# proto: An R Package for Prototype Programming

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#### Abstract

proto is an R package which facilitates a style of programming known as prototype programming. Prototype programming is a type of object oriented programming in which there are no classes. proto is simple yet retains the object oriented features of delegation (the prototype counterpart to inheritance) and object oriented dispatch. proto can be used to organize the concrete data and procedures in statistical studies and other applications without the necessity of defining classes while still providing convenient access to an object oriented style of programming. Furthermore, it can be used in a class-based style as well. The key design goals of the package are to integrate into R while providing nothing more than a thin layer on top of it. This paper describes how these goals are achieved within proto and provides examples of the use of the "proto" package.

Keywords: prototype programming, delegation, inheritance, clone, object orientated, S3, R.

## 1. Introduction

## 1.1. Object Oriented Programming in R

The R system for statistical computing (R Development Core Team 2005, http://www.R-project.org/) ships with two systems for object oriented programming referred to as S3 and S4. With the increased interest in object oriented programming within R over the last years additional object oriented programming packages emerged. These include the R.oo package (Bengtsson 2003) and the OOP package (Chambers and Lang 2001, http://www.omegahat.org/00P/). All these packages have the common thread that they use classes as the basis of inheritance. When a method is sent to an object the class of the object is examined and that class determines the specific function that is executed. In prototype programming there are no classes making it simple yet it retains much of the power of class-based programming. In the fact, proto is so simple that there is only one significant new routine name, proto. The other routines are just the expected support routines such as as.proto to coerce objects to proto objects, \$ to access and set proto object components and is.proto to check whether an object is a proto object. In addition, dot.proto will generate a graphical ancestor tree showing the parent-child relationships among generated proto objects.

### 1.2. History

The concept of prototype programming (Lieberman 1986; Taivalsaari 1996; Noble, Taivalsaari, and Moore 1999) has developed over a number of years with the Self language (Agesen, Bak, Chambers, Chang, Hölzle, Maloney, Smith, and Ungar 1992) being the key evolved programming language to demonstrate the concept. In statistics, the Lisp-based LispStat programming language (Tierney 1990) was the first and possibly only statistical system to feature prototype programming.

Despite having been developed over 20 years ago, and some attempts to enter the mainstream (e.g. Newtonscript on the Newton computer, which is no longer available, and Javascript where it is available but whose domain of application largely precluses use of prototype programming) prototype programming is not well known due to lack of language support in popular programming

languages such as C and Java. It tends to be the domain of research languages or Lisp.

Thus the the availability of a popular language, R <sup>1</sup>, that finally does provide the key infrastructure is an important development.

This work grew out of the need to organize multiple scenarios of model simulations in ecological modelling (Petzoldt 2003) and was subsequently generalized to the present package. A number of iterations of the code, some motivated by the ever increasing feature set in R, resulted in a series of utilities and ultimately successive versions of an R package developed over the last year. An initial version used R lists as the basis of the package. Subsequently the package was changed to use R environments. The first version to use environments stored the receiver object variable in a proxy parent environment which was created on-the-fly at each method call. The present version of proto passes the receiver object through the argument list, while hiding this from the caller. It defines the proto class as a subclass of the S3 environment class so that functionality built into R for the environment class is automatically inherited by the proto class.

#### 1.3. Benefits

The key benefit of the **proto** package is to provide access to a style of programming that has not been conveniently accessible within R or any other mainstream language today.

One particularly important use of the **proto** package in statistical applications is to wrap the code and data for each run of a particular study into an object. This provides a desirable level of organization yet gives the benefits of inheritance and object oriented dispatch without the overhead of having to create classes. Such concrete applications are the key domain of prototype programming. In contrast abstractions, such as a time series class, a matrix class, etc. are well suited to class-based programming although **proto** can be used in such applications as well since, as shown later, it is powerful enough to encompass class-based object-oriented programming even though classes are not a primitive notion in **proto**.

### 1.4. Overview

It is assumed that the reader has some general familiarity with object oriented programming concepts and with R.

The paper will proceed primarily by example focusing on illustrating the package proto through such demonstration. The remainder of the paper is organized as follows: Section 2 explains how "proto" objects are created and illustrates the corresponding methods for setting and getting components. It further discusses how object oriented delegation (the prototype programming analogue of inheritance) is handled and finally discusses the internals of the package. This section uses small examples chosen for their simplicity in illustrating the concepts. In Section 3 we provide additional examples of prototype programming in action. Three examples are shown. The first involves smoothing of data. Secondly we demonstrate the calculation of correlation confidence intervals using classical (Fisher Transform) and modern (bootstrapping) methods. Lastly we demonstrate the development of a binary tree as would be required for a dendogram. Section 4 gives a few summarizing remarks. Finally, an appendix provides a reference card that summarizes the functionality contained in proto in terms of its constituent commands.

 $<sup>^{1}</sup>$ Some indications of the popularity of R are the high volume mailing lists, international development team, the existence of over 400 addon packages, conferences and numerous books and papers devoted to R.

## 2. The class "proto" and its methods

## 2.1. Creation of "proto" objects

In this section we shall show, by example, the creation of two prototype objects and related operations. The simple idea is that each "proto" object is a set of components: functions (methods) and variables, which are tightly related in some way.

A prototype object is created using the constructor function **proto** (see Appendix A at the end of this paper or **proto** package online help for complete syntax of commands):

```
addProto <- proto( x = rnorm(5), add = function(.) sum(.$x) )</pre>
```

In this simple example, the proto function defines two components: a variable x and a method add. The variable x is a vector of 5 numbers and the method sums those numbers. As shown with the add method in this example, formal argument lists of methods must always have a first argument of dot (i.e. .) which signifies the object on which the method is operating. The dot refers to the current object in the same way that a dot refers to the current directory in UNIX. Within the method one must refer to other variables and methods in the object by prefacing each with .\$. For example, in the above we write sum(.\$x). Finally, note that the data and the method are very closely related. Such close coupling is important in order to create an easily maintained system.

To illustrate the usage of proto, we first load the package and set the random seed to make the examples in this paper exactly reproducible.

```
> library(proto)
> set.seed(123)
```

Then, we create the proto object from above and call its add method.

```
> addProto <- proto(x = rnorm(5), add = function(.) sum(.$x))
> addProto$add()
[1] 0.9678513
```

We also create another object, addProto2 with a different x vector and invoke its add method too.

```
> addProto2 <- addProto$proto(x = 1:5)
> addProto2$add()
```

[1] 15

In the examples above, we created a prototype object addProto and then called its add method as just explained. The notation addProto\$add tells the system to look for the add method in the addProto object. In the expression addProto\$add, the proto object to the left of the dollar sign, addProto here, is referred to as the receiver object. This expression also has a second purpose which is to pass the receiver object implicitly as the first argument of add. Note that we called add as if it had zero arguments but, in fact, it has one argument because the receiver is automatically and implicitly supplied as the first argument. In general, the notation object\$method(arguments) is used to invoke the indicated method of the receiver object using the object as the implicit first argument along with the indicated arguments as the subsequent arguments. As with the addProto example, the receiver object not only determines where to find the method but also is implicitly passed to the method through the first argument. The motivation for this notation is to relieve the user of specifying the receiver object twice: once to locate the method in the object and a second

time to pass the object itself to the method. The \$ is overloaded by the proto class to automatically do both with one reference to the receiver object. Even though, as with the addProto example, the first first argument is not listed in the call it still must be listed among the formal arguments in the definition of the method. It is conventional to use a dot . as the first formal argument in the method/function definition. That is, we call add using addProto\$add() displaying zero arguments but we define add in addProto displaying one argument add <- function(.), the dot.

In this example, we also created a second object, addProto2, which has the first object, addProto as its parent. Any reference to a component in the second object that is unsuccessful will cause search to continue in the parent. Thus the call addProto2\$add() looks for add in addProto2 and not finding it there searches its parent, addProto, where it is, indeed, found. add is invoked with the receiver object, addProto2, as the value of dot. The call addProto2\$add() actually causes the add in addProto to run but it still uses the x from addProto2 since dot (.) is addProto2 here and add references .\$x. Note that the reference to .\$x in the add found in addProto does not refer to the x in addProto itself. The x in addProto2 has overridden the x in its parent. This point is important so the reader should take care to absorb this point.

This simple example already shows the key elements of the system and how delegation (the prototype programming term for inheritance) works without classes.

We can add new components or replace components in an object and invoke various methods like this:

```
> addProto2$y <- seq(2, 10, 2)
> addProto2$x <- 1:10
> addProto2$add3 <- function(., z) sum(.$x) + sum(.$y) + sum(z)
> addProto2$add()

[1] 55
> addProto2$add3(c(2, 3, 5))

[1] 95
> addProto2$y

[1] 2 4 6 8 10
```

In this example, we insert variable y into the object addProto2 with a value of seq(2,10,2), reset variable x to a new value and insert a new method, add3. Then we invoke our two methods and display y. Again, note that in the case of protoAdd2\$add the add method is not present in protoAdd2 and so search continues to the parent addProto where it is found.

### 2.2. Internals

So far, we have used simple examples to illustrate the basic manipulation of objects: construction, getting and setting components and method invocation. We now discuss the internals of the package and how it relates to R constructs. proto is actually an S3 class which is a subclass of the environment class. Every proto object is an environment and its class is c("proto", "environment"). The \$ accessor is similar to the same accessor in environments except it will use the R get function to search up parent links if it cannot otherwise find the object (unlike environments). When accessing a method, \$ automatically supplies the first argument to the method unless the object is that or super. that is a special variable which proto adds to every proto object denoting the object itself. super is also added to every proto object and is the parent of that. that and super are normally used within methods of an object to refer to other components of the same or parent object, respectively, as opposed to the receiver (.). For example, suppose we want add in addProto2 to add the elements of x together and the elements of y together and then add these two sums. We could redefine add like this:

```
> addProto2$add <- function(.) super$add(.) + sum(.$y)</pre>
```

making use of the add already defined in the parent. One exception should be noted here. When one uses super, as above, or that to specify a method then the receiver object must be explicitly specified in argument one (since in those cases the receiver is possibly different than super or that so the system cannot automatically supply it to the call.)

Setting a value is similar to the corresponding operation for environments except that any function, i.e method, which is inserted has its environment set to the environment of the object into which it is being inserted. This is necessary so that such methods can reference that and super using lexical scoping.

#### 2.3. Traits

Let us look at the definition of a child object once again. addProto is the previously defined parent object and the expression addProto (x = 1:5) defines a child object of addProto and assigns it to variable addProto2a.

```
> addProto2a <- addProto$proto(x = 1:5)
> addProto2a$add()
[1] 15
```

That is, proto can be used to create a new child of an existing object by writing the parent object on the left of the \$ and proto on its right. Any contents to be added to the new child are listed in arguments of proto as shown.

For example, first let us create a class-like structure. In the following Add is an object that behaves very much like a class with an add method and a method init which constructs new objects. In the line creating object add1 the expression Add $\sinit(x = 1:5)$  invokes the init constructor of the receiver object Add. The method init has an argument of x = 1:5 which defines an x variable in the add1 object being instantiated. We similarly create another object add2.

```
> Add <- proto(add = function(.) sum(.$x), init = function(., x) .$proto(x = x))
> add1 <- Add$init(x = 1:5)
> add1$add()

[1] 15
> add2 <- Add$init(x = 1:10)
> add2$add()

[1] 55
```

An object which contains only methods and variables that are intended to be shared by all its children (as opposed to an object whose purpose is to have its own methods and variables) is known as a *trait*. It is similar to a class in class-based object oriented programming. Note that the objects add1 and add2 have the trait Add as their parent. We could implement subclass-like and superclass-like objects by simply defining similar trait objects to be the parent or child of Add. For example, suppose we want a class which calculates the sum of the logarithms of the data. We could define:

```
> Logadd <- Add$proto(logadd = function(.) log(.$add()))
> logadd1 <- Logadd$init(1:5)
> logadd1$logadd()
```

#### [1] 2.70805

Here the capitalized objects are traits. Logadd is a trait. It is a child of Add which is also a trait. logadd1 is an ordinary object, not a trait. One possible design is to create a tree of traits and other objects in which the leaves are ordinary objects and the remaining nodes are traits. This would closely correspond to class-based object oriented programming.

Note that the delegation of methods from one trait to another as in init which is inherited by Logadd from Add is nothing more than the same mechanism by which traits delegate methods to objects since, of course, traits are just objects no different from any other object other than by the conventions we impose on them. This unification of subclassing and instantiation beautifully shows the simplification that prototype programming represents.

#### 2.4. Utilities

The fact that method calls automatically insert the first argument can be used to good effect in leveraging existing R functions while allowing an object-oriented syntax.

For example, 1s() can be used to list the components of proto objects:

```
> addProto$ls()
[1] "add" "super" "that" "x"
Functions like:
> addProto$str()
> addProto$print()
> addProto$as.list()
```

show additional information about the elements. eapply can be used to explore more properties such as the length of each component of an object:

### > addProto\$eapply(length)

Another example of some interest in any object oriented system which allows multiple references to one single object is that object identity can be tested using the respective base function:

#### > addProto\$identical(addProto2)

#### [1] FALSE

It is important to notice here, that proto has no code that is specific to ls, str or any of the other ordinary R functions listed. We are simply making use of the fact that obj\$fun(...) is transformed into get("fun", obj)(obj, ...) by the proto \$ operator. For example, in the case of addProto\$ls() the system looks for ls in object addProto. It cannot find it there so it looks to its parent, which is the global environment. It does not find it there so it searches the remainder of the search path, i.e. the path shown by running the R command search(), and finally finds it in the base package, invoking it with an argument of addProto. Since all proto objects are also environments ls(addProto) interprets addProto as an environment and runs the ls command with it. In the ls example there were no arguments other than addProto, and even that one was implicit, but if there were additional arguments then they would be passed as shown in the eapply and identical examples above.

### 2.5. Plotting

The dot.proto function can be used to create output that is readable by the GraphViz dot program (Ganser and North 2000) to create visual representations of ancestor trees (figure 1). It takes two arguments, a proto object (or an environment) and a file name:

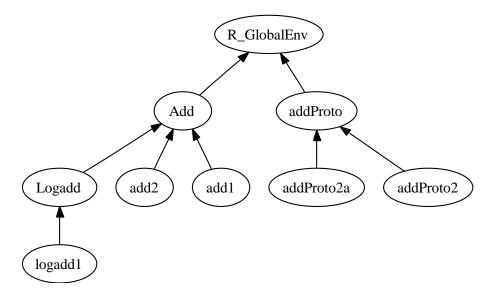


Figure 1: Ancestor tree generated using dot.proto. Edges point from child to parent.

```
> dot.proto(file = "test.dot")
> system("dot -Tps test.dot -o test.ps")
```

Both arguments can be omitted in which case the current environment and the standard output are the defaults.

## 3. Examples

## 3.1. Smoothing

In the following we create a proto object named oo containing a vector of data x (generated from a simulated autoregressive model) and time points tt, an intermediate result x.smooth, some plotting parameters codexlab, ylab, pch, col and three methods smooth, plot and residuals which smooth the data, plot the data and calculate residuals, respectively. We also define ..x.smooth which holds intermediate results. Names beginning with two dots prevent them from being delegated to children. If we override x in a child we would not want an out-of-sync x.smooth. Note that the components of an object can be specified using a code block in place of the argument notation we used previously in the proto command.

Having defined our proto object we can inspect it, as shown below, using print which is automatically invoked if the name of the object, oo, is entered on a line by itself. In this case, there is no proto print method so we inherit the environment print method which displays the environment hash code. Although it produces too much output to show here, we could have displayed a list of the entire contents of the object oo via oo\$as.list(all.names = TRUE). We can get a list of the names of the components of the object using oo\$ls(all.names = TRUE) and will look at the contents of one component, oo\$pch.

```
> 00
<environment: 016B806C>
attr(,"class")
[1] "proto"
                   "environment"
> oo$1s(all.names = TRUE)
 [1] "..x.smooth" "col"
                                 "pch"
                                               "plot"
                                                             "residuals"
 [6] "smooth"
                   "super"
                                 "that"
                                               "tt"
                                                             "x"
[11] "xlab"
                   "ylab"
> oo$pch
[1] "."
```

Let us illustrate a variety of manipulations. We will set up the output to plot 2 plots per screen using mfrow. We change the plotting symbol, smooth the data, invoke the plot method to display a plot of the data and the smooth and then plot the residuals in the second plot (figure 2).

```
> par(mfrow = c(1, 2))
> oo$pch <- 20
> oo$smooth()
> oo$plot()
> plot(oo$residuals(), type = "1")
```

Now let us illustrate the creation of a child object and delegation. We create a new child object of oo called oo.res. We will override the x value in its parent by setting x in the child to the value of the residuals in the parent. We will also override the pch and ylab plotting parameters. We will return to 1 plot per screen and run plot using the oo.res object as the receiver invoking the smooth and plot methods (which are delegated from the parent oo) with the data in the child (figure 3).

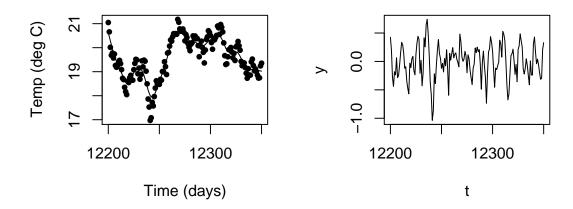


Figure 2: Data and smooth from oo\$plot() (left) and plot of oo\$residuals() (right).

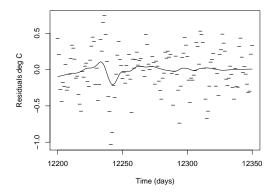


Figure 3: Output of oo.res\$plot(). oo.res\$x contains the residuals from oo.

```
> oo.res <- oo$proto(pch = "-", x = oo$residuals()$y, ylab = "Residuals deg C") > par(mfrow = c(1, 1)) > oo.res$smooth() > oo.res$plot()
```

Now we make use of delegation to change the parent and child in a consistent way with respect to certain plot characteristics. We have been using a numeric time axis. Let us interpret these numbers as the number of days since the Epoch, January 1, 1970, and let us also change the plot colors.

```
> oo$tt <- oo$tt + as.Date("1970-01-01")
> oo$xlab <- format(oo.res$tt[1], "%Y")
> oo$col = c("blue", "red")
```

We can introduce a new method, splot, into the parent oo and have it automatically inherited

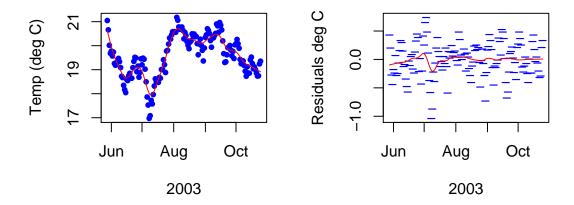


Figure 4: Plotting options and splot function applied to both parent (left) and child (right) object

by its children. In this example it smooths and then plots and we use it with both oo and oo.res (figure 4).

```
> oo$splot <- function(., ...) {
+    .$smooth(...)
+    .$plot()
+ }
> par(mfrow = c(1, 2))
> oo$splot(bass = 2)
> oo.res$splot()
```

Numerous possibilities exist to make use of the mechanisms shown, so one may create different child objects, apply different smoothing parameters, overwrite the smoothing function with a lowess smoother and finally compare fits and residuals.

Now lets change the data and repeat the analysis. Rather than overwrite the data we will preserve it in oo and create a child oos to hold an analysis with sinusoidal data.

```
> oos <- oo$proto(expr = {
+     tt <- seq(0, 4 * pi, length = 1000)
+     x <- sin(tt) + rnorm(tt, 0, 0.2)
+ })
> oos$splot()
```

Lets perform the residual analysis with oos. We will make a deep copy of oo.res, i.e. duplicate its contents and not merely delegate it, by copying oo.res to a list from which we create the duplicate, or cloned, proto object (figure 5 and 6):

```
> oos.res <- as.proto(oo.res$as.list(), parent = oos)
> oos.res$x <- oos$residuals()$y
> oos.res$splot()
```

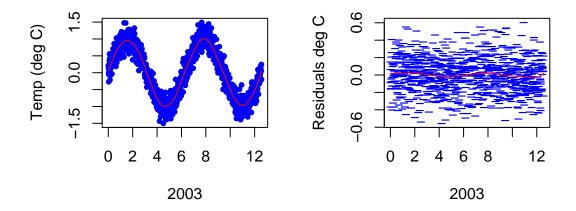


Figure 5: Smoothing of sinusoidal data (left) and of their residuals (right)

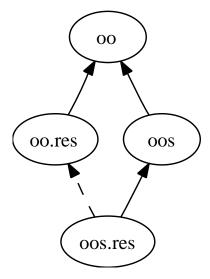


Figure 6: Cloning (dashed line) and delegation (solid line). Edges point from child to parent.

We have delegated variables and methods and overridden both. Thus, even with such a simple analysis, object orientation and delegation came into play. The reader can plainly see that smoothing and residual analysis were not crucial to the example and this example could be replaced with any statistical analysis including likelihood or other estimation techniques, time series, survival analysis, stochastic processes and so on. The key aspect is just that we are performing one-of analyses and do not want to set up an elaborate class infrastructure but just want to directly create objects to organize our calculations while relying on delegation and dispatch to eliminate redundancy.

## 3.2. Correlation, Fisher's Transform and Bootstrapping

The classical approach to confidence intervals for the correlation coefficient is to assume normality

of the underlying data and then use Fisher's transform to transform the correlation coefficient to an approximately normal random variable. Fisher showed that with the above assumption, transforming the correlation coefficient using the hyperbolic arc tangent function approximately yields a random variable distributed with an  $\frac{N(p,1)}{\sqrt{(n-3)}}$  distribution. The transformed random variable can be used to create normal distribution confidence intervals and the procedure can be back transformed to get confidence intervals for the original correlation coefficient.

The modern approach to confidence intervals for the correlation coefficient is to use bootstrapping. This does not require the assumption of normality of the underlying distribution and requires no special purpose theory devoted solely to the correlation coefficient,

Let us calculate the 95% confidence intervals using Fisher's transform first. We use GNP and Unemployed from the Longley data set. First we retrieve the data set and extract the required columns into x. Then we set n to the number of cases and pp to the percentiles of interest. Finally we calculate the sample correlation and create a function to calculate the confidence interval using Fisher's Transform. This function not only returns the confidence interval but also stores it in CI in the receiver object.

Now let us repeat this analysis using the bootstrapping approach. We derive a new object longley.ci.boot as child of longley.ci, setting the number of replications, N, and defining the procedure, ci which does the actual bootstrap calculation.

```
> longley.ci.boot <- longley.ci$proto({
+     N <- 2000
+     ci <- function(.) {
+         corx <- function(idx) cor(.$x[idx, ])[1, 2]
+         samp <- replicate(.$N, corx(sample(.$n, replace = TRUE)))
+         (.$CI <- quantile(samp, .$pp))
+    }
+ })</pre>
```

In the example code below the first line runs the Fisher Transform procedure and the second runs the bootstrap procedure. Just to check that we have performed sufficient bootstrap iterations we rerun it in the third line, creating a delegated object on-the-fly running its ci method and then immediately throwing the object away.

From the three runs we see that the bootstrap confidence intervals are narrower than the Fisher Transform intervals despite the reduced assumptions and no special purpose theory. The fact that 8,000 replications give roughly the same result as 2,000 replications satisfies us that we have used a sufficient number of replications.

```
> longley.ci$ci()
[1] 0.1549766 0.8464304
> longley.ci.boot$ci()
```

```
2.5% 97.5%

0.2413246 0.8228509

> longley.ci.boot$proto(N = 8000)$ci()

2.5% 97.5%

0.2551993 0.8298349
```

We now have the results stored in two objects nicely organized for the future. Note, again, that despite the simplicity of the example we have used the features of object oriented programming, coupling the data and methods that go together, while relying on delegation and dispatch to avoid duplication.

### 3.3. Dendograms

In Gentleman (2002) there is an example of creating a binary tree for use as a dendogram. Here we directly define a binary tree with no setup at all. To keep it short we will create a binary tree of only two nodes having a root whose left branch points to a leaf. The leaf inherits the value and incr components from the root. The attractive feature is that the leaf be defined as a child of the parent using that\$proto before the parent is even finished being defined. Compared to the cited \$\$4\$ example there is no need to create an extra class to introduce the required level of indirection.

tree is the root node of the tree. It has four components. A method incr which increments the value component, a .Name, the value component itself and the left branch ..left ..left is itself a proto object which is a child of tree. The leaf inherits the value component from its parent, the root. As mentioned, at the time we define ..left we have not even finished defining tree yet we are able to reference the yet to be defined parent via that as proto makes that available to inserted components as they are inserted.

Although this is a simple structure we could have embedded additional children into root and leaf and so on recursively making the tree or dendogram arbitrarily complex.

Let us do some computation with this structure. We display the value fields in the two nodes, increment the value field in the root and then display the two nodes again to show that the leaf changed too.

```
> cat("root:", tree$value, "leaf:", tree$..left$value, "\n")
root: 3 leaf: 3
> tree$incr(1)
> cat("root:", tree$value, "leaf:", tree$..left$value, "\n")
root: 4 leaf: 4
```

If we increment value in leaf directly (see the example below where we increment it by 10) then it receives its own copy of value so from that point on leaf no longer inherits value from root. Thus incrementing the root by 5 no longer increments the value field in the leaf.

```
> tree$..left$incr(10)
> cat("root:", tree$value, "leaf:", tree$..left$value, "\n")
root: 4 leaf: 14
> tree$incr(5)
> cat("root:", tree$value, "leaf:", tree$..left$value, "\n")
root: 9 leaf: 14
```

## 4. Summary

The package **proto** provides an S3 subclass of the **environment** class for constructing and manipulating object oriented systems without classes. It can also emulate classes even though classes are not a primitive structure. Its key design goals are to provide as simple and as thin a layer as practically possible while giving the user convenient access to this alternate object oriented paradigm. This paper describes, by example, how prototype programming can be carried out in R using **proto** and illustrates such usage. Delegation, cloning traits and general manipulation are all reviewed by example.

## Computational details

The results in this paper were obtained using R 2.1.0 with the package **proto** 0.2–3. R itself and the **proto** package are available from CRAN at http://CRAN.R-project.org/. The GraphViz software is available from http://www.graphviz.org.

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## A. Reference Card

#### Creation

prot

proto(., expr, envir, ...) embeds the components specified in expr and/or ... into the proto object or environment specified by envir. A new object is created if envir is omitted. The parent of the object is set to . . The parent object, ., defaults to the parent of envir or the current environment if envir is missing. expr and ... default to empty specifications. The returned object will contain that and super variables referring to the object itself and the parent of the object, respectively.

#### Coercion

as.proto

If x is a proto object or environment then x is returned as a proto object with the values of that and super inserted in the case of an environment or refreshed in the case of a proto object. If x is a list then additional arguments are available: as.proto(x, envir, parent, FUN, all.names, ...). Each component of x is copied into envir. envir may be an environment or proto object. If it is missing a new proto object is created. If all.names = FALSE then only list components whose names do not begin with a dot are copied. If FUN is specified then, in addition, only list components v for which FUN(v) is TRUE are copied. If parent is specified then the resulting proto object will have that parent. Otherwise, it will have the parent of envir if envir was specified. If neither are specified the parent defaults to the current environment.

#### Standard methods

- \$ obj\$x searches proto object obj for x. If the name x does not begin with two dots then ancestors are searched if the name is not found in obj. If x is a variable or if obj is super or that then x is returned. Otherwise, the call obj\$x(...) is equivalent to the call get("x", obj)(obj, ...). If it is desired to return a method as a value rather than in the context of a call then use get("x", obj) (or obj[["x"]] x is known to be directly in obj) rather than \$ syntax.
- \$<- obj\$x <- value sets x in proto object obj to value creating x if
  not present. If obj is super then a side effect is to set the parent of
  obj to value.</pre>

is.proto(x) returns TRUE if x is a proto object and othewise returns FALSE.

#### Utilities

dot.proto

dot.proto(e, file, control) creates a input file which for the GraphViz dot program which generates an ancestor tree among all proto objects in environment or proto object e. e defaults to the current environment and file defaults to the standard output. control is an optional list of display parameters. The components are include and arrow.from.child. include is a string of GraphViz dot commands to be included in the output. arrow.from.child causes arrows to point from children to parents if TRUE (default) and otherwise from parents to children.