Textual description of makecdfenv

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1 Introduction

The *makecdfenv* package is part of the Bioconductor¹ project. It can be used to create cdf environments from Affymetrix chip description files to be used with the package *affy*.

For many *CDF* files a pre-assembled package is available from http://www.bioconductor. org/data/cdfenvs/cdfenvs.html. The following notes explain what to do if you want to create your own.

2 Creating a CDF package

To install packages, you need certain tools besides R itself. They are generally available on most Unix platforms. Under MS-Windows, if you use the precompiled binary distribution of R from CRAN, you will need to install some extra tools, like the *source package installation files* and *Perl*. If you think this is too complicated, you can contribute the package to Bioconductor and we will create a windows binary for you. Alternatively, you may skip this section and proceed to section 3

Let's say your CDF file is called eggplantgenome.cdf and is in your working directory. To make a package, simply write

¹http://www.bioconductor.org/

R> make.cdf.package("eggplantgenome.cdf")

This will create a subdirectory eggplantgenomecdf in your working directory, which contains the package. Please consult the help page for make.cdf.package to find out about further options.

Now, open a terminal with an operating system shell, and write

> R CMD INSTALL eggplantgenomecdf

This will install the package into your R. You are now ready to use the affy package to process eggplant genechips.

3 Creating CDF environments

If you do not choose to use the package creation mechanism, you can still produce a data structure (in R lingo, it is called *environment*) that can be used by the affy package. Let's say your CDF file is called eggplantgenome.cdf and is in your working directory. Simply write

R> eggplantcdf = make.cdf.env("eggplantgenome.cdf")

Please consult the help page for make.cdf.env to find out about further options.

4 Naming of CDF packages and environments

What should you call your package or environment?

make.cdf.package chooses a default name by stripping all non-letters and nonnumbers from the CDF file name, and converting everything to lower case. In many cases, this will work.

You can obtain the CDF name that is associated with a CEL file through

R> pname <- cleancdfname(whatcdf("mycelfile.cel"))

Then call the package making function this way:

```
R> make.cdf.package("eggplantgenome.cdf", packagename=pname)
```

Similarly, if you work with the environments directly (as described in section 3) simply replace the word eggplantcdf with what ever the value of pname is.

In instances of AffyBatch, the cdfName slot gives the name of the appropriate CDF file for the arrays that are represented in the intensity slot. The functions read.celfile, read.affybatch, and ReadAffy extract the CDF filename from the CEL files that they read. The function cleancdfname converts Affymetrix' CDF name to the name that is used in Bioconductor. Here are two examples:

```
> cat("HG_U95Av2 is", cleancdfname("HG_U95Av2"), "\n")
HG_U95Av2 is hgu95av2cdf
> cat("HG-133A is", cleancdfname("HG-133A"), "\n")
HG-133A is hg133acdf
```

The method getCdfInfo takes as an argument AffyBatch and returns the appropriate environment. If x is an AffyBatch, this function will look for an environment with name cleancdfname(x@cdfName).

5 Location-ProbeSet Mapping

On Affymetrix GeneChip arrays, several probes are used to represent genes in the form of probe sets. From a *CEL* file we get for each physical location, or cel, (uniquely identified by its x and y coordinates) an intensity. The *CEL* file also contains the name of the *CDF* file needed for the location-probe-set mapping. The *CDF* files store the name of the probe set related to each location on the array. We store this mapping information in R environments, such as the ones produced by make.cdf.env or contained in the packages made by make.cdf.package.

In affy, the x and y coordinates are internally stored as one number i. The mapping between (x, y) and i is provided by the functions i2xy(i) and xy2i that are contained in each CDF package. They are very simple:

$$i = ys_x + x + 1$$

$$x = (i - 1) \%\% s_x$$

$$y = (i - 1) \%/\% s_x$$

where s_x is the side length of the chip (measured in number of probes) in x-direction.