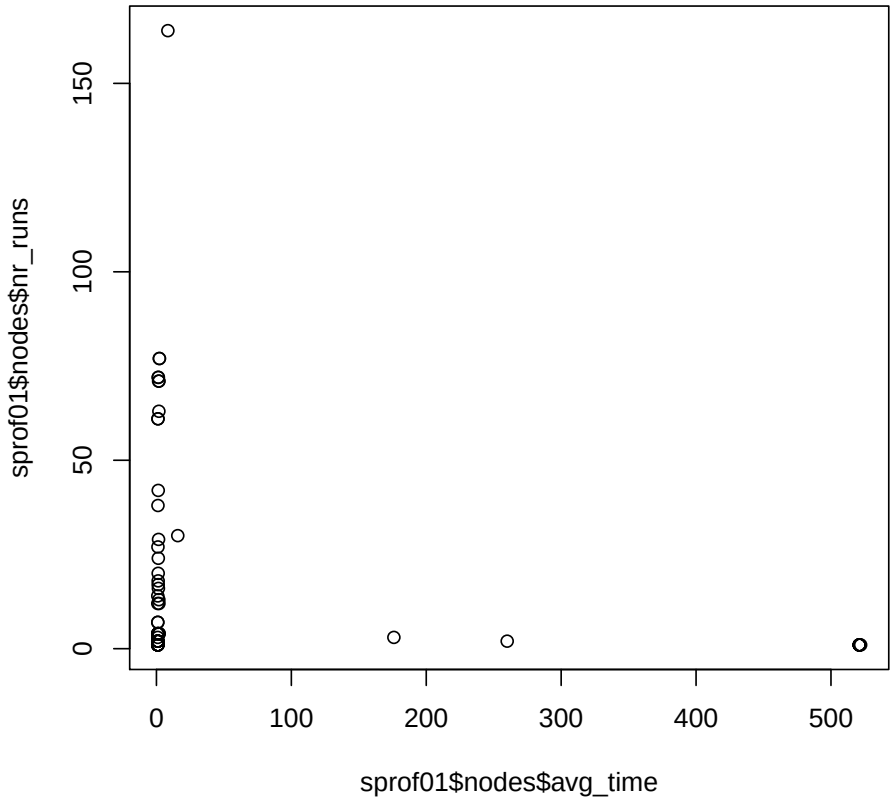


FIGURE 6. Run length information.

1.2.1. *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. In *sprof*, we are 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: colour by class – redo. Bundle colour index with colour?

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 very helpful tags. These can be used in a flexible way.

ToDo: add class by keyword

Input

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

Input

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

Output

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

Input

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

Nodes by package, colours by RColorBrewer: see fig. 7 on the next page.

Input

```
oldpar <- par(mfrow=c(3,2))
if (require(RColorBrewer)) colpack <-
        brewer.pal(length(levels(sprof01$nodes$icol)), "Paired") else
        colpack <- rainbow(length(levels(sprof01$nodes$icol)))
plot_nodes(sprof01, which=1:6, col=colpack)
par(oldpar)
```

If you want to, you can use your own classification to group variables.

Input

```
x_apply <- c("apply", "lapply", "vapply", "sapply")
x_as <- c("as.list", "as.data.frame", "as.list.data.frame",
        "as.character", "as.list.default","as.name")
```

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

Input

```
nodeclass <- rep("x_nn", sprof01$info$nrnodes)
nodeclass[sprof01$nodes$name %in% x_apply] <- "x_apply"
nodeclass[sprof01$nodes$name %in% x_as] <- "x_as"
```

or use assignments on the fly

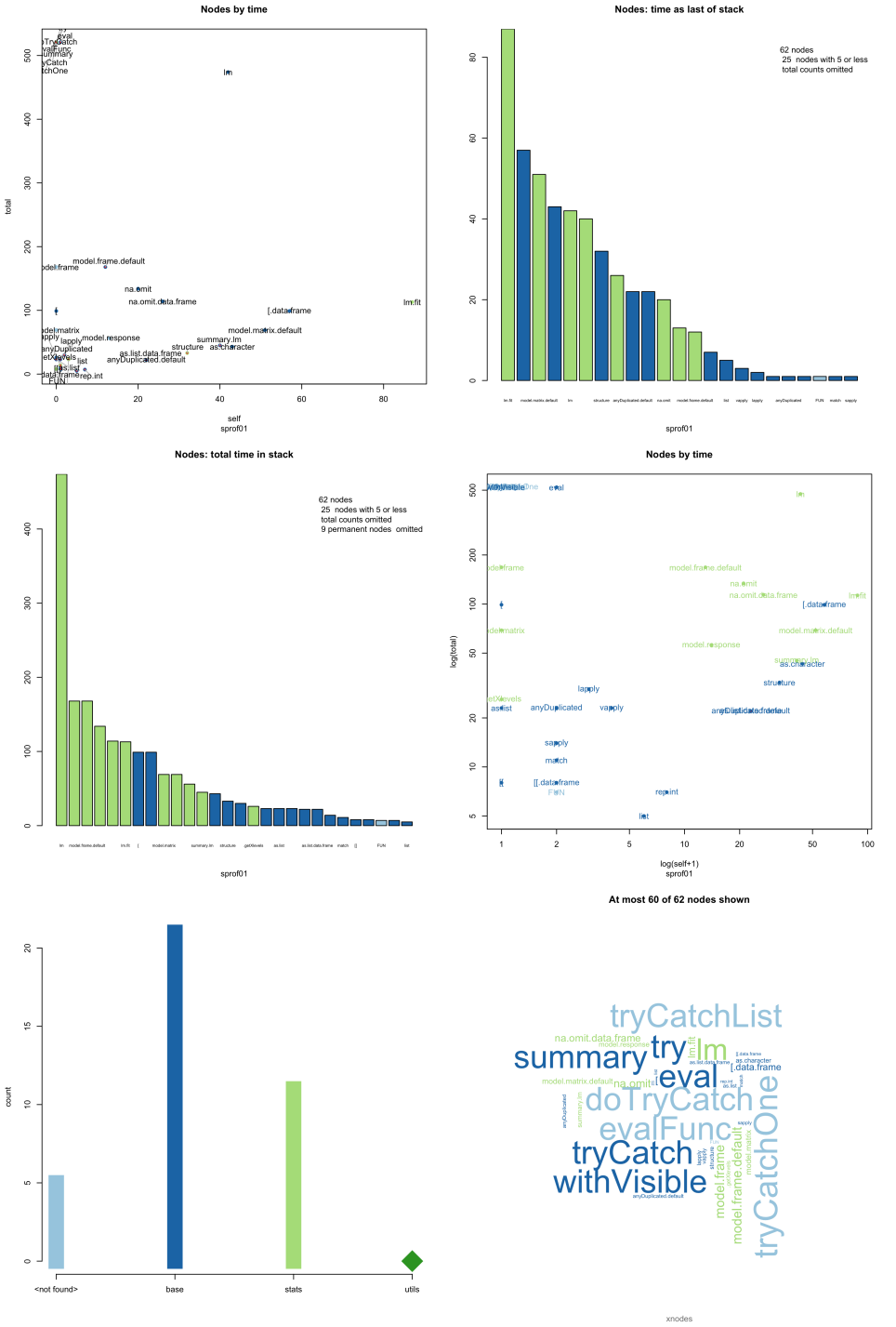


FIGURE 7. 12 18 Nodes by package, colours by RColorBrewer.

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"
nodeclass[sprof01$nodes$name %in%
        c("lm", "lm.fit", "summary.lm")
                ] <- "x_lm"
nodeclass[sprof01$nodes$name =="<Anonymous>"] <- "x_Anon"
nodeclass[sprof01$nodes$name == "Sweave"] <- "x_Sweave"
nodeclass[sprof01$nodes$name %in% c( "summary" ,"summary.lm")] <-
         "x_summary"
```

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

adds a sticky colour attribute. For interpretation, you should choose your preferred colour palette, for example

Input
```
nodeclasscol <- c("red", "green", "blue", "yellow",
        "cyan", "magenta", "purple",
        "brown", "aquamarine")

# gold cyan4 aquamarine pink violet orchid hotpink salmon turquoise1
```

ToDo: Defaults by class
ToDo: classes need separate colour palette, distinct from package or keyword.

Nodes by class, user defined colours: see fig. 8 on the facing page.

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

ToDo: check/fix colour for wordcloud

Nodes by class: default colour selection: see fig. 9 on page 26.

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

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.

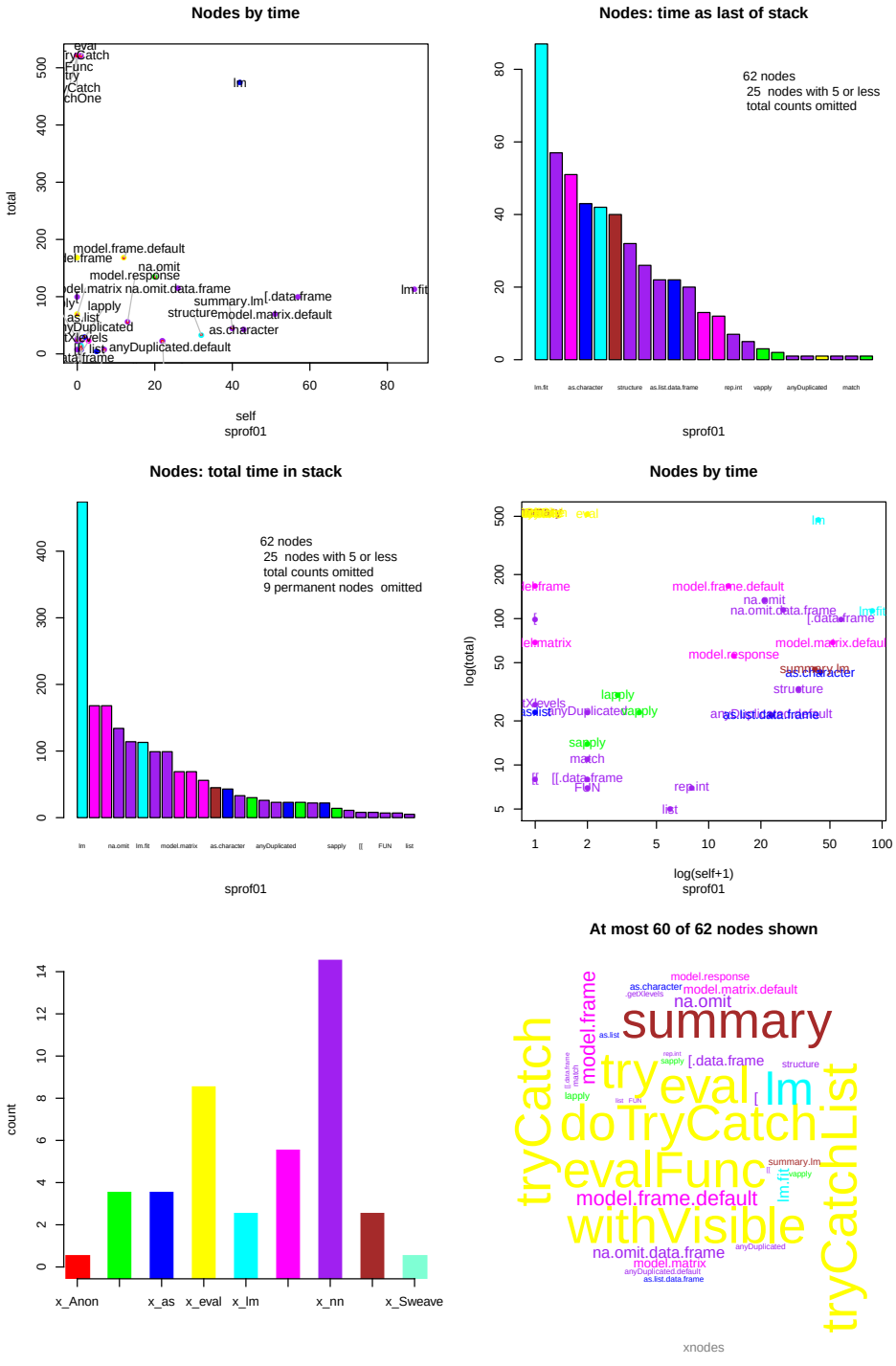


FIGURE 8. 8 12 Nodes by class. User defined colours.

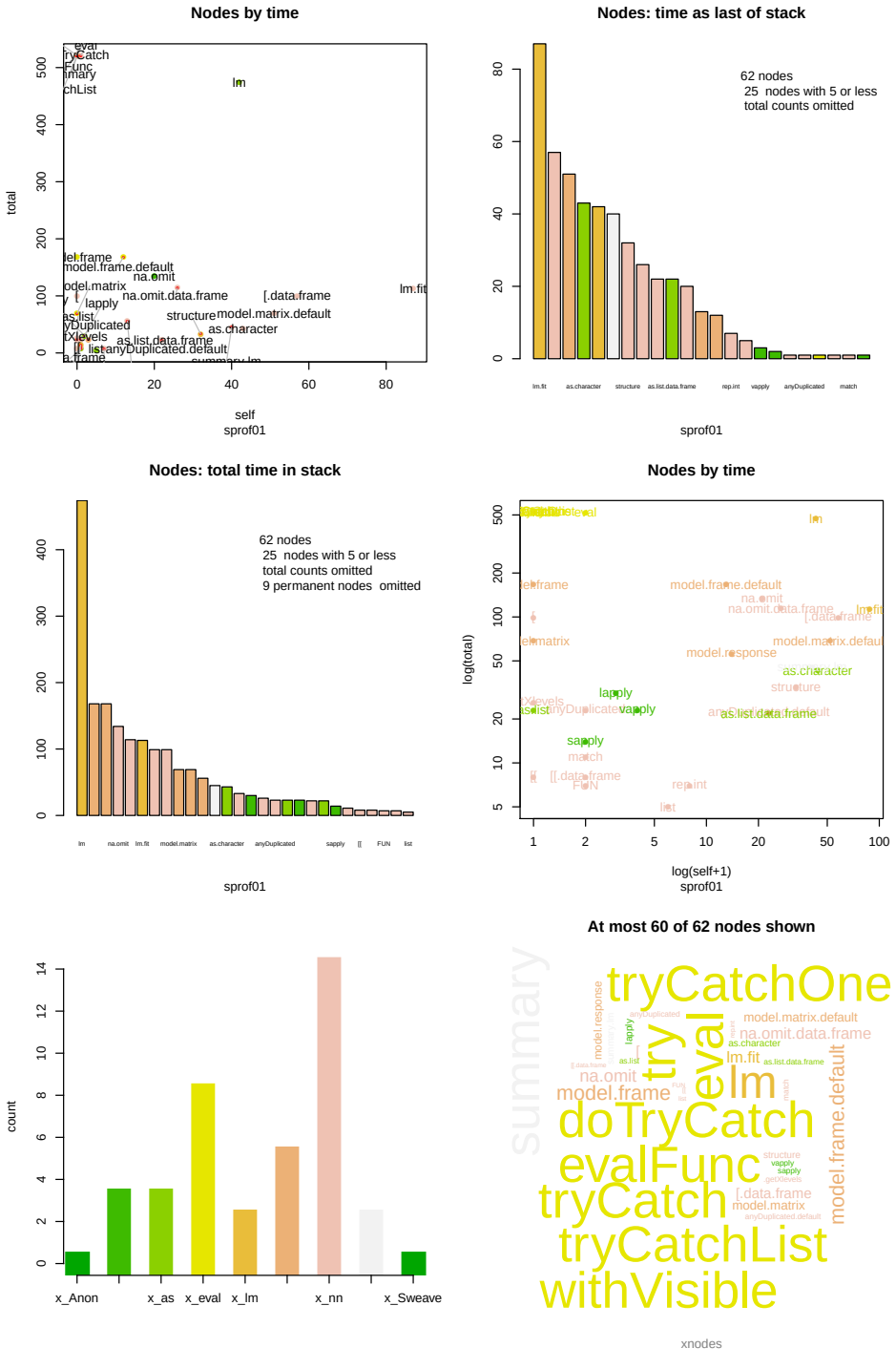


FIGURE 9. 8 12 Nodes by class: default colour selection

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 12.

ToDo: add attributes to stacks, and discuss scope

ToDo: sorting/arranging stacks

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: re-think: sort stacks

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.

Note: after sanitising, stacks may have an empty node list. However they should not be removed as long as they may still be present in the profile.

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. (This may change.)

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.

ToDo: data frame conversion?

ToDo: complete matrix conversion

TABLE 5. Extraction and conversion routines

profiles_matrix()	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

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.

ToDo: check and stabilise colour linking

See fig. 10 on the facing page for a summary of nodes by stack and profile.

Input

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

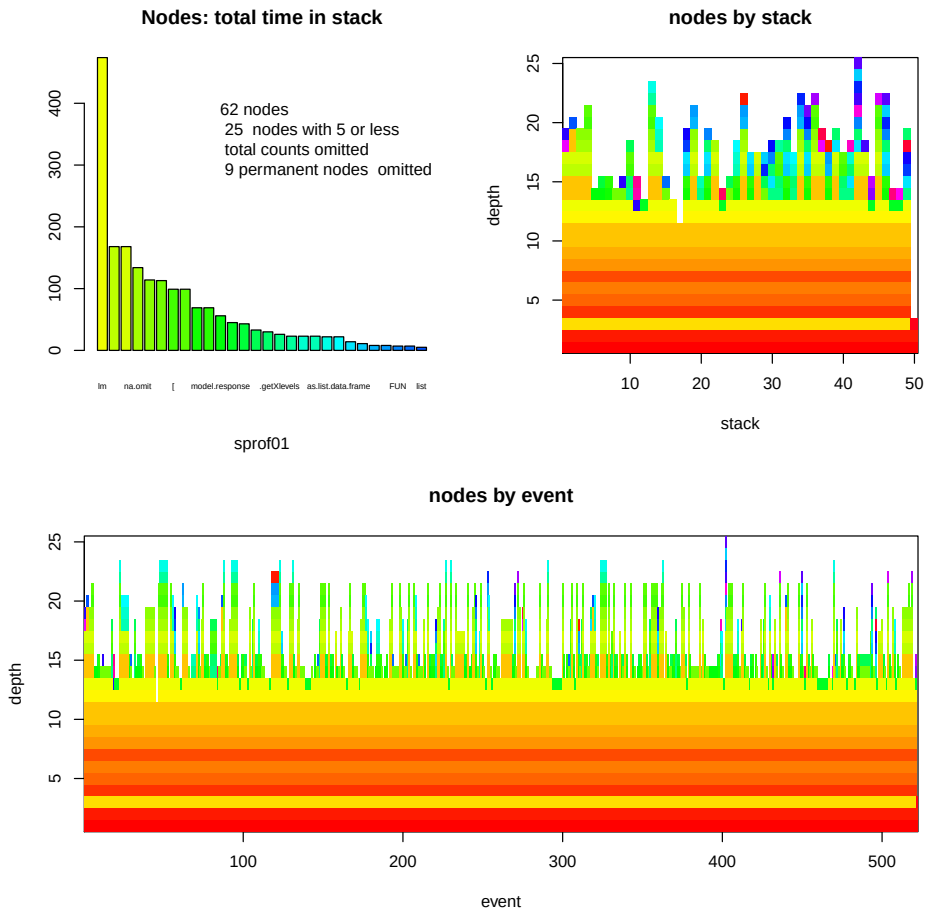


FIGURE 10. Nodes by stack and profile

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.

Input

```
stacks_nodes <- list.as.matrix(sprof01$stacks$nodes)
nrnodes <- apply(stacks_nodes,1,function(x) {length(unique(x))})
cat("nr unique nodes per stack level\n")
```

Output

```
nr unique nodes per stack level
```

Input
```
nrnodes
```

Output
```
 [1]  1  1  2  2  2  2  2  2  2  2  2  2  4 11 12 10 10 16  9  8  6  8
[23]  3  2  2
```

Nr. of unique nodes by stack level: See fig. 11.

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")
```

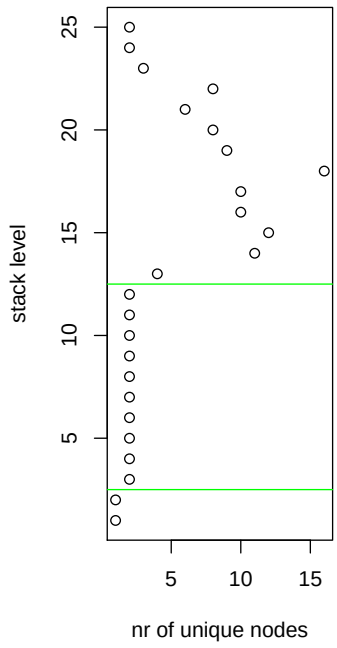


FIGURE 11. Nr of unique nodes by stack level.

ToDo: check and synchronise

We will come to finer tools in section 2.4 on page 41 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 three leaves us to start with five roots on the next level in our example.

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.

Input
```
summary(sprof01$stacks$refcount)
```

Output
```
   Min. 1st Qu.  Median    Mean 3rd Qu.    Max.
    1.0     1.0     2.0    10.4    12.0    87.0
```

ToDo: 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.

Input
```
summary(sprof01$stacks$refcount[sprof01$stacks$refcount>2])
```

Output
```
   Min. 1st Qu.  Median    Mean 3rd Qu.    Max.
   3.00    7.75   16.00   24.30   40.50   87.00
```

Input
```

```

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. In our example, *lm* or *summary.lm* may be convinient markers.

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 36.

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

ToDo: colours. recolour. Propagate colour to graph.

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 46 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[1]. *graph* on its side has an undocumented feature: it needs *Rgraphviz* to handle graph attributes[2]. We 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 *"graphNEL"* format which is shared by *graph* and *Rgraphviz*. *Rgraphviz* is hidden in the use of *plot*(). So here is a bare foot approach (Call graph derived from profile information: See fig. 12 on the next page.). A more sophisticated function implementation is in section 3 on page 46.

Input

```
#6 sprofadjNEL02
library(graph)
sprof01adjNEL <- as(adjacency(sprof01),"graphNEL")
plot(sprof01adjNEL,  main="sprof01: graph layout example",
```

[1]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.
[2]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.

```
        sub=sprof01$info$id,
        attrs=list(node=list(cex=4, fontsize=40,  shape="ellipse")),
        cex.main=2)
rm(sprof01adjNEL) # used for look-ahead example only
```

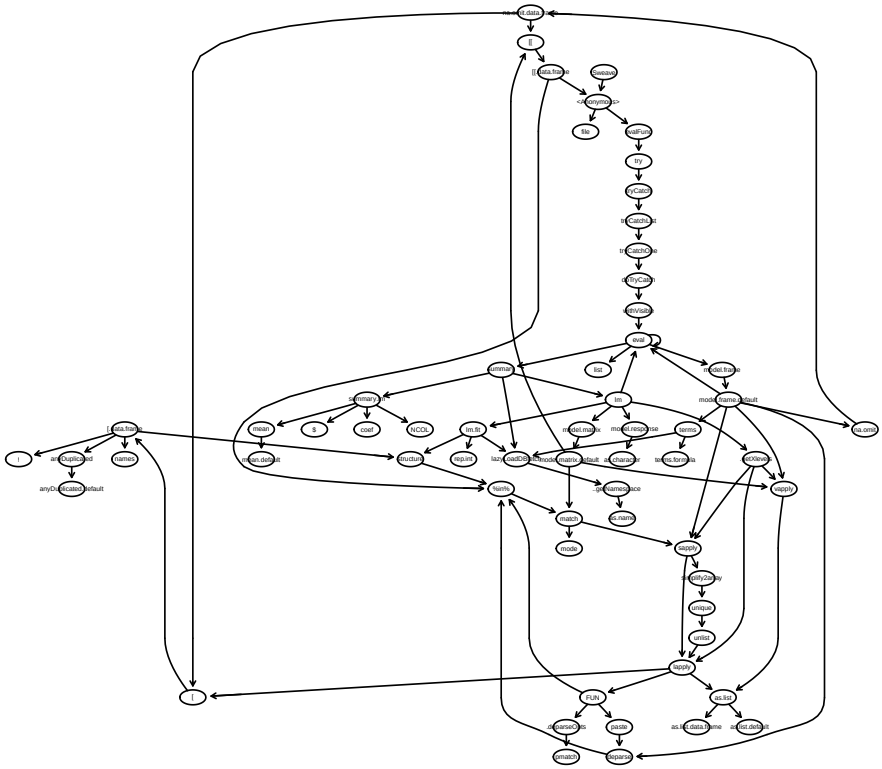


FIGURE 12. Call graph derived from profile information

ToDo: log scale?

2.3. **Cheap thrills.** Before starting additional inspection, the data better be trimmed. Trimming routines are in section 2.3.1 on page 36, but the data structure is robust enough to allow manual intervention as used here.

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

Input

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

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

Input

```
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, and instead of cutting off the stacks we replace them by an empty mark.

ToDo: should this be NA or NULL?

Input

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

ToDo: handle empty stacks and zero counts gracefully

ToDo: add a purge function

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 *info* section is inconsistent at this point. Another reason to call *updateRprof*().

Input

```
strx(sprof02$info)
```

Output

```
##strx: sprof02$info
'data.frame':        1 obs. of  9 variables:
$ id : chr "sprof02: trimmed"
$ date : POSIXct, format: "2013-09-10 18:00:07"
$ nrnodes : int 62
$ nrstacks : int 50
$ nrrecords : int 522
$ sample.interval: num 0.001
$ sampling.time : num 0.522
$ ctllines : chr "sample.interval=1000"
$ ctllinenr : num 1
```

Input

```
nodes<- sprof02$nodes[,-8]; rownames(nodes) <- NULL
prxt(nodes,
        caption="sprof02, before update",
        label="tab:sprof02info1",
        digits=c(0,0,0,2,0,2,0,2),
        zero.print=" . "
        )
```

Table 7: sprof02, before update

	name	self.time	self.pct	total.time	total.pct	nr_runs	avg_time
1	!	2	0.38	2	0.03	2	1.00
2	..getNamespace	.	0.00	1	0.01	1	1.00
3	.deparseOpts	2	0.38	4	0.05	4	1.00
4	.getXlevels	.	0.00	26	0.34	20	1.30
5	[	.	0.00	99	1.29	72	1.38
6	[.data.frame	57	10.92	99	1.29	72	1.38

7	[[	.	0.00	8	0.10	4	2.00
8	[[.data.frame	1	0.19	8	0.10	4	2.00
9	%in%	1	0.19	4	0.05	4	1.00
10	<Anonymous>	6	1.15	522	6.79	3	176.00
<cut>		⋮	⋮	⋮	⋮	⋮	⋮
53	terms	.	0.00	2	0.03	2	1.00
54	terms.formula	1	0.19	1	0.01	1	1.00
55	try	.	0.00	521	6.78	1	521.00
56	tryCatch	.	0.00	521	6.78	1	521.00
57	tryCatchList	.	0.00	521	6.78	1	521.00
58	tryCatchOne	.	0.00	521	6.78	1	521.00
59	unique	3	0.57	4	0.05	4	1.00
60	unlist	.	0.00	1	0.01	1	1.00
61	vapply	3	0.57	23	0.30	16	1.44
62	withVisible	.	0.00	521	6.78	1	521.00

Input

Input

```
sprof02 <- updateRprof(sprof02)
sprof02$info$id <- "sprof02 updated"
```

Input

```
strx(sprof02$info)
```

Output

```
##strx: sprof02$info
'data.frame':        1 obs. of  10 variables:
$ id : chr "sprof02 updated"
$ date : POSIXct, format: "2013-09-10 18:00:07"
$ nrnodes : int 62
$ nrstacks : int 50
$ nrrecords : int 522
$ sample.interval: num 0.001
$ sampling.time : num 0.522
$ ctllines : chr "sample.interval=1000"
$ ctllinenr : num 1
$ date_updated : POSIXct, format: "2013-09-10 18:00:14"
```

Input

```
nodes<- sprof02$nodes[,-8]; rownames(nodes) <- NULL
prxt(nodes,
        caption="sprof02, after update",
        label="tab:sprof02info2",
        #digits=c(0,0,0,2,0,2,0,2,0,2),
        digits=c(0,0,0,2,0,2,0,2),
        zero.print=" . "
)
```

Table 8: sprof02, after update

	name	self.time	self.pct	total.time	total.pct	nr_runs	avg_time
1	!	1	0.23	1	0.06	1	1.00
2	..getNamespace	.	0.00	1	0.06	1	1.00
3	.deparseOpts	2	0.46	4	0.25	4	1.00
4	.getXlevels	.	0.00	26	1.64	20	1.30
5	[	.	0.00	98	6.17	71	1.38
6	[.data.frame	57	13.16	98	6.17	71	1.38
7	[[	.	0.00	8	0.50	4	2.00
8	[[.data.frame	1	0.23	8	0.50	4	2.00
9	%in%	1	0.23	4	0.25	4	1.00
10	<Anonymous>	6	1.39	6	0.38	2	3.00
<cut>		⋮	⋮	⋮	⋮	⋮	⋮
53	terms	.	0.00	1	0.06	1	1.00
54	terms.formula	1	0.23	1	0.06	1	1.00
55	try	.	0.00	.	0.00	.	0.00
56	tryCatch	.	0.00	.	0.00	.	0.00
57	tryCatchList	.	0.00	.	0.00	.	0.00
58	tryCatchOne	.	0.00	.	0.00	.	0.00
59	unique	3	0.69	4	0.25	4	1.00
60	unlist	.	0.00	1	0.06	1	1.00
61	vapply	3	0.69	23	1.45	16	1.44
62	withVisible	.	0.00	.	0.00	.	0.00

Input

Nodes by stack and profile: see fig. 13 on the facing page.

Input

```
#8 8
shownodes(sprof02)
```

2.3.1. *Trimming.* Note: trimming may be supported by the graph packages. If you are more familiar with your graph package, you may prefer to handle trimming there.

To decide about trimming, we can use the displays shown above, or we can use some statistics. Global trimming works on the stack level. To make life easy, we can imbed the stack information in a matrix and use marginals.

Input

```
stackm01 <- list.as.matrix(sprof01$stacks$nodes)
nr_uniquenodes01 <- apply(stackm01,1, function(x) { length(unique(x))})
names( nr_uniquenodes01)<- rownames(nr_uniquenodes01,FALSE,"L")
cat("Nr of unique nodes by level\n")
```

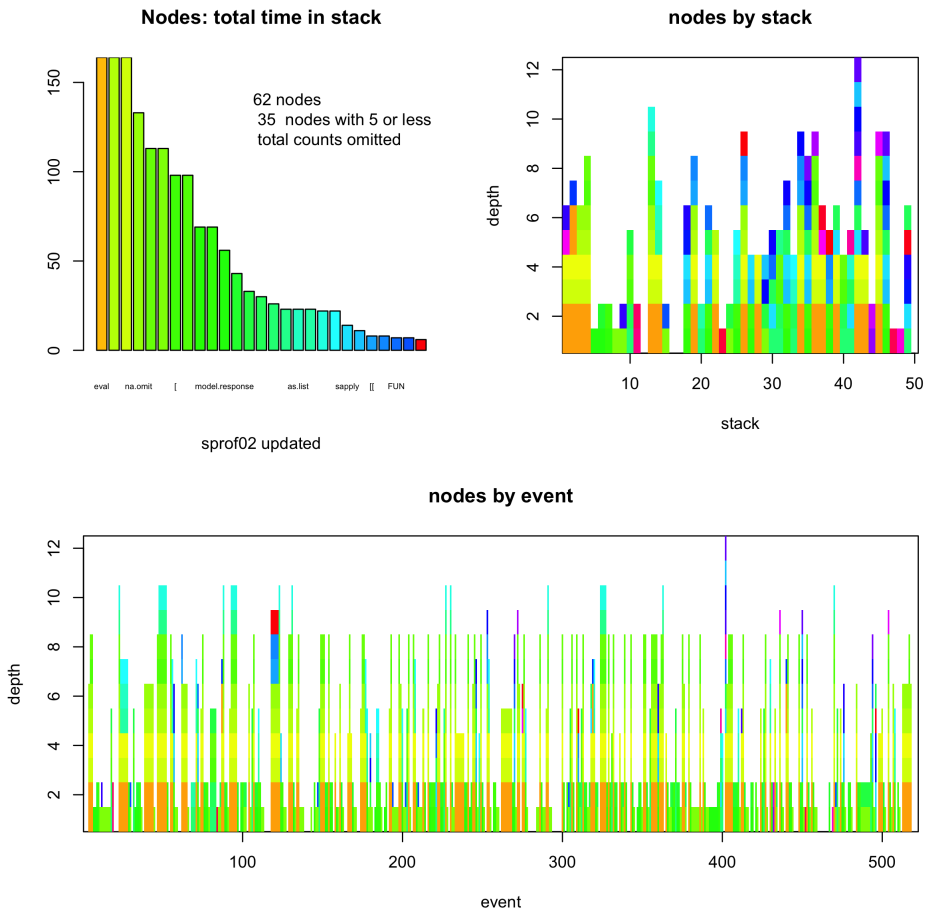


FIGURE 13. sprof02: Nodes by stack and profile

Output

```
Nr of unique nodes by level
```

Input

```
nr_uniquenodes01
```

Output

```
 L1  L2  L3  L4  L5  L6  L7  L8  L9 L10 L11 L12 L13 L14 L15 L16 L17 L18
  1   1   2   2   2   2   2   2   2   2   2   2   4  11  12  10  10  16
L19 L20 L21 L22 L23 L24 L25
  9   8   6   8   3   2   2
```

Input

```
rm(stackm01, nr_uniquenodes01)
```

Input

ToDo: This section needs to be reworked

```
rsprof01Tr <- trimstacks(sprof01, level=11)
```

After trimming, it is worth to inspect the result and to see whether additional trimming is helpful. Function *roots_sprof* is handy.

Input
```
roots_sprof(sprof01,stacks=rsprof01Tr)
```

Output
```
summary    <NA>
     50      NA
```

Here after trimming we have one root node. This may be esthetically pleasing, but it is more informative to note that *summary* is a common root, and then cut it off (thus fragmenting the graph into components).

Input
```
rsprof02Tr <- trimstacks(sprof01, level=12)
roots_sprof(sprof01,stacks=rsprof02Tr)
```

Output
```
            lm lazyLoadDBfetch      summary.lm            <NA>
            29              27              51              NA
```

Now this is revealing. We know how our test example was contructed. So we are prepared to find *lm* and *summary.lm* at prominent positions. But it is surprising to find *lazyLoadDBfetch* as dominant node, and with a frequency comparable to that of *lm*.

ToDo: move to other section

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. Surgery can also been done conveniently on the source level, using simple replace statements.*

The surgery functions will be moved to the package in the next release.

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 469 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.

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

Other candidates are:
`"as.list" "vapply" "model.frame.default" -> "<model_as.list>"`
or
`"as.list" "vapply" "model.matrix.default" -> "<model_matrix_as.list>"`

ToDo: Implement. Currently best handled on source=text level

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".

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

Input

```
addnode <- function(nodes, newnode, warn = options("warn"))
{
        i <- match(newnode, nodes$name, nomatch=0)
        if (i==0){
                nodes$name <- as.character(nodes$name)
                nodes <- rbind(nodes,NA)
                i <- length(nodes$name)
                nodes$name[i] <- newnode
                rownames(nodes) <- nodes$name
        if (as.logical(warn))
                message("addnode: node added. An updateRprof() may be necessary.")
        }
        return(nodes)
}
```

Input

```
sprof03 <- sprof02; sprof03$info$id <- "sprof03: surgery"
nodes <- addnode(sprof03$nodes, "<.getXlevels_[>", warn=FALSE)
nodes <- addnode(nodes, "<model_as.list>", warn=FALSE)
sprof03$nodes <- addnode(nodes, "<model_matrix_as.list>", warn=FALSE)
```

ToDo: xreplacenodes: improve implement

So far, we use factor indices only.

ToDo: clean up factor handling

Input

```
xwhere <- function(oldseq, x){
        x <- unlist(x)
        firstpos <- 1; lastpos <-  length(x) -  length(oldseq) +1

        if (lastpos <1) return(0)
        l <- length(oldseq)-1
        while (firstpos <= lastpos) {
                if ( isTRUE (
                all.equal(oldseq , x[firstpos:(firstpos+l)], check.attributes=FALSE)
                ) )
                        {return(firstpos) }
                firstpos <- firstpos+1
                }
```

```
        return(0)
}
```

Input

```
xreplace <- function(oldseq, newseq, x){
#! handle multiple replacements
        wh <- xwhere(oldseq,x)
        if (wh) {
        l <- length(oldseq)-1
        if (wh>1) x1 <- c(x[1:(wh-1)], newseq) else x1 <- newseq
        if (wh+l < length(x)) x <- c(x1, x[ -(1:(wh + l))] ) else x <- x1
        return(x)
        } else return(x)
}
```

Input

```
nodenames2index <- function(names, sprof){
if (is.character(names))
                { sapply(names, function(x) {match(x,sprof$nodes$name)})
                  #if (is.na(trimnode)) return(ts)
                } else names
}
```

Input

```
xreplacenodes <- function(sprof, oldseq, newseq)
{
        if (is.character(oldseq)) oldseq <- nodenames2index(oldseq,sprof)
        if (is.character(newseq)) newseq <- nodenames2index(newseq,sprof)
        stacks <- sprof$stacks$nodes
        sapply(stacks, function(x){xreplace(oldseq, newseq, unlist(x))})
}
```

ToDo: stackssrc, collstacksdictrev etc. now out of date

Input

```
sprof03$stacks$nodes <- xreplacenodes(sprof03,
        c(".getXlevels", "lapply", "["),
        "<.getXlevels_[>")
sprof03$stacks$nodes <- xreplacenodes(sprof03,
        c("model.frame.default", "vapply", "as.list"),
         "<model_as.list>")
sprof03$stacks$nodes <- xreplacenodes(sprof03,
        c("model.matrix.default", "vapply", "as.list"),
         "<model_matrix_as.list>")
sprof03 <- updateRprof(sprof03)
#asfactormodel(sprof03$stacks$nodes, sprof03$nodes$name)
```

ToDo: warn about undefined vars, e.g. rle, class, ...

ToDo: replace by sprof03

ToDo: fix null name

This surgery gives no additional information on the profile data. It only helps the graph representation, by extending the one-step information given by the edges to two or omre step information at some critical points.

We use a prepared sanitised version of our data set.

Input

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

Applying our old classification:

Input

```
nodeclass <- rep("x_nn", length(sprof04$nodes$name))
nodeclass[sprof04$nodes$name %in% x_apply] <- "x_apply"
nodeclass[sprof04$nodes$name %in% x_as] <- "x_as"
nodeclass[sprof04$nodes$name %in%
        c("eval",  "evalFunc",
                "try", "tryCatch", "tryCatchList", "tryCatchOne",
                "doTryCatch", "withVisible")
                ] <- "x_eval"
nodeclass[sprof04$nodes$name %in%
        c("model.frame", "model.matrix.default","model.frame.default",
         " model.response", "model.matrix", "model.response")
                ] <- "x_model"
nodeclass[sprof04$nodes$name %in%
        c("lm", "lm.fit", "summary.lm")
                ] <- "x_lm"
nodeclass[sprof04$nodes$name =="<Anonymous>"] <- "x_Anon"
nodeclass[sprof04$nodes$name == "Sweave"] <- "x_Sweave"
nodeclass[sprof04$nodes$name %in% c( "summary" ,"summary.lm")] <-
         "x_summary"
sprof04$nodes$icol <-as.factor(nodeclass)
```

now gives

Input

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

Sanitised nodes by class, user defined colours: see fig. 14 on the following page.

Surgery is a means to clarify the graph structure. It should be applied with some sense. Some collusions are so intuitive that they can be ignored without surgery.

Keep in mind Simpson's paradox. If you upgrade the weights after surgery, a single node may be split to different containers, thus seemingly reducing the weights.

ToDo: add smart surgery with memory for attributing resources.

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 to the profile information. This gives a detailed information about the presence time of each node, by stack level.

ToDo: use sprof02 or sprof03?

Input

```
profile_nodes <- profiles_matrix(sprof02)
profile_nodes_rle<- rrle(profile_nodes, collapseNA=FALSE)
```

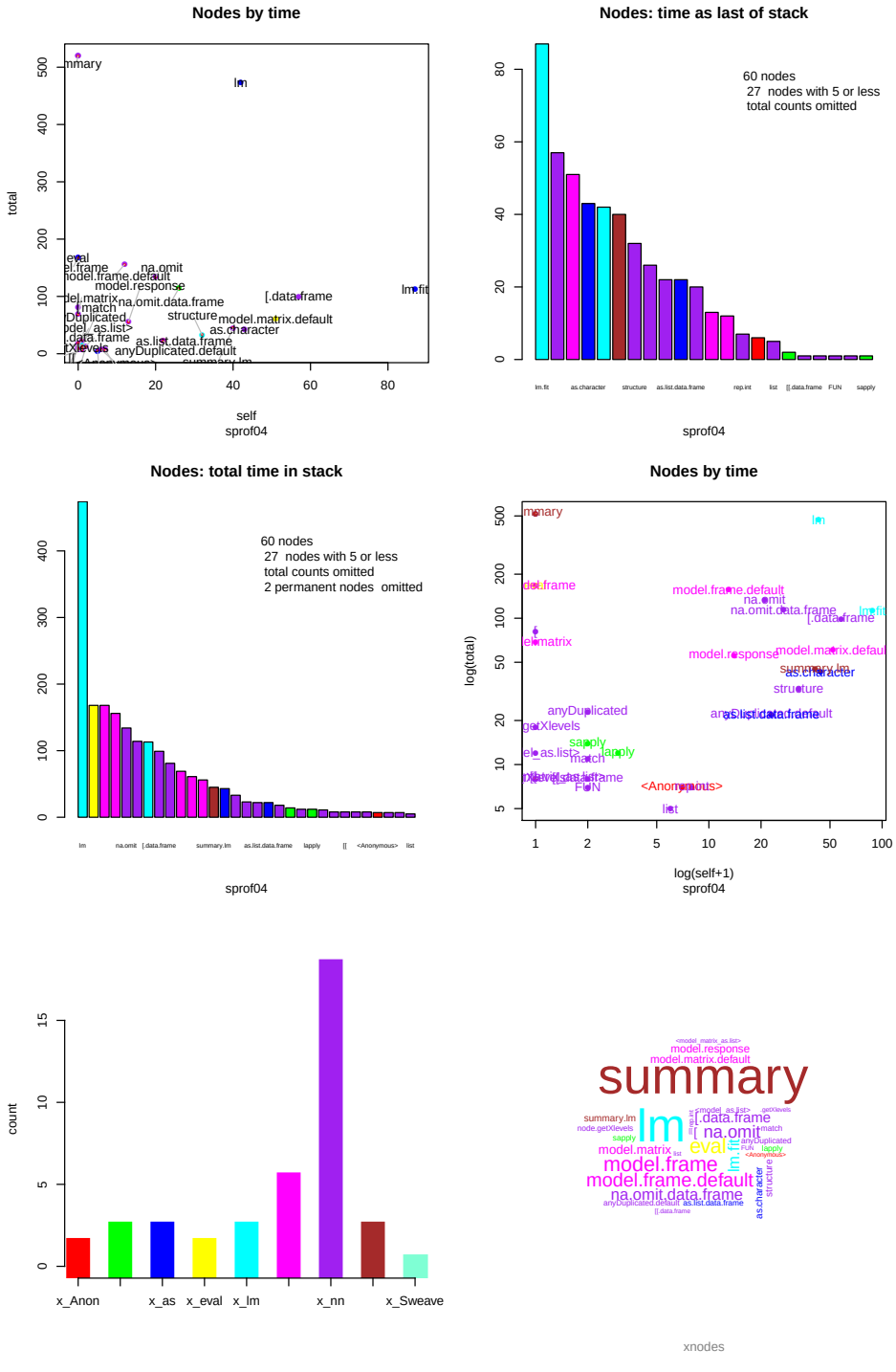


FIGURE 14. Sanitised. Nodes by class. User defined colours.

```
#!NA needs special case in run length handling.

strx(profile_nodes_rle, list.len=5)
```

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
..$ 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.

Input

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

Output

```
##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]*.

Input

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

ToDo: Warning: data structure still under discussion

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

ToDo: add summary for NA

ToDo: add marginals and conditionals. Provide function node_summary.

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

Output

```
     run_length
level  1  2  3  4  5  6  7 Sum
  1    .  .  .  .  .  .  .   .
  2    .  .  .  .  .  .  .   .
  3   40 17  7  4  2  6  1  77
  4    .  .  .  .  .  .  .   .
  5    .  .  .  .  .  .  .   .
  6    .  .  .  .  .  .  .   .
  7    .  .  .  .  .  .  .   .
  8    .  .  .  .  .  .  .   .
  9    .  .  .  .  .  .  .   .
  10   .  .  .  .  .  .  .   .
  11   .  .  .  .  .  .  .   .
  12   .  .  .  .  .  .  .   .
  Sum 40 17  7  4  2  6  1  77
```

Input

```
amt <- nodesrunlength(sprof02)
prxt(amt,
         caption=paste0("Marginal statistics on nodes by run length, ",
                 "sorted by total time used"),
        label="tab:pramt4",
        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

	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
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
.getXlevels	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>	1	.	.	.	1	.	.	2	6	3.00
.deparseOpts	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
mode	2	.	.	.	.	.	.	2	2	1.00
names	2	.	.	.	.	.	.	2	2	1.00
pmatch	2	.	.	.	.	.	.	2	2	1.00
!	1	.	.	.	.	.	.	1	1	1.00
..getNamespace	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

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

ToDo: hack. keep length in nodesrun-length

From the summary information, `nr_runs` and `avg_time` are included in the node information by default. We can use this information to enhance graphical displays. See: nodes marked by run length and run count, fig. 19 on page 55.

`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 lenght.

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.

ToDo: table: node #runs min median run length max

3. Graph package

What we have achieved so far can be seen from the graph representations. For our purposes, an edge table is most convenient. To allow for edge attibutes, we can use an R `data.frame` as provided by

We can make use of any graph mapping package. Unfortunately, each seem to have its own calling convention. So we habe to do some translation.

Input

```
#12 6
library(graph)
oldpar <- par(mfrow=c(1,2))
        plotviz(sprof01)
        plotviz(sprof02)
par(oldpar)
```

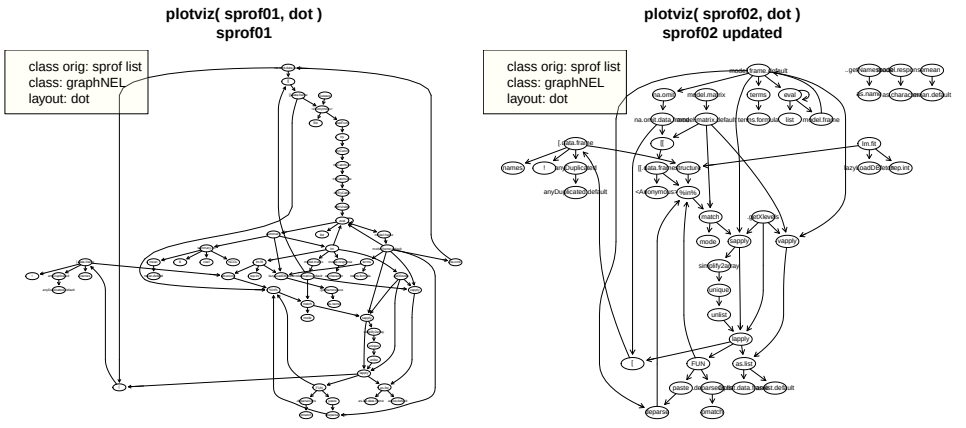


Figure 15. Sprof graph, before and after trimming.

See fig. 15 on the preceding 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 to.

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. (Sprof graph, after trimming, and after trimming and some sanitising, see fig. 16.)

ToDo: fix null name

Input

```
#12 6
sprof04 <- readRprof("RprofsRegressionExpl03.out", id="sprof04")
oldpar <- par(mfrow=c(1,2))
plotviz(sprof02)
plotviz(sprof04)
par(oldpar)
```

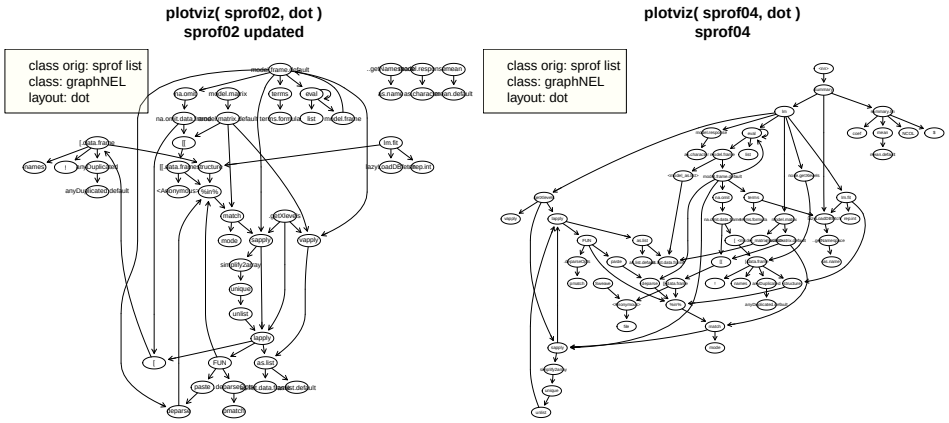


FIGURE 16. Sprof graph, after trimming, and after trimming and some sanitising.

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.

The functions in this chapter are not included in the package to avoid dependency of *sprof* on *graph* and other graph packages.

Input

```
library(graph)
library(Rgraphviz)
```

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

Input

```
as_graphNEL_sprof <- function(sprof, weight=TRUE){

        a04<-adjacency(sprof)

        rnames <- rownames(a04)

        if (!weight) {
        dimold <- dim(a04);a04 <- as.numeric(a04); dim(a04) <- dimold
        rownames(a04)<- rnames; colnames(a04)<- rnames;
        } #! define lwd first

        el04 <- edgedf(a04)
        el04$lwd  <- rkindex(el04$count, maxindex=7, ties.method="min")

        a04NEL <- as(a04,"graphNEL")
        nodeDataDefaults(a04NEL, "shape") <- "ellipse"
        nodeDataDefaults(a04NEL, "cex") <- 0.6
        nodeDataDefaults(a04NEL, "weigth") <- 1
        nodeDataDefaults(a04NEL, "fill") <- "green"
        nodeDataDefaults(a04NEL, "col") <- "yellow"

        a04NEL <- layoutGraph(a04NEL)

        nodeRenderInfo(a04NEL) <- list(shape="ellipse")
        nodeRenderInfo(a04NEL) <- list(cex=0.6, shape="ellipse")
        nodeRenderInfo(a04NEL) <- list(weight=1)
        #nodeRenderInfo(a04NEL) <- list(color="yellow")
        nodeRenderInfo(a04NEL) <- list(fill="yellow", col="blue")

        edgeDataDefaults(a04NEL,"lwd") <- 1
        edgeDataDefaults(a04NEL,"col") <- "grey"

        #nodeRenderInfo(a04NEL) <- list(weight=1)

        #edgeRenderInfo(a04NEL) <- list(lwd=el04$lwd)
        #edgeRenderInfo(a04NEL)$lwd <- el04$lwd
        for (i in 1:length(el04$lwd))
        {edgeRenderInfo(a04NEL)$lwd[i] <- el04$lwd[i]}
        a04NEL
}
```

As *as_graphNEL_sprof*, the following function is not included in the package to avoid dependency of *sprof* on *graph* and other graph packages.

Input

```
# source('~/projects/rforge/sintro/pkg/sprof/R/pgn.R', chdir = TRUE)
plot_graphNEL_sprof <- function(sprof,
        layoutType = "dot",
        nodeattrs, # = c("default", "time", "runs"),
        fill_list = NULL,
        lwd_list = NULL,
        srclegend="topleft",
        attrlegend="bottomright",
        main = NULL, sub = NULL,...)
{
        main1 <- deparse(substitute(sprof))
        xsubid <- NULL
        y<-NULL
        class1 <- NULL

        if (missing(nodeattrs)) nodeattrs <- "default"
        #nodeattrs <- match.arg(nodeattrs)
        # cat(nodeattrs)

        if (inherits(sprof, "sprof")) {
                class1 <- paste0("class orig: ", paste(class(sprof), collapse=" "))
                main1 <- sprof$info$id
                graphNEL <-
                        as_graphNEL_sprof(sprof, weight=FALSE)
                y <- layoutType
        } else graphNEL <- sprof

         if (!is.null(sub)) sub <- as.character(sub)

        main=paste0(main, "\n plot_graphNEL_sprof( ",
                main1 , ", ",
                layoutType, " )\n", xsubid)

# functions
        legattributes <- function(legnd=NULL,
                nodecol=NULL,  nodelwd=NULL,
                edgcol=NULL, edglwd="frequency")
        {
                #
                llegnd<- legnd
                if (!is.null(nodecol)) {
                        nodecol <- paste0("node col: ",nodecol, collapse="")
                        if (is.null(llegnd))  llegnd <- nodecol else
                                llegnd <- paste0(llegnd,"\n", nodecol,collapse="")
                }
                if (!is.null(nodelwd)) {
                        nodelwd <- paste0("node lwd: ", nodelwd, collapse="")
                        if (is.null(llegnd)) llegnd <- nodelwd else
                                llegnd <- paste0(llegnd,"\n", nodelwd, collapse="")
```

```
        }
        if (!is.null(edgcol)) {
                edgcol <- paste0("edge col: ", edgcol, collapse="\n")
                if (is.null(llegnd)) llegnd <- edgcol else
                        llegnd <- paste0(llegnd,"\n", edgcol,collapse="")
        }
        if (!is.null(edglwd)) {
                edglwd <- paste0("edge lwd: ", edglwd, collapse="")
                if (is.null(llegnd)) llegnd <- edglwd else
                        llegnd <- paste0(llegnd, "\n",edglwd,collapse="")
        }
        if (!is.null(attrlegend)){
        if (!is.null(llegnd))
        {        llegnd=paste0(llegnd,"\n")

        legend(attrlegend,
                        legend= llegnd,
                        bg="#FFFFE0",
                        seg.len=0,
                        bty="n",
                        text.font=3)
                #bg="#0000"
                }
                }
} # legattributes

graphNEL <- layoutGraph(graphNEL, layoutType=layoutType)

nodeDataDefaults(graphNEL, "shape") <- "ellipse"

if (nodeattrs == "runs"){
        amt <- nodesrunlength(sprof, clean=FALSE)

        #sprof$nodes$self.time -> fill
        fill_list <- heat.colors(12)[
                rkindex(-amt[,"avg_time"],
                        pwr=0.5, maxindex=12, ties.method="min")]
        names(fill_list) <- sprof$nodes$name

        #sprof$nodes$total.time -> lwd
        lwd_list  <-  rkindex(amt[,"nr_runs"],
                pwr=0.5, maxindex=7, ties.method="min")
        names(lwd_list) <- sprof$nodes$name

        #strx(nodeRenderInfo(graphNEL))
        nDD3 <- (nodeRenderInfo(graphNEL))

        nodeDataDefaults(graphNEL, "shape") <- "ellipse"

        nodeRenderInfo(graphNEL) <- list(lwd=lwd_list,
                fill=fill_list,
                col="#0000FF80",
                shape="ellipse",
```

```
                weight=1)
        #strx(nodeRenderInfo(graphNEL))
} else if (nodeattrs == "time"){
        # node attributes
        #sprof$nodes$self.time -> fill
        fill_list <- heat.colors(12)[
                rkindex(-sprof$nodes$self.time,
                        pwr=0.5, maxindex=12, ties.method="min")]
        names(fill_list) <- sprof$nodes$name

        #sprof$nodes$total.time -> lwd
        lwd_list  <-  rkindex(sprof$nodes$total.time,
                pwr=0.5, maxindex=7, ties.method="min")
        names(lwd_list) <- sprof$nodes$name

        #strx(nodeRenderInfo(graphNEL))
        nDD3 <- (nodeRenderInfo(graphNEL))
}   else {
        # graphNEL <- layoutGraph(graphNEL, layoutType=layoutType)
        fill_list <- NULL
        lwd_list <-  NULL
}
nodeDataDefaults(graphNEL, "shape") <- "ellipse"

nodeRenderInfo(graphNEL) <- list(
        lwd=lwd_list,
        fill=fill_list,
        col="#0000FF80",
        shape="ellipse",
        weight=1)

# edge attributes
# strx(edgeRenderInfo(graphNEL))
nER01 <- edgeRenderInfo(graphNEL)
edgeRenderInfo(graphNEL) <- list(col= "#80808080")
# strx(edgeRenderInfo(graphNEL))

renderGraph(graphNEL)

if (!is.null(srclegend))
        legend(srclegend,
                legend=c( class1,
                        paste0("class: ",
                                paste(class(graphNEL), collapse=" ")),
                        paste0("layout: ", layoutType)),
                bg="#FFFFE040",
                seg.len=0
                )#"lightyellow" =#FFFFE0

 if (nodeattrs == "runs"){
         title(main=paste0("nodes (run length)\n", main)        )
         legattributes(nodecol="avg run len, pwr=0.5",
                        nodelwd="nr runs, pwr=0.5")
```

```
        } else if (nodeattrs == "time"){
                title(main=paste0("nodes (time)\n", main))
                 legattributes( nodecol="self.time, pwr=0.5",
                        nodelwd="total.time, pwr=0.5")
        } else {
                title(main=main)
                legattributes( )
        }
        title( sub=sprof$info$id, col.sub=grey(0.5))
        invisible(graphNEL)
}# %:  plot_graphNEL_sprof(sprof02)
```

Input

```
plot_graphNEL_sprof(sprof04)
```

ToDo: remove global colour; implement local colour

Rgraphviz/graph basic plot: see fig. 17.

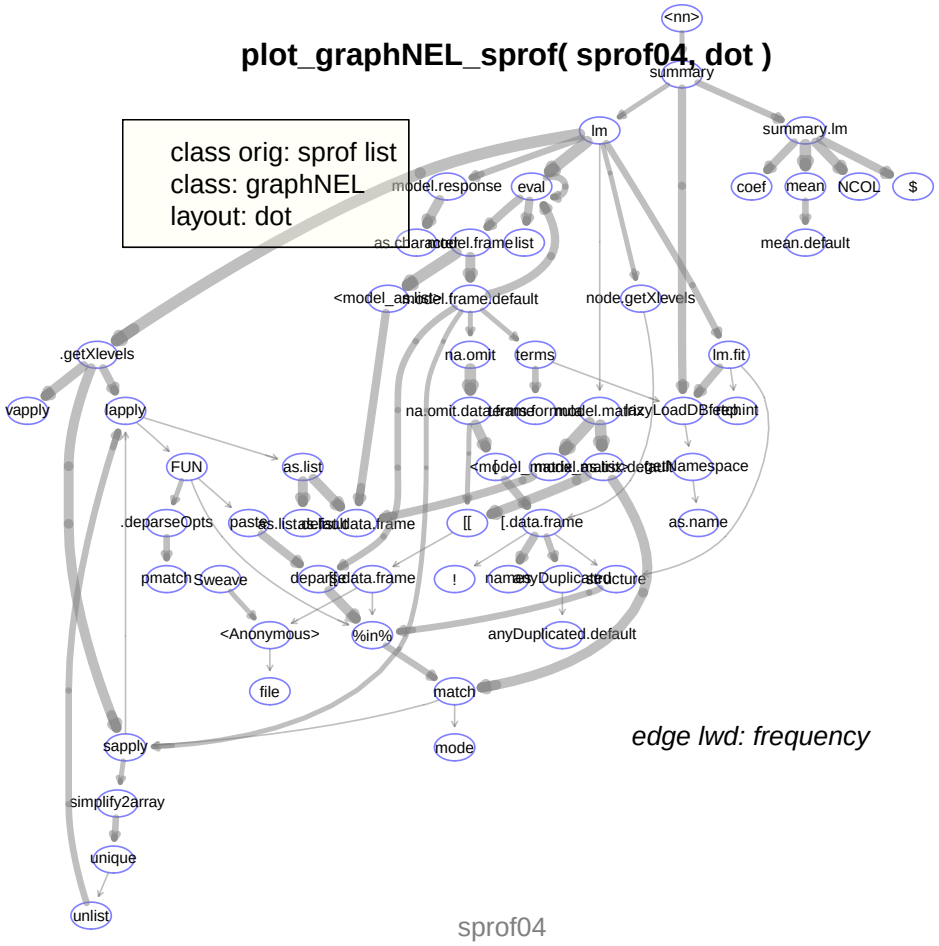


FIGURE 17. Rgraphviz/graph basic plot

To use attributes on nodes and edges, we need *Rgraphviz*.

ToDo: merge with as_graphNEL_sprof

Rgraphviz/graph plot with attributes: see fig. 18. This plot gives us the traditional view, highlighting nodes and edges with overall time presence.

Input

```
#6 6
plot_graphNEL_sprof(sprof04, nodeattrs="time")
```

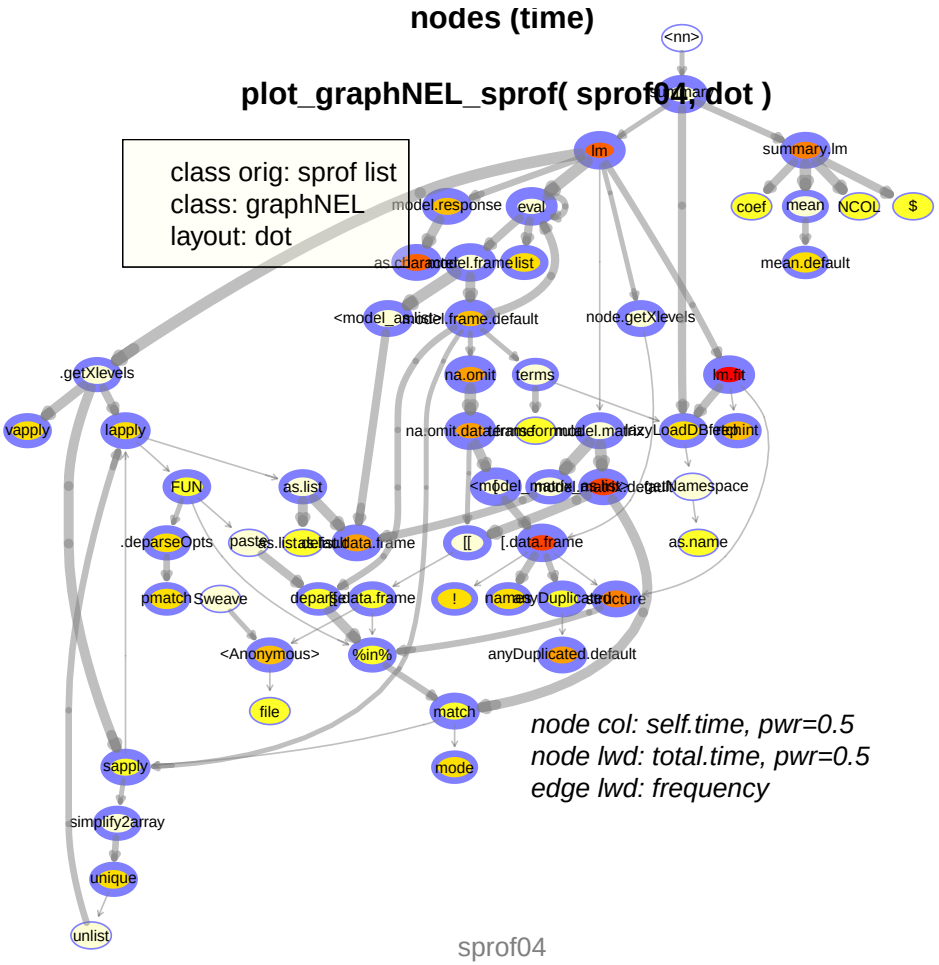


FIGURE 18. Rgraphviz/graph plot with attributes

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 shapes 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 41.

Nodes marked by run length and run count: see fig. 19 on the facing page.

Input

```
#6 6
library(Rgraphviz)
plot_graphNEL_sprof(sprof04, nodeattrs="runs")
```

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.

The bottom line is: graph layout is best left to graph representation packages. You can help by trimming and surgery. As far as attributes are concerned, the current recommendation is to start with run time lengths and run time frequencies by node and level:

```
plot_graphNEL_sprof( x, nodeattrs = "runs")
```

Here for comparison is a run-down of the previous stages of our example, using this plot.

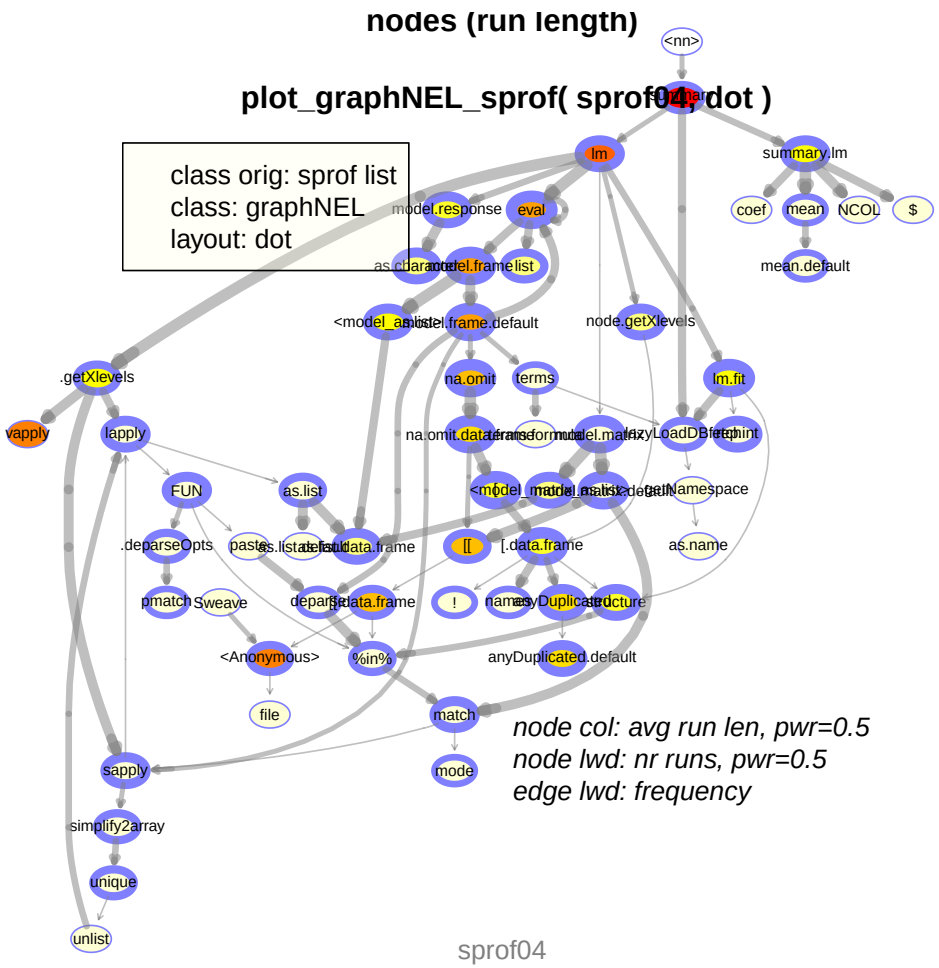
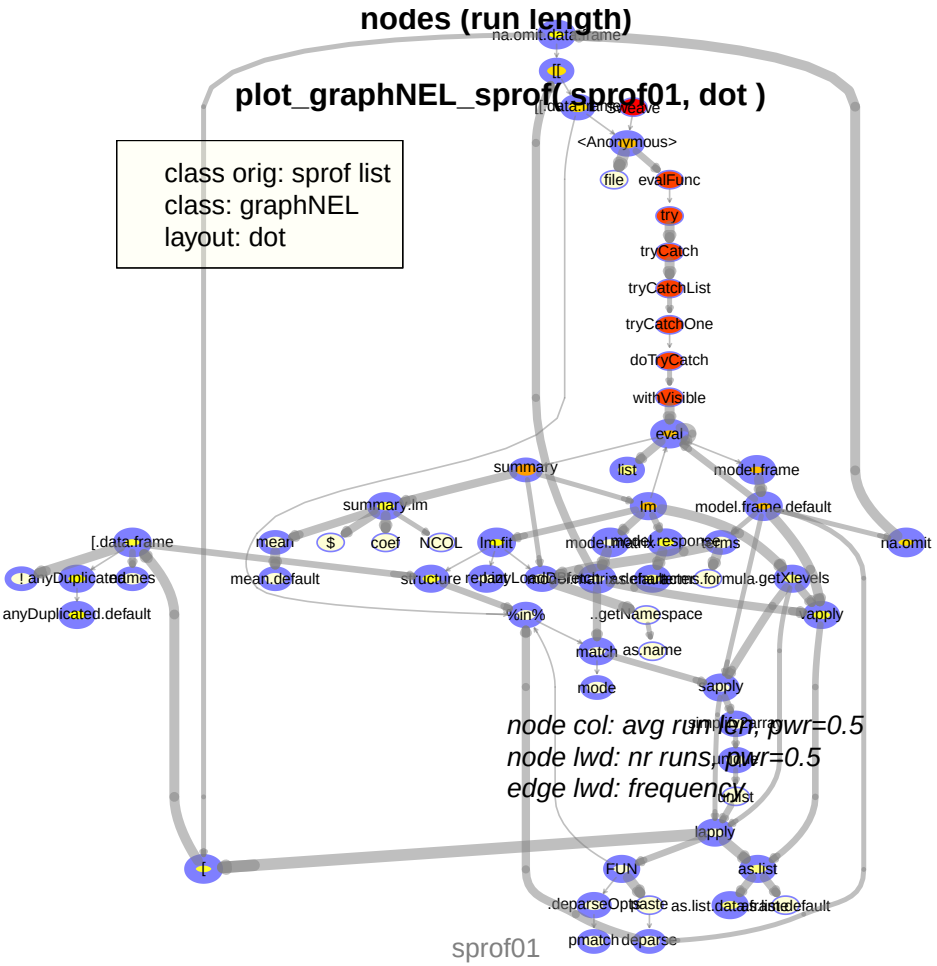


FIGURE 19. nodes (run length)

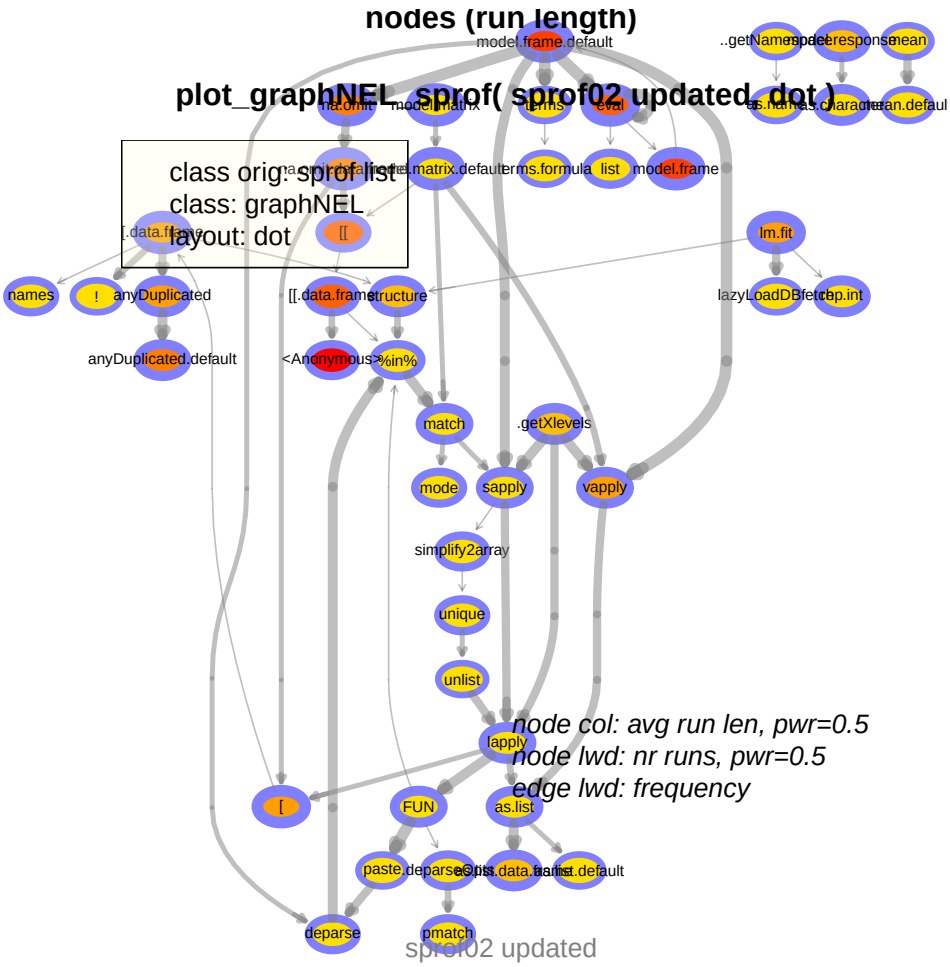
Input

```
plot_graphNEL_sprof(sprof01, nodeattrs="runs")
```


nodes (run length)
plot_graphNEL_sprof(sprof01, dot)
class orig: sprof list
class: graphNEL
layout: dot
node col: avg run length, pwr=0.5
node lwd: nr runs, pwr=0.5
edge lwd: frequency
sprof01

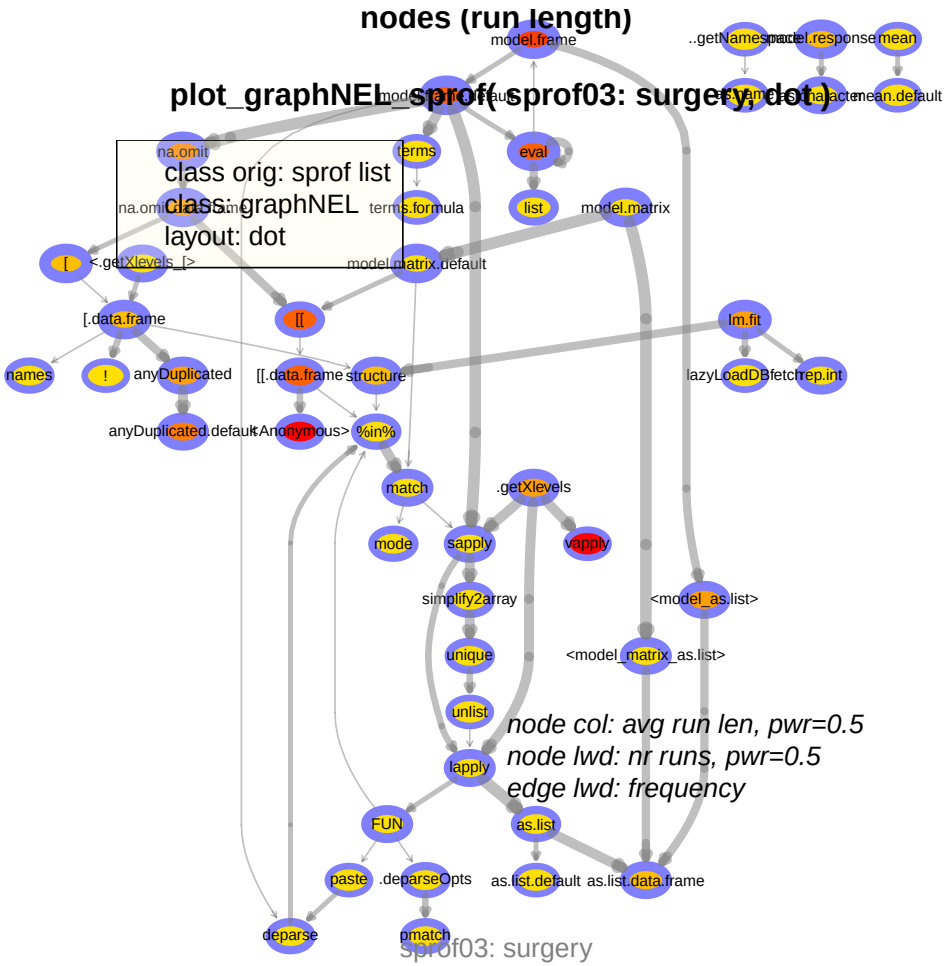
Input

```
plot_graphNEL_sprof(sprof02, nodeattrs="runs")
```



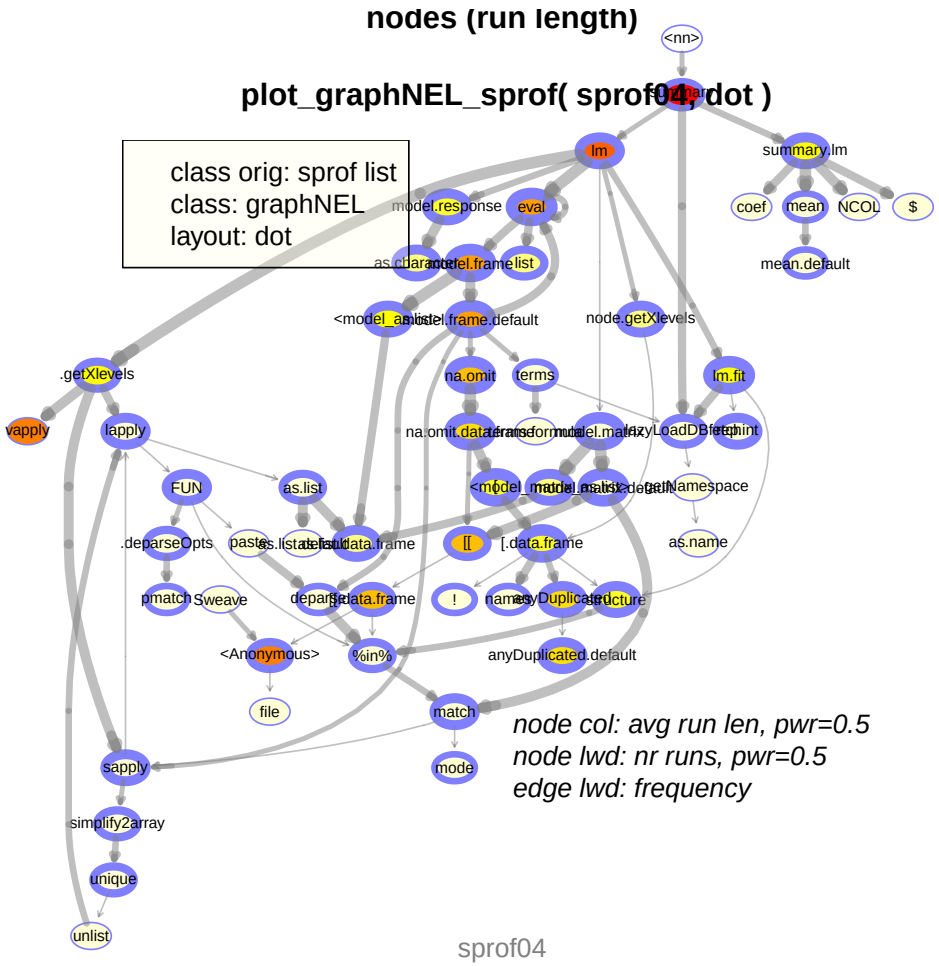
Input

```
plot_graphNEL_sprof(sprof03, nodeattrs="runs")
```



Input

```
plot_graphNEL_sprof(sprof04, nodeattrs="runs")
```



sprof04

Now it is time for detailed inspection. We just give hints how to start. Our example is very simple. In out example, we just use two functions, `lm` and `summary`. We should not be surprised to find them at prominent position with dominating weight. The `lm` and `summary.lm` are clearly separated and can be inspected separately.

In the `lm` branch, `eval` is prominent. In the overall structure of R, `eval` is the central interpreter. So it is expected to be prominent in all records.

Sorry. This example may be outdated at the time you read it. This is from R *3.0.1. Chances are that it will be outdated soon when the* R *core people read it. 2013-08-31*

`.getXlevels` is surprising. We have a plain vanilla real numbers problem. Why should it occur at all, and why at a prominent place? It is inside `lm`. But since R is open source, we can inspect it. We find `.getXlevels` at exactly one place near the end of the source of `lm`:

```
z$xlevels <- .getXlevels(mt, mf)
```

`z` is returned as result, and the help file tells us:

`xlevels` *(only where relevant) a record of the levels of the factors used in fitting.*

In a pure regression problem, it is not relevant. But there is no conditional code at this place. `xlevels` is always calculated (at the cost of some time, and always returning a named list of length 0). Ok. So we found a point where someone could look for an improvement.

We leave additional steps as an exercise. For example: look at the handling of NAs. Why (and where) do we spend time to handle NAs in a problem where there are no NAs at all?

4. Standard output

For a reference, here are the standard functions.

Input
```
sprof <- sprof01
```

4.1. **Print.** We omit the (lengthy) print output here and just give the commands as a reference.

Input
```
print_nodes(sprof)
```

Input
```
print_stacks(sprof)
```

Input
```
print_profiles(sprof)
```

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

4.2. **Summary.**

Input
```
summary_nodes(sprof)
```

Input
```
summary_stacks(sprof)
```

Input
```
summary_profiles(sprof)
```

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*.

Input
```
plot_nodes(sprof)
```

Input
```
plot_stacks(sprof)
```

Input
```
plot_profiles(sprof)
```

The *plot.sprof()* method for *sprof* objects concatenates these three functions, see fig. 20. Using the plot functions above allows better control and will be preferred. *shownodes()* may be a sufficient summary, see fig. 10 on page 29.

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

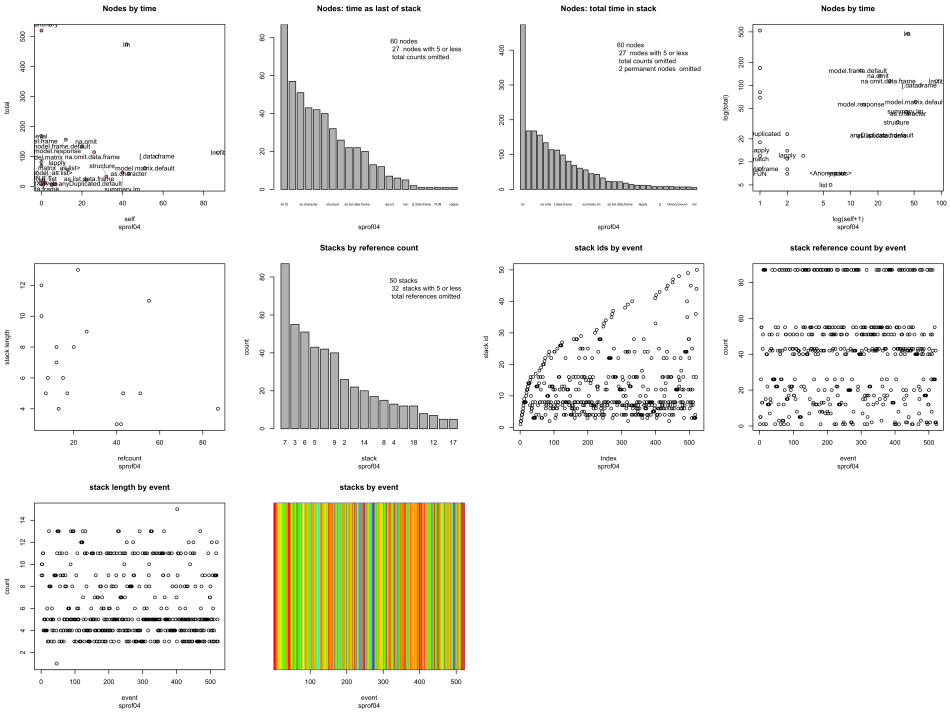


FIGURE 20. *plot.sprof(sprof04)*

5. More Graphs

Note: This section is collecting experiments with various graph packages. So far, none of the experiments looks too promising. This section 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.

ToDo: by graph package: preferred input format?

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*** in the next statement with your profile information.

ToDo: use attributes. Edge width should be easy.

Input
```
sprof <- sprof02
```

ToDo: include information from stack connectivity.

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

Input
```
sprofadj02 <- 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
```
sprofedgel02 <- edgedf(sprofadj02)
sprofedgel02$lwd  <- rkindex(sprofedgel02$count,
        maxindex=12,
        ties.method="min")
```

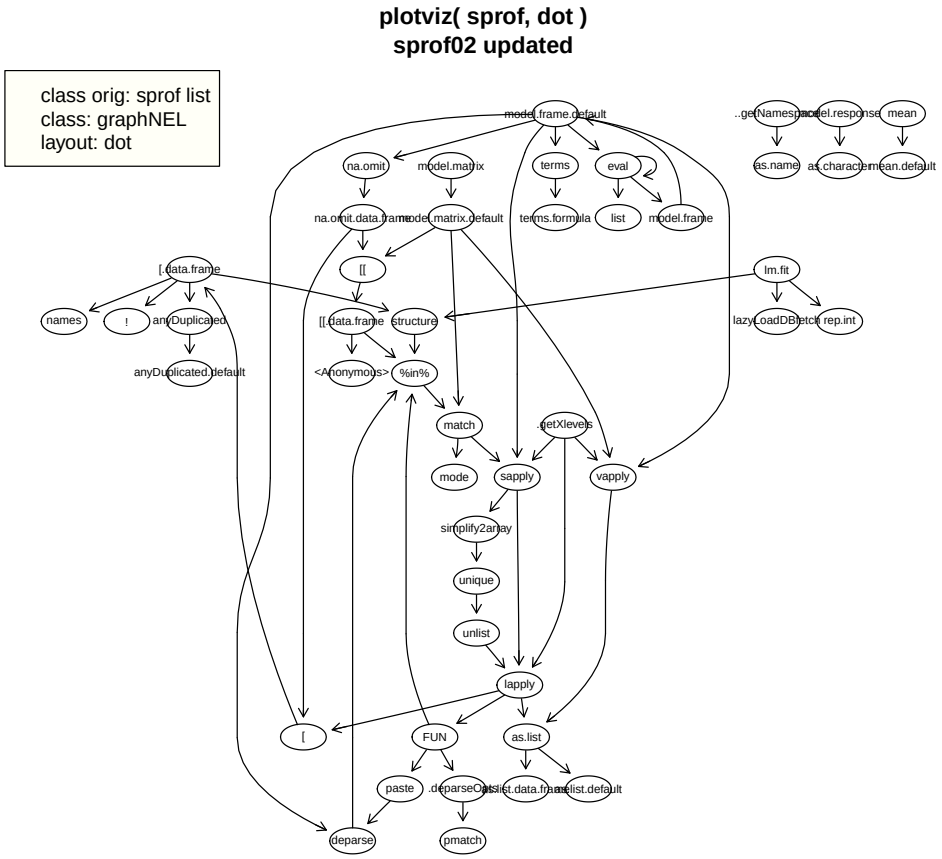
ToDo: add usage of sprofedgel02

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`.

Input

```
library(graph)
plotviz(sprof)
```



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)` The output of `igraph` gives problems when rendered with Adobe Acrobat on OS X Maverick. These examples have been moved to the `demo` section. You can call them using

`demo(topic=<Item>, package = "sprof")`.

Input

```
spdemo <- demo(,"sprof")
prxt(spdemo$results[, c("Item","Title")])
```

	Item	Title
1	igraphFrRein	igraph Fruchtermann-Reingold layout
2	igraphFrReingrid	igraph Fruchtermann-Reingold grid layout
3	igraphKK	igraph kamada kawai layout
4	igraphKK2	igraph kamada kawai layout alternate
5	igraphauto	igraph auto layout
6	igraphcircle	igraph circle layout
7	igraphgraphopt	igraph graphopt layout
8	igraphlgl	igraph lgl layout
9	igraphrandom	igraph random layout
10	igraphsphere	igraph sphere layout
11	igraphspring	igraph spring layout
12	igraphsvd	igraph svd layout

ToDo: propagate

5.1.3. *network package.*

Input

```
library(network)
as_network_sprof <- function(sprof) {
        sprofadj02 <- adjacency(sprof)
        nwsprof02 <- as.network(sprofadj02)
        network.vertex.names(nwsprof02) <-
                rownames(sprofadj02) # not honoured by plot
        return(nwsprof02)
}
```

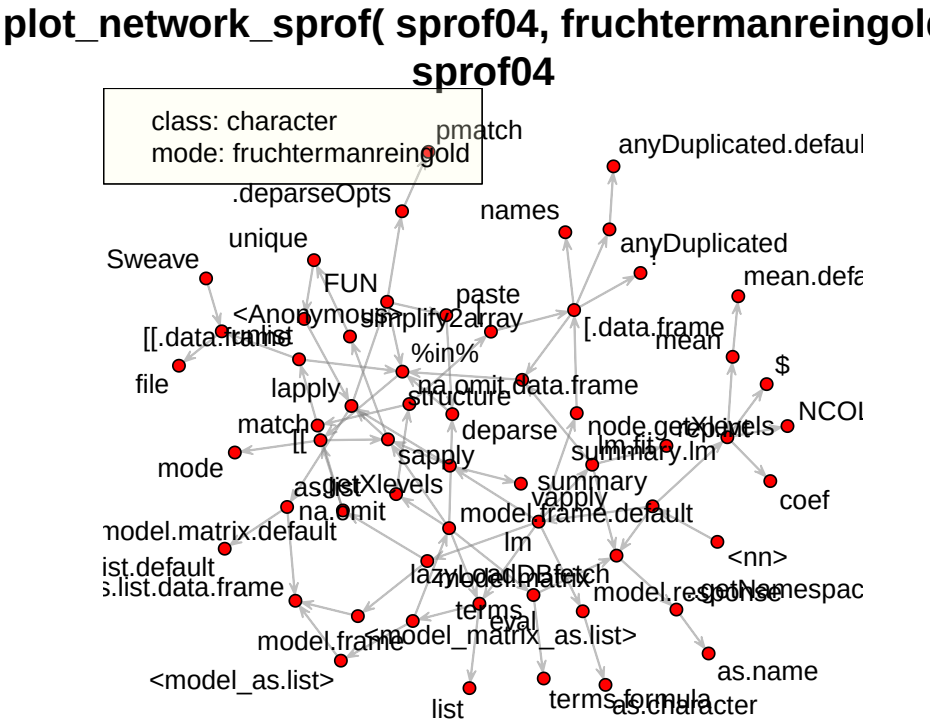
Input

```
plot_network_sprof <- function( nwsprof,
        mode = "fruchtermanreingold",
        main=NULL, label=NULL, sub=NULL,...)
{         classnwsprof <- class(nwsprof)
        xid <- deparse(substitute(nwsprof))
        xsubid <- NULL

        if (inherits(nwsprof, "sprof")) {
                xsubid <- nwsprof$info$id
                nwsprof <-  as_network_sprof(nwsprof)
                }
        if (!is.null(label)) {
                warning("plot_network_sprof: explicit label supplied, but will use vertex names")
                strx(label)
                strx(network.vertex.names(nwsprof))
                }

        if (!is.null(sub)) sub <- as.character(sub)
                main <-
                        paste0(main,
                                "\n plot_network_sprof( ", xid ,", ", mode, " )\n",
                                xsubid)

        plot( nwsprof,
                label = network.vertex.names(nwsprof),
                main= main,
                mode = mode,
                edge.len=2,
                edge.col="#80808080",
                sub= sub,
                cex.main=1.5,...)

        legend("topleft",
                legend=c(
                        paste0("class: ",paste(class(classnwsprof), collapse=" ")),
                        paste0("mode: ",mode)),
                bg="#FFFFE040",
                seg.len=0
```

```
                    )
}
```

Input

```
#6 6
plot_network_sprof(sprof04)
```

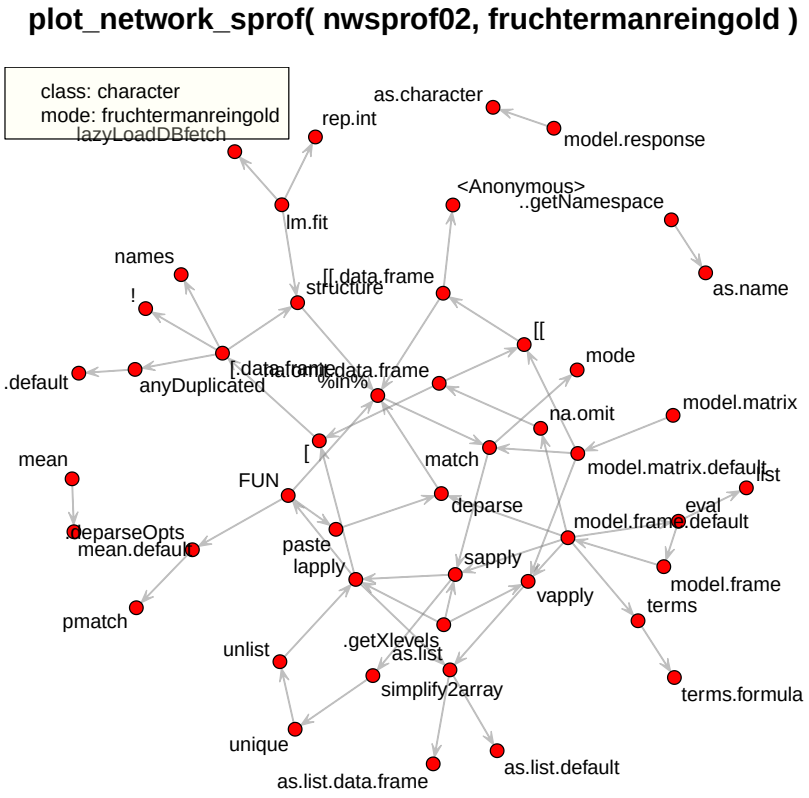


Input

```
nwsprof02 <- as.network(sprofadj02) # names is not imported
plot_network_sprof(nwsprof02, label=rownames(sprofadj02))
```

Output

```
##strx: label
chr [1:47] "!" "..getNamespace" ".deparseOpts" ...
##strx: network.vertex.names(nwsprof)
chr [1:47] "!" "..getNamespace" ".deparseOpts" ...
```



ToDo: maximum edge.lwd?

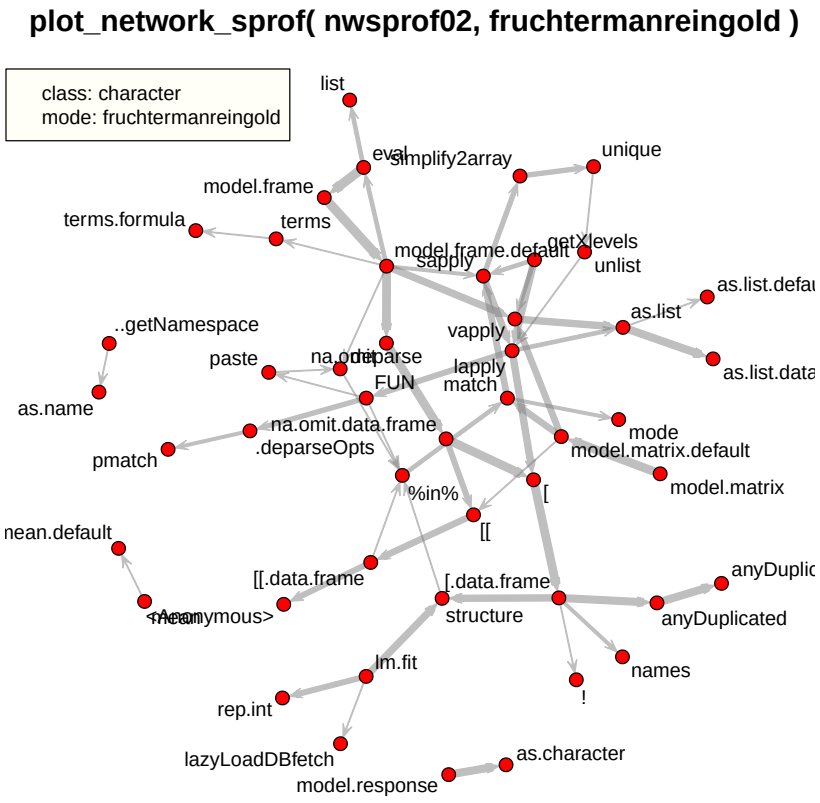
Experiments to include weight.

Input

```
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)
```

Output

```
##strx: label
chr [1:47] "!" "..getNamespace" ".deparseOpts" ...
##strx: network.vertex.names(nwsprof)
chr [1:47] "!" "..getNamespace" ".deparseOpts" ...
```



5.1.4. *Rgraphviz package.* Package 'Rgraphviz' was removed from the CRAN repository.

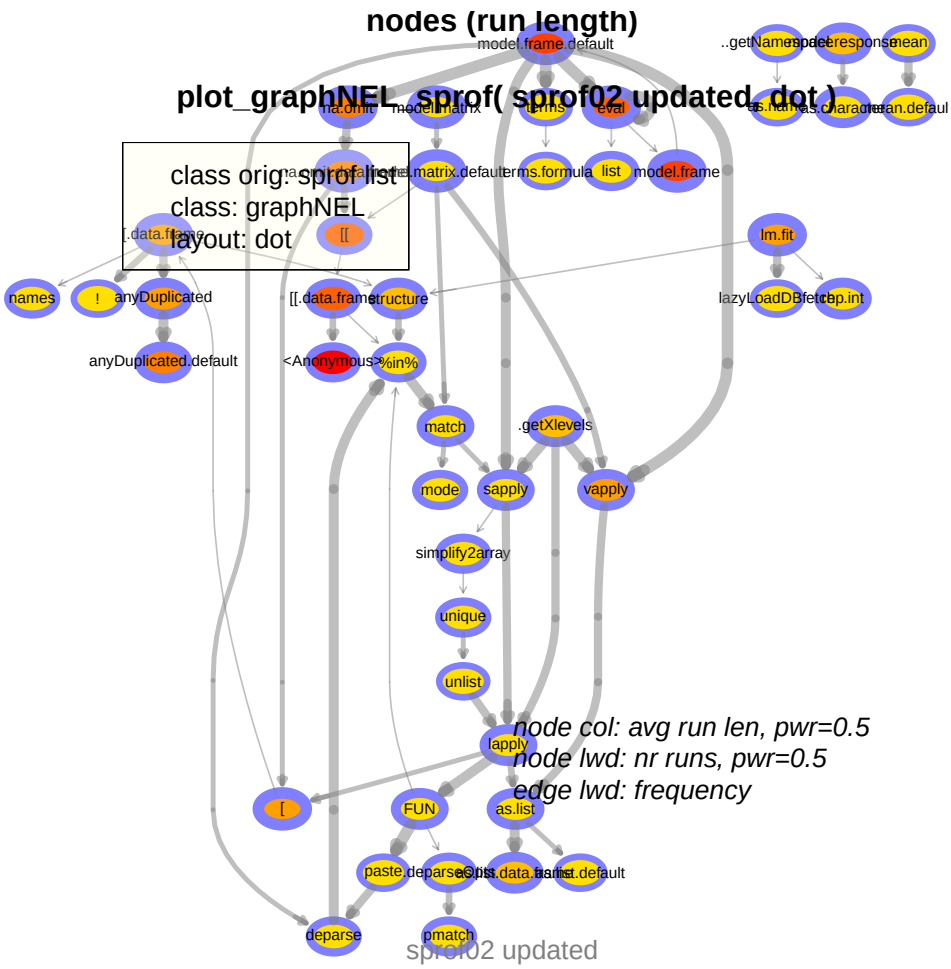
This package is now available from Bioconductor only, see `http://www.bioconductor.org/packages/release/bioc/html/Rgraphviz.html`.

Input

```
library(Rgraphviz)
```

Input

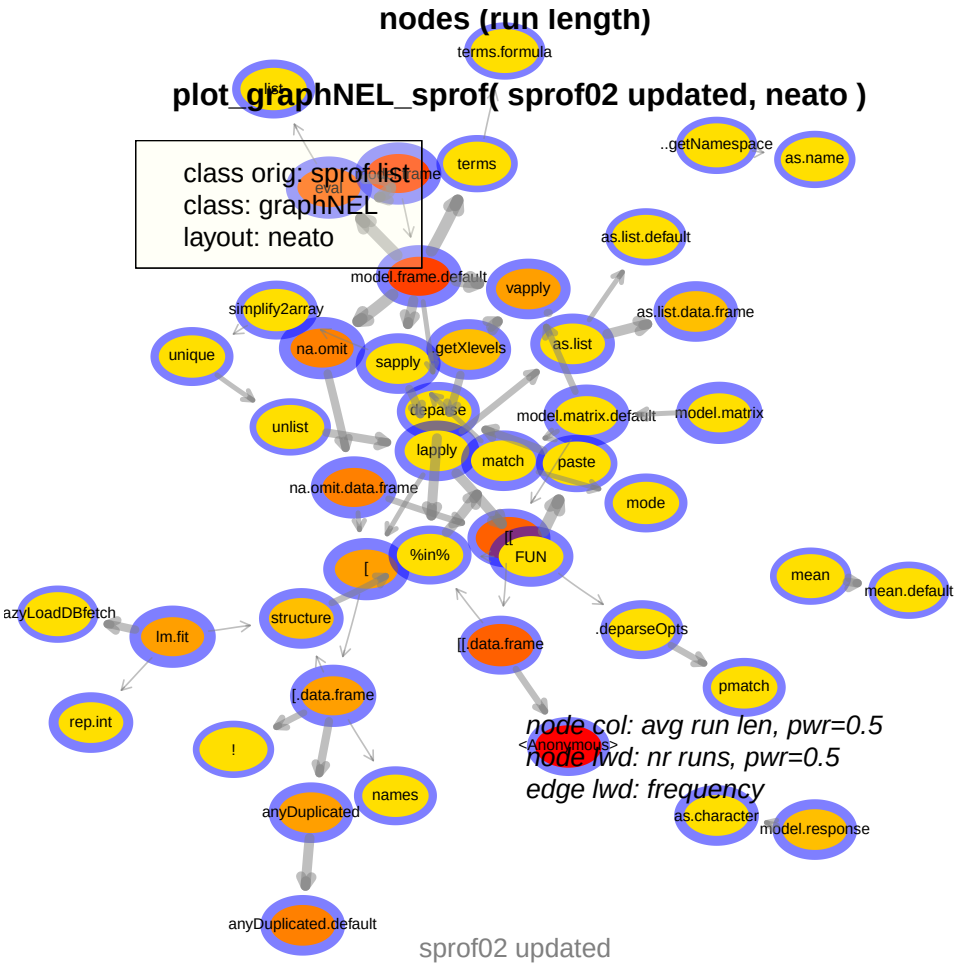
```
# 6 6
plot_graphNEL_sprof(sprof02,  layoutType="dot", nodeattrs="runs")
```



node col: avg run len, pwr=0.5
node lwd: nr runs, pwr=0.5
edge lwd: frequency

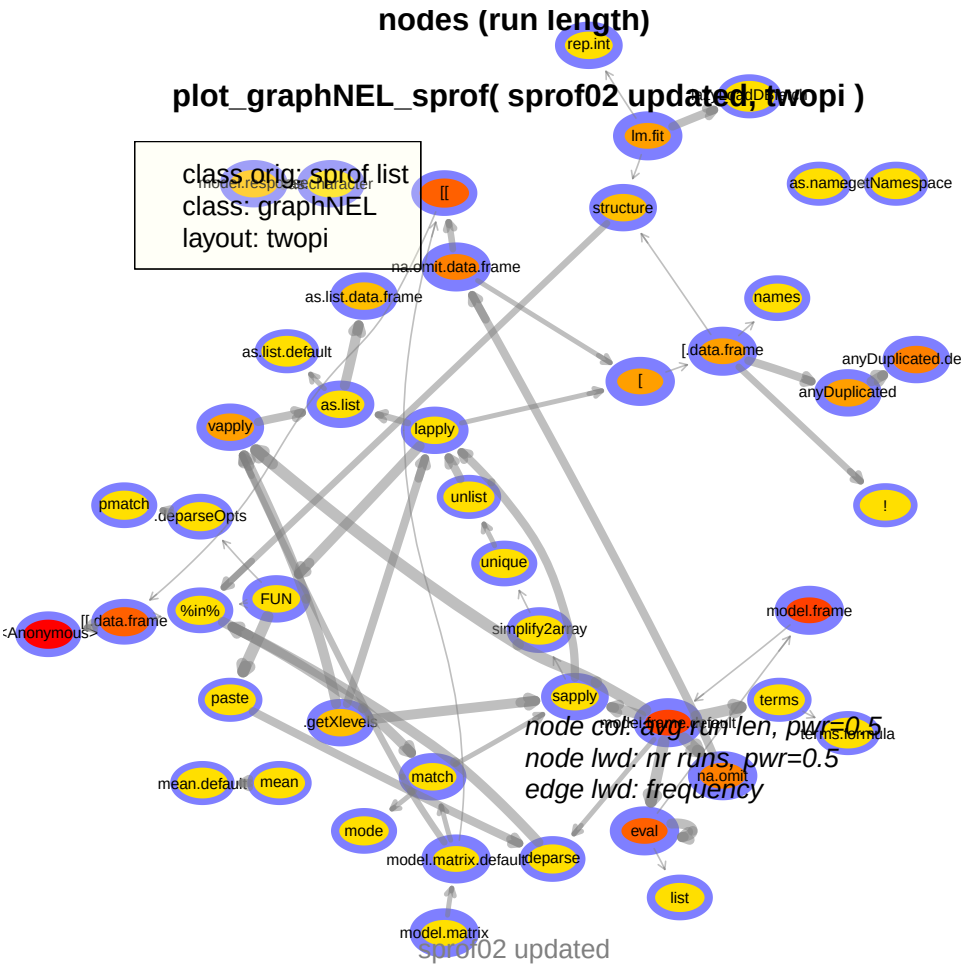
Input

```
#6 6
plot_graphNEL_sprof(sprof02,
        layoutType="neato", nodeattrs="runs", sub=sprof$info$id)
```



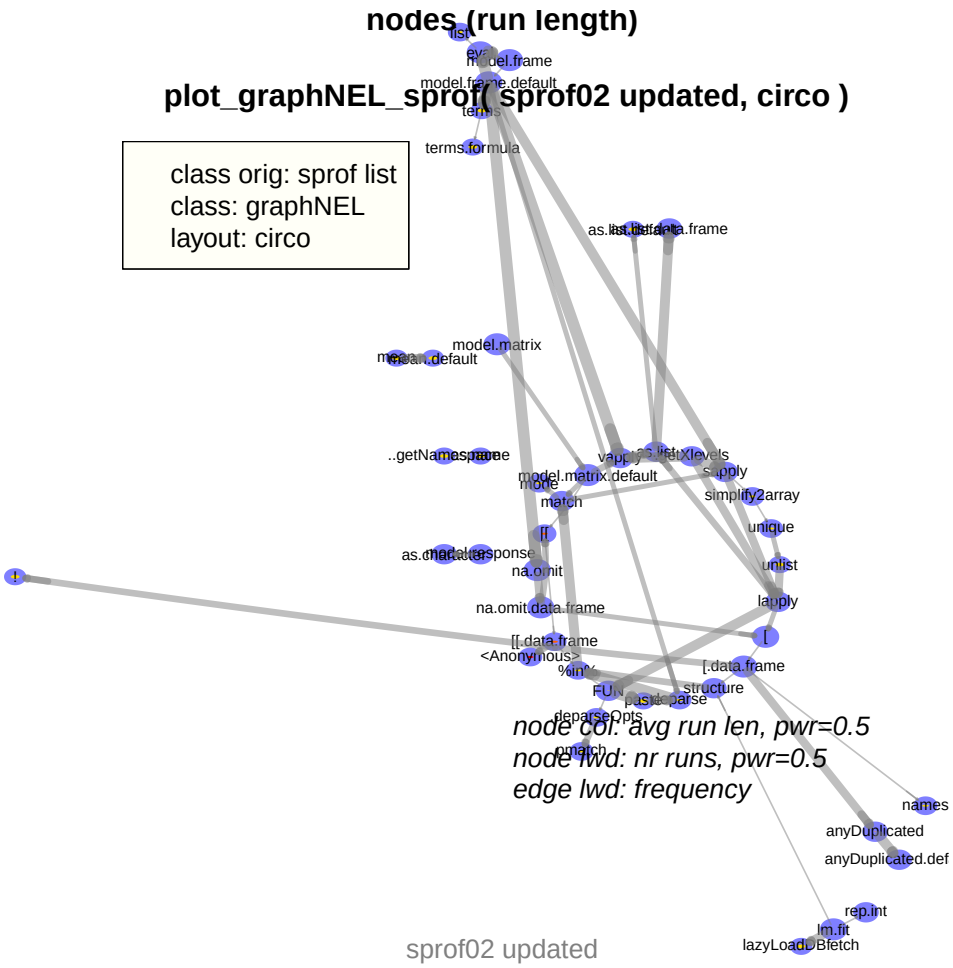
Input

```
#6 6
plot_graphNEL_sprof(sprof02,
        layoutType="twopi", nodeattrs="runs", sub=sprof$info$id)
```



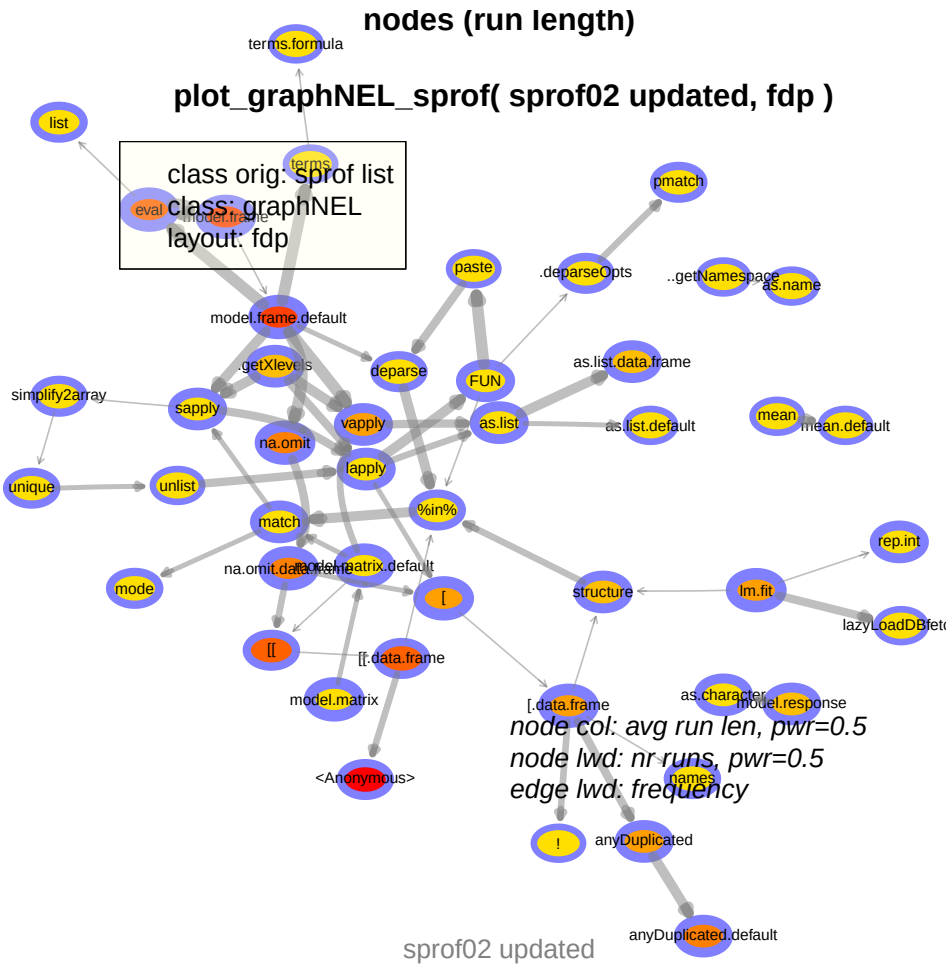
Input

```
#6 6
plot_graphNEL_sprof(sprof02,
        layoutType="circo", nodeattrs="runs", sub=sprof$info$id)
```



Input

```
#6  6
plot_graphNEL_sprof(sprof02,
        layoutType="fdp", nodeattrs="runs", sub=sprof$info$id)
```



6. Template

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

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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/C/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, sprof 0.1-0, wordcloud 2.4, xtable 1.7-1
- Loaded via a namespace (and not attached): BiocGenerics 0.6.0, 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:

`$HeadURL: svnssh://gsawitzki@svn.r-forge.r-project.org/svnroot/sintro/pkg/sprof/vignettes/sprofiling.Rnw +`

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

`$Id: sprofiling.Rnw 238 2013-09-08 17:22:45Z gsawitzki $`

`$Revision: 238 $`

`$Date: 2013-09-08 19:22:45` $0200(Sun, 08Sep2013)+$

`$Name: $`

`$Author: gsawitzki $`

7. XXX – LOST & FOUND

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.

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