

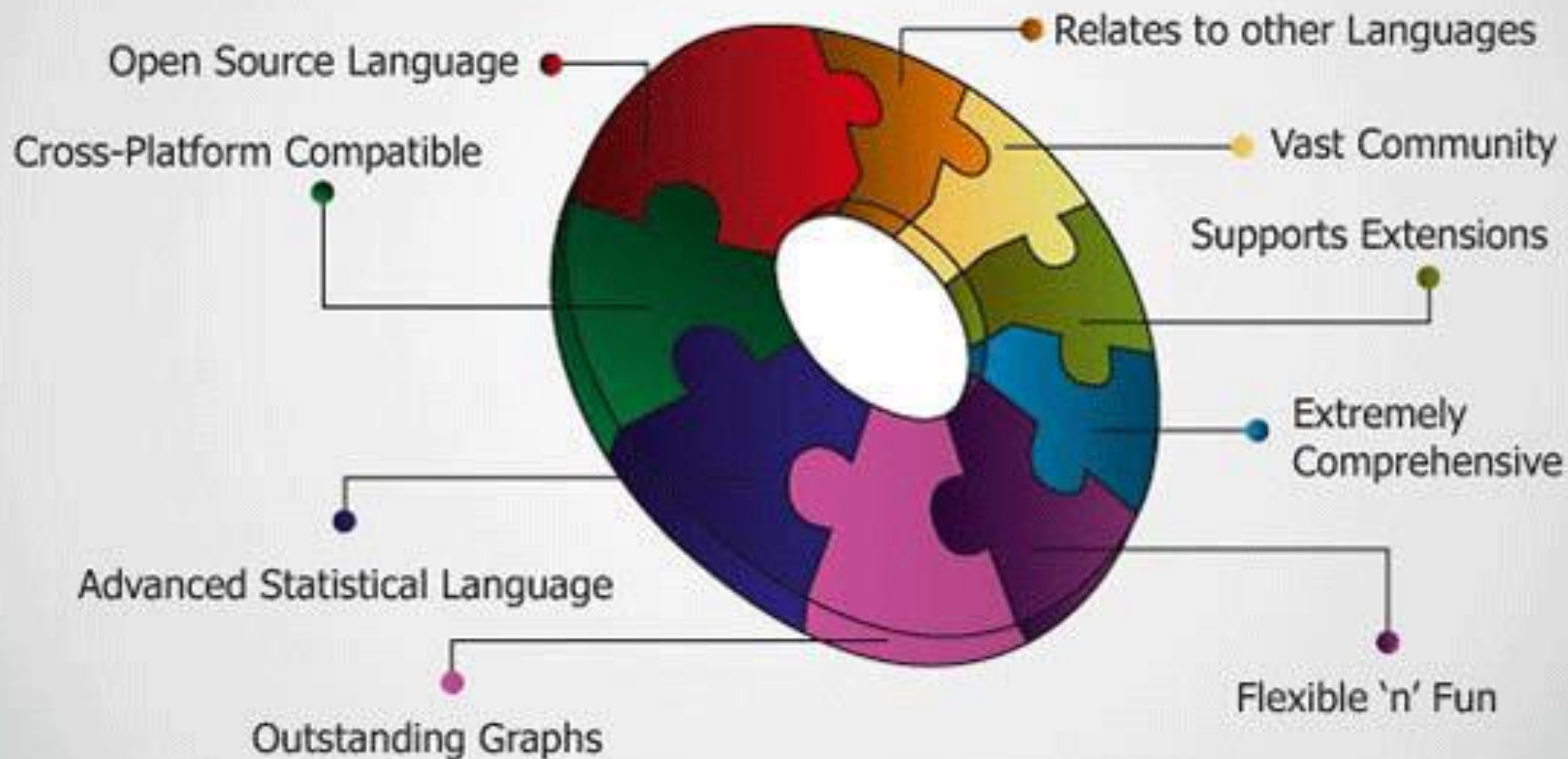
Some useful web content and information related to the programming language R

Magnus Fontes

May 2017

Why Learn R?

edureka!



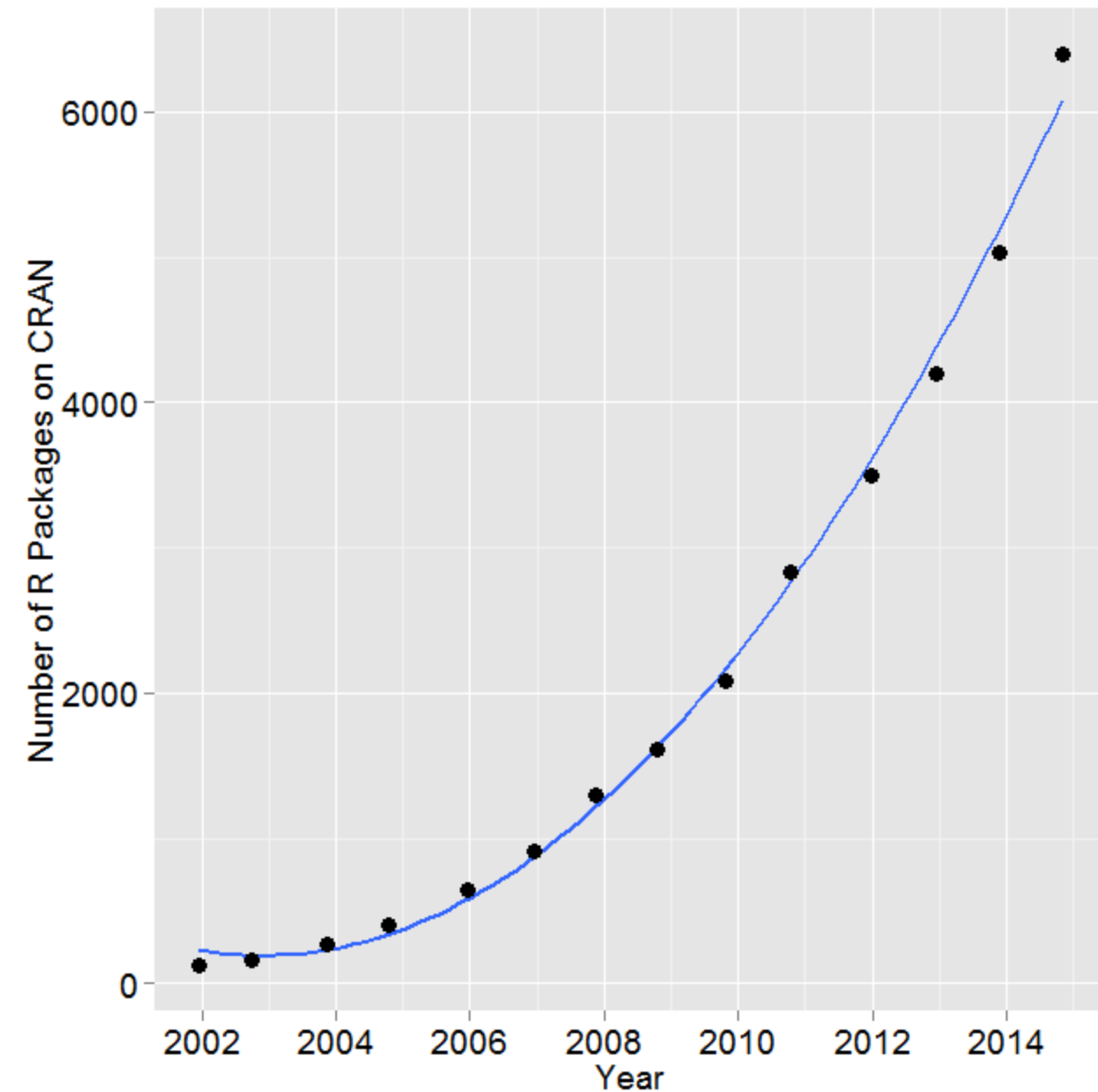


Programming tools: Adventures with R

A guide to the popular, free statistics and visualization software that gives scientists control of their own data analysis.

Sylvia Tippmann

29 December 2014 | Clarified: 13 February 2015



- Versatile
- Packages-Community
- Reproducibility
- Open and free
- Development

“During 2015 alone, R added more functions/procs than SAS Institute has written in its entire history.”

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FEATURE

60+ R resources to improve your data skills



By Sharon Machlis

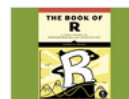
Executive Editor, Online & Data Analytics, Computerworld | OCT 3, 2016 4:08 AM PT



This list was originally published as part of the [Computerworld Beginner's Guide to R](#) but has since been expanded to also include resources for advanced beginner and intermediate users. If you're just starting out with R, I recommend first heading to the [Beginner's Guide PDF download](#).

These [websites](#), [videos](#), [blogs](#), [social media/communities](#), [software](#) and [books/ebooks](#) can help you do more with R; my favorites are listed in

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Book review -- The Book of R: A First Course in Programming and Statistics



Microsoft ratchets up its R enthusiasm



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6 new Jobs for R users – from all over the world (2017-05-24)

May 24, 2017

By Tal Galili

R jobs

To post your R job on the next post Just visit this link and post a new R job to the R community. You can post a job for free (and there are also “featured job” options available for extra exposure). Current R jobs Job

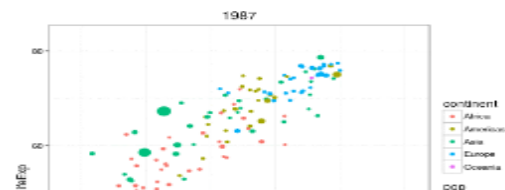
seekers: please follow the links below to learn more and apply for your R job of interest: Featured Jobs Full-Time Data Scientist...

[Read more »](#)

Create smooth animations in R with the tweenr package

May 23, 2017

By David Smith



Rcpp 0.12.11: Loads of goodies

May 23, 2017

By Thinking inside the box

The elevent update in the 0.12.* series of Rcpp landed on CRAN yesterday following the initial upload on the weekend, and the Debian package and Windows...

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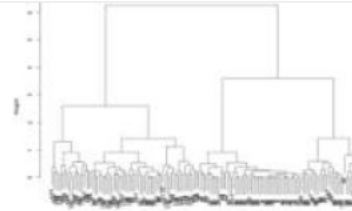
[Newest 'bioinformatics' Questions - Stack Overflow](#)

<https://stackoverflow.com/questions/tagged/bioinformatics>I was working in 'Genome of The Netherlands' GoNL where I served as a **Bioinformatics** researcher for a year. There were 2 Nature publications among many ...

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[R: K-Means Clustering](#)

<https://stat.ethz.ch/R-manual/R-devel/library/stats/html/kmeans.html> ▼

an **R** object of class "**kmeans**", typically the result of `ob <- kmeans(...)` method. character: may be abbreviated. "centers" causes fitted to return **cluster** centers ...

[Quick-R: Cluster Analysis](#)

www.statmethods.net/advstats/cluster.html ▼

R has an amazing variety of functions for **cluster** analysis. In this section, I will ... **K-means clustering** is the most popular partitioning method. It requires the ...

[How to Perform K-Means Clustering in R Statistical Computing ...](#)

<https://www.youtube.com/watch?v=sAtnX3UJyN0> ▼

Sep 29, 2013 - Uploaded by Influxity

In this video I go over how to perform **k-means clustering** using **r** statistical

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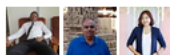
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K Means Clustering in R

December 28, 2015

By Teja Kodali

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(This article was first published on [DataScience+](#), and kindly contributed to R-bloggers)

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Hello everyone, hope you had a wonderful Christmas! In this post I will show you how to do k means clustering in R. We will use the `iris` dataset from the `datasets` library.

What is K Means Clustering?

K Means Clustering is an unsupervised learning algorithm that tries to cluster data based on their similarity. Unsupervised learning means that there is no outcome to be predicted, and the algorithm just tries to find patterns in the data. In k means clustering, we have to specify the number of clusters we want the data to be grouped into. The algorithm randomly assigns each observation to a cluster, and finds the centroid of each cluster. Then, the algorithm iterates through two steps:

- Reassign data points to the cluster whose centroid is closest.
- Calculate new centroid of each cluster.

These two steps are repeated till the within cluster variation cannot be reduced any further. The within cluster variation is calculated as the sum of the euclidean distance between the data points and their respective cluster centroids.

Exploring the data

The `iris` dataset contains data about sepal length, sepal width, petal length, and petal width of flowers of different species. Let us see what it looks like:

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New series: R and big data (concentrating on Spark and sparklyr)
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kmeans {stats}

R Documentation

K-Means Clustering

Description

Perform k-means clustering on a data matrix.

Usage

```
kmeans(x, centers, iter.max = 10, nstart = 1,  
       algorithm = c("Hartigan-Wong", "Lloyd", "Forgy",  
                     "MacQueen"), trace=FALSE)  
## S3 method for class 'kmeans'  
fitted(object, method = c("centers", "classes"), ...)
```

Arguments

- x**
numeric matrix of data, or an object that can be coerced to such a matrix (such as a numeric vector or a data frame with all numeric columns).
- centers**
either the number of clusters, say k , or a set of initial (distinct) cluster centres. If a number, a random set of (distinct) rows in **x** is chosen as the initial centres.
- iter.max**
the maximum number of iterations allowed.
- nstart**
if **centers** is a number, how many random sets should be chosen?
- algorithm**
character: may be abbreviated. Note that "Lloyd" and "Forgy" are alternative names for one algorithm.
- object**
an R object of class "kmeans", typically the result `ob` of `ob <- kmeans(...)`.
- method**
character: may be abbreviated. "centers" causes `fitted` to return cluster centers (one for each input point) and "classes" causes `fitted` to return a vector of class assignments.
- trace**
logical or integer number, currently only used in the default method ("Hartigan-Wong"): if positive (or true), tracing information on the progress of the algorithm is produced. Higher values may produce more tracing information.



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Download and Install R

Precompiled binary distributions of the base system and contributed packages, **Windows and Mac** users most likely want one of these versions of R:

- [Download R for Linux](#)
- [Download R for \(Mac\) OS X](#)
- [Download R for Windows](#)

R is part of many Linux distributions, you should check with your Linux package management system in addition to the link above.

Source Code for all Platforms

Windows and Mac users most likely want to download the precompiled binaries listed in the upper box, not the source code. The sources have to be compiled before you can use them. If you do not know what this means, you probably do not want to do it!

- The latest release (Friday 2017-04-21, You Stupid Darkness) [R-3.4.0.tar.gz](#), read [what's new](#) in the latest version.
- Sources of [R alpha and beta releases](#) (daily snapshots, created only in time periods before a planned release).
- Daily snapshots of current patched and development versions are [available here](#). Please read about [new features and bug fixes](#) before filing corresponding feature requests or bug reports.
- Source code of older versions of R is [available here](#).
- Contributed extension [packages](#)

Questions About R

- If you have questions about R like how to download and install the software, or what the license terms are, please read our [answers to frequently asked questions](#) before you send an email.

What are R and CRAN?

R is 'GNU S', a freely available language and environment for statistical computing and graphics which provides a wide variety of statistical and graphical techniques: linear and nonlinear modelling, statistical tests, time series analysis, classification, clustering, etc. Please consult the [R project homepage](#) for further information.

CRAN is a network of ftp and web servers around the world that store identical, up-to-date, versions of code and documentation for R. Please use the CRAN [mirror](#) nearest to you to minimize network load.

Submitting to CRAN

To "submit" a package to CRAN, check that your submission meets the [CRAN Repository Policy](#) and then use the [web form](#).

If this fails, upload to <ftp://CRAN.R-project.org/incoming/> and send an email to CRAN@R-project.org following the policy. Please do not attach submissions to emails, because this will

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Bring R to the web

“Open” software

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Dynamic Documents for R

R Markdown is an authoring format that enables easy creation of dynamic documents, presentations, and reports from R. It combines the core syntax of [markdown](#) (an easy to write plain text format) with embedded R code chunks that are run so their output can be included in the final document.

R Markdown documents are fully reproducible (they can be automatically regenerated whenever underlying R code or data changes).

R Markdown has many available output formats including [HTML](#), [PDF](#), [MS Word](#), [Beamer](#), [HTML5 slides](#), [Tufte handouts](#), [books](#), [dashboards](#), and [websites](#).

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Literate Computing

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by RStudio

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A web application framework for R

Turn your analyses into interactive web applications

No HTML, CSS, or JavaScript knowledge required

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Here is a Shiny app

Shiny apps are easy to write. No web development skills are required.

Number of bins in histogram (approximate):

20

ui.R

server.R

shinyUI(bootstrapPage(

Iris k-means clustering

X Variable

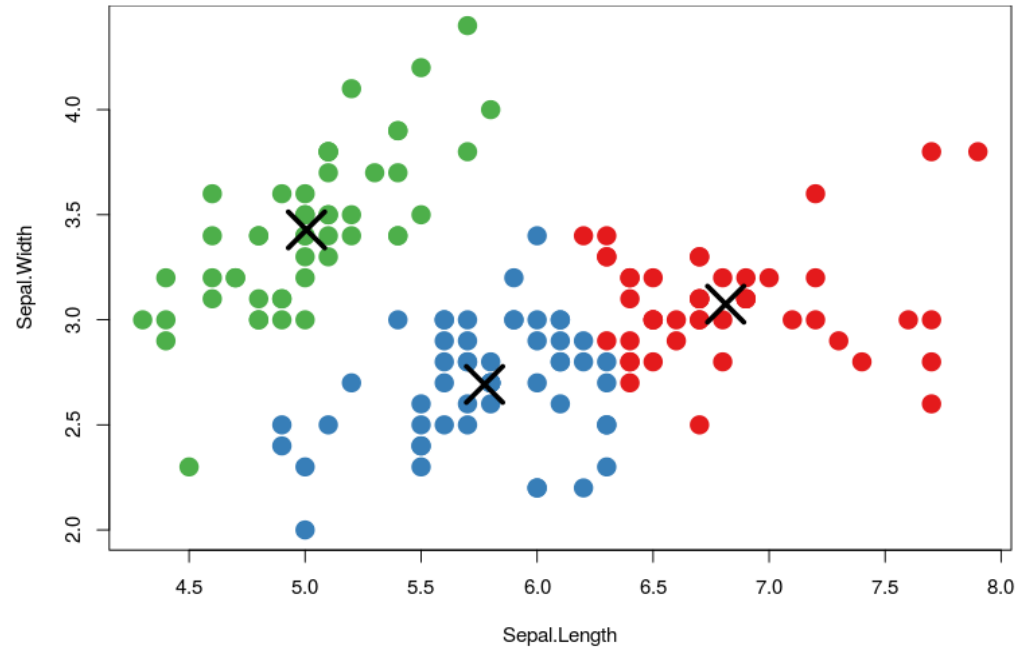
Sepal.Length

Y Variable

Sepal.Width

Cluster count

3



server.R

ui.R

show below

```
function(input, output, session) {  
  
  # Combine the selected variables into a new data frame  
  selectedData <- reactive({  
    iris[, c(input$xcol, input$ycol)]  
  })  
  
  clusters <- reactive({  
    kmeans(selectedData(), input$clusters)  
  })  
  
  output$plot1 <- renderPlot({  
    palette(c("#E41A1C", "#377EB8", "#4DAF4A", "#984EA3",  
              "#FF7F00", "#FFFF33", "#A65628", "#F781BF", "#999999"))  
  
    par(mar = c(5.1, 4.1, 0, 1))  
    plot(selectedData(),  
          col = clusters()$cluster,  
          pch = 20, cex = 3)  
    points(clusters()$centers, pch = 4, cex = 4, lwd = 4)  
  })  
}
```

Standardized whole-blood transcriptional profiling enables the deconvolution of complex induced immune responses

Interactive Application provided by
The Milieu Interieur Consortium

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Principal Component Analysis on the gene expression signature.

for more information, check the Instructions tab.

Select PCA Stimuli Space

[IFN-B, IFN-G, IL-1B, TNFa] based ▼

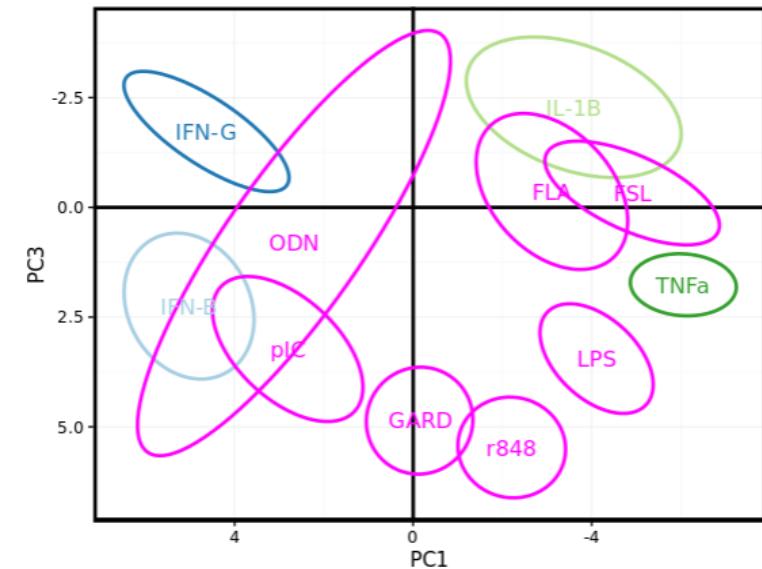
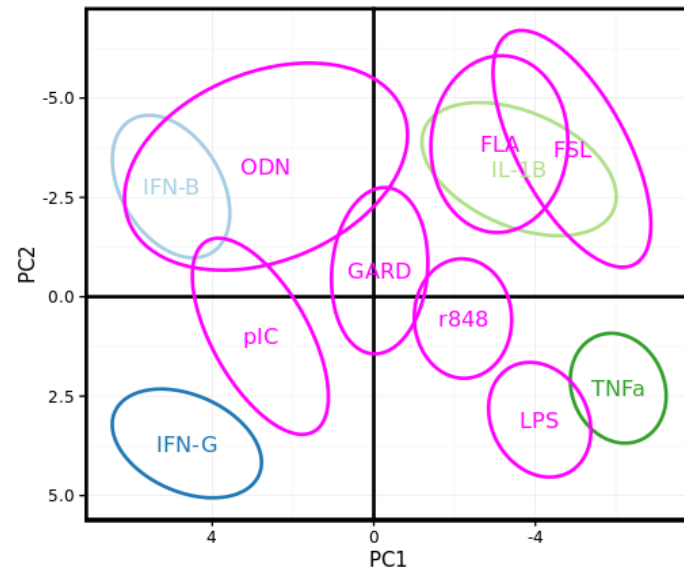
Fig.1A

Fig.1B

Fig.S4A

Fig.S4B

Fig.S5



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Bioconductor provides tools for the analysis and comprehension of high-throughput genomic data. Bioconductor uses the R statistical programming language, and is open source and open development. It has two releases each year, [1383 software packages](#), and an active user community. Bioconductor is also available as an [AMI](#) (Amazon Machine Image) and a series of [Docker](#) images.

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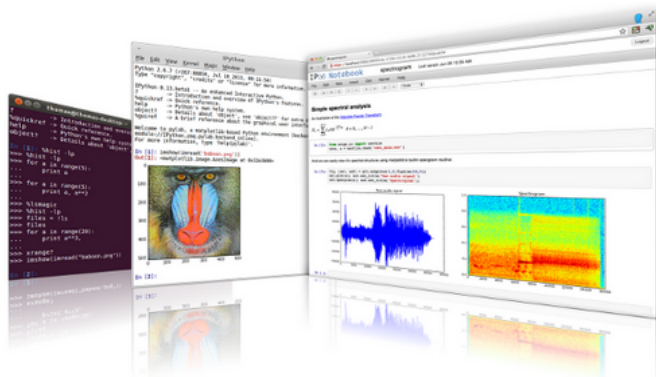


IP[y]: IPython Interactive Computing

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IPython provides a rich architecture for interactive computing with:

- A powerful interactive shell.
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To get started with IPython in the Jupyter Notebook, see our [official example collection](#). Our [notebook gallery](#) is an excellent way to see the many things you can do with IPython while learning about a variety of topics, from basic programming to advanced statistics or quantum mechanics.

To learn more about IPython, you can download our [talks and presentations](#), or read our [extensive documentation](#). IPython is open source (BSD license), and is used by a range of [other projects](#); add your project to that list if it uses IPython as a library, and please don't forget to [cite the project](#).

IPython supports Python 2.7 and 3.3 or newer. Our older 1.x series supports Python 2.6 and 3.2.

Jupyter and the future of IPython

IPython is a growing project, with increasingly language-agnostic components. IPython 3.x was the last monolithic release of IPython, containing the notebook server, qtconsole, etc. As of IPython 4.0, the

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June 20th - June 24th 2017, Berkeley, CA.

[CfP](#) closes on March 25th 31st!

Julia is part of **Google Summer of Code 2017!**

Check out [our project ideas](#) and get involved!

[Julia](#) is a high-level, high-performance dynamic programming language for numerical computing. It provides a sophisticated compiler, [distributed parallel execution](#), numerical accuracy, and an [extensive mathematical function library](#). Julia's Base library, largely written in Julia itself, also integrates mature, best-of-breed open source C and Fortran libraries for [linear algebra](#), [random number generation](#), [signal processing](#), and [string processing](#). In addition, the Julia developer community is contributing a number of [external packages](#) through Julia's built-in package manager at a rapid pace. [IJulia](#), a collaboration between the [Jupyter](#) and Julia communities, provides a powerful browser-based graphical notebook interface to Julia.

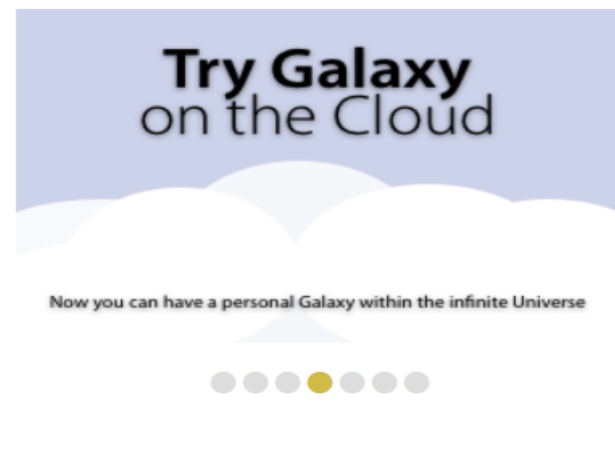
Les programmes Julia sont organisés autour de la [distribution multiple](#) (multiple dispatch) ; en définissant les fonctions et en les surchargeant pour différentes combinaisons de types d'arguments, qui peuvent également être définis par l'utilisateur. Pour une discussion plus approfondie de la justification et des avantages de Julia sur d'autres systèmes, voir les points suivants ou lire [l'introduction](#) dans le [manuel en ligne](#).

JuliaCon 2016, la conférence annuelle sur Julia, a eu lieu du 21 au 25 juin au MIT. Voici une vidéo aléatoire de notre liste de lecture YouTube de ces présentations. Cliquez sur l'icône de sélection pour voir les autres vidéos.

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[NGS: Gemini](#)
[NGS: Assembly](#)
[NGS: Chromosome Conformation](#)
[NGS: Mothur](#)
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[BEDTools](#)
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[EMBOSS](#)
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Galaxy is an open source, web-based platform for data intensive biomedical research. If you are new to Galaxy [start here](#) or consult our [help resources](#). You can install your own Galaxy by following the [tutorial](#) and choose from thousands of tools from the [Tool Shed](#).



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The [Galaxy Team](#) is a part of the [Center for Comparative Genomics and Bioinformatics](#) at [Penn State](#), the [Department of Biology](#) and at [Johns Hopkins University](#) and the [Computational Biology Program](#) at [Oregon Health & Science University](#).

This instance of Galaxy is utilizing infrastructure generously provided by the [CyVerse](#) at the [Texas Advanced Computing Center](#), with support from the [National Science Foundation](#).

Classification of pediatric acute lymphoblastic leukemia by gene expression profiling

Mary E. Ross, Xiaodong Zhou, Guangchun Song, Sheila A. Shurtleff, Kevin Girtman, W. Kent Williams, Hsi-Che Liu, Rami Mahfouz, Susana C. Raimondi, Noel Lenny, Anami Patel, James R. Downing

Blood 2003 102:2951-2959; doi:10.1182/blood-2003-01-0338

Example Data: Affymetrix chip 22282 ProbeIDs and 132 Samples

NEOPLASIA

Classification of pediatric acute lymphoblastic leukemia by gene expression profiling

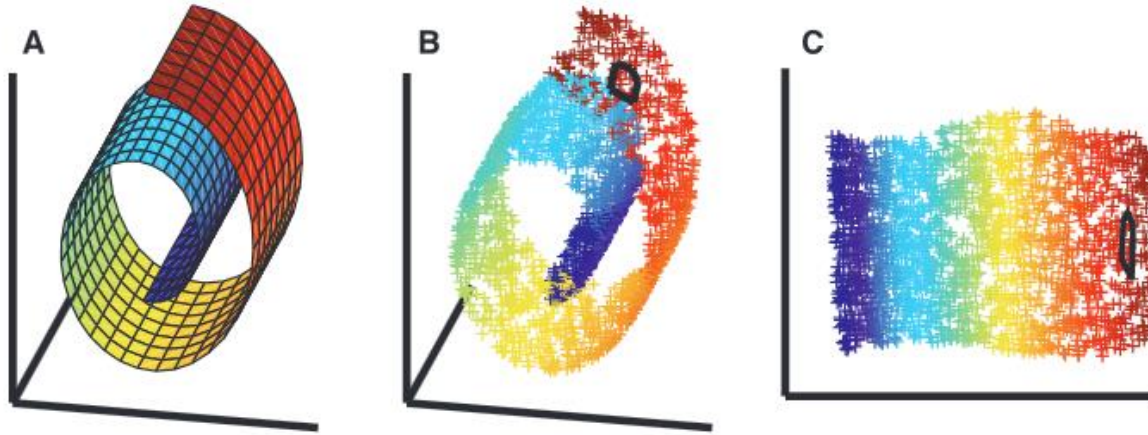
Mary E. Ross, Xiaodong Zhou, Guangchun Song, Sheila A. Shurtleff, Kevin Girtman, W. Kent Williams, Hsi-Che Liu, Rami Mahfouz, Susana C. Raimondi, Noel Lenny, Anami Patel, and James R. Downing

Contemporary treatment of pediatric acute lymphoblastic leukemia (ALL) requires the assignment of patients to specific risk groups. We have recently demonstrated that expression profiling of leukemic blasts can accurately identify the known prognostic subtypes of ALL, including T-cell lineage ALL (T-ALL), *E2A-PBX1*, *TEL-AML1*, *MLL* rearrangements, *BCR-ABL*, and hyperdiploid karyotypes with more than 50 chromosomes. As the next step toward developing this methodology into a frontline diagnostic tool, we have now

analyzed leukemic blasts from 132 diagnostic samples using higher density oligonucleotide arrays that allow the interrogation of most of the identified genes in the human genome. Nearly 60% of the newly identified subtype discriminating genes are novel markers not identified in our previous study, and thus should provide new insights into the altered biology underlying these leukemias. Moreover, a proportion of the newly selected genes are highly ranked as class discriminators, and when incorporated into class-predict-

ing algorithms resulted in an overall diagnostic accuracy of 97%. The performance of an array containing the identified discriminating genes should now be assessed in frontline clinical trials in order to determine the accuracy, practicality, and cost effectiveness of this methodology in the clinical setting. (Blood. 2003; 102:2951-2959)

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LLE ALGORITHM

1. Compute the neighbors of each data point, $\vec{X}_i$.
2. Compute the weights W_{ij} that best reconstruct each data point $\vec{X}_i$ from its neighbors, minimizing the cost in eq. (1) by constrained linear fits.
3. Compute the vectors $\vec{Y}_i$ best reconstructed by the weights W_{ij} , minimizing the quadratic form in eq. (2) by its bottom nonzero eigenvectors.

Locally Linear Embedding LLE

S. T. Roweis and L. K. Saul,
Nonlinear Dimensionality Reduction by Locally Linear Embedding,
 Science Vol 290, 2000

R package lle

Package 'lle'

February 20, 2015

Type Package

Title Locally linear embedding

Version 1.1

Date 2012-03-21

Author Holger Diedrich, Dr. Markus Abel (Department of Physics,
 University Potsdam)

Maintainer Holger Diedrich <holgerdiedrich@gmx.net>

Stochastic Neighbor Embedding

Original distances

$$p_{j|i} = \frac{\exp(-\|x_i - x_j\|^2 / 2\sigma_i^2)}{\sum_{k \neq i} \exp(-\|x_i - x_k\|^2 / 2\sigma_i^2)},$$

$$q_{j|i} = \frac{\exp(-\|y_i - y_j\|^2)}{\sum_{k \neq i} \exp(-\|y_i - y_k\|^2)}.$$

Distances in reduced space

$$q_{ij} = \frac{(1 + \|y_i - y_j\|^2)^{-1}}{\sum_{k \neq l} (1 + \|y_k - y_l\|^2)^{-1}}.$$

The cost function C is given by

$$C = \sum_i KL(P_i || Q_i) = \sum_i \sum_j p_{j|i} \log \frac{p_{j|i}}{q_{j|i}},$$

Original SNE

tSNE

Kullback Leibler
Relative Entropy

Package 'Rtsne'

June 30, 2016

Type Package

Title T-Distributed Stochastic Neighbor Embedding using a Barnes-Hut Implementation

Version 0.11

Description An R wrapper around the fast T-distributed Stochastic Neighbor Embedding implementation by Van der Maaten.

License BSD_3_clause + file LICENSE

URL <https://github.com/jkrijthe/Rtsne>

Imports Rcpp (>= 0.11.0)

Journal of Machine Learning Research 1 (2008) 1-48

Submitted 4/00; Publ

Visualizing Data using t-SNE

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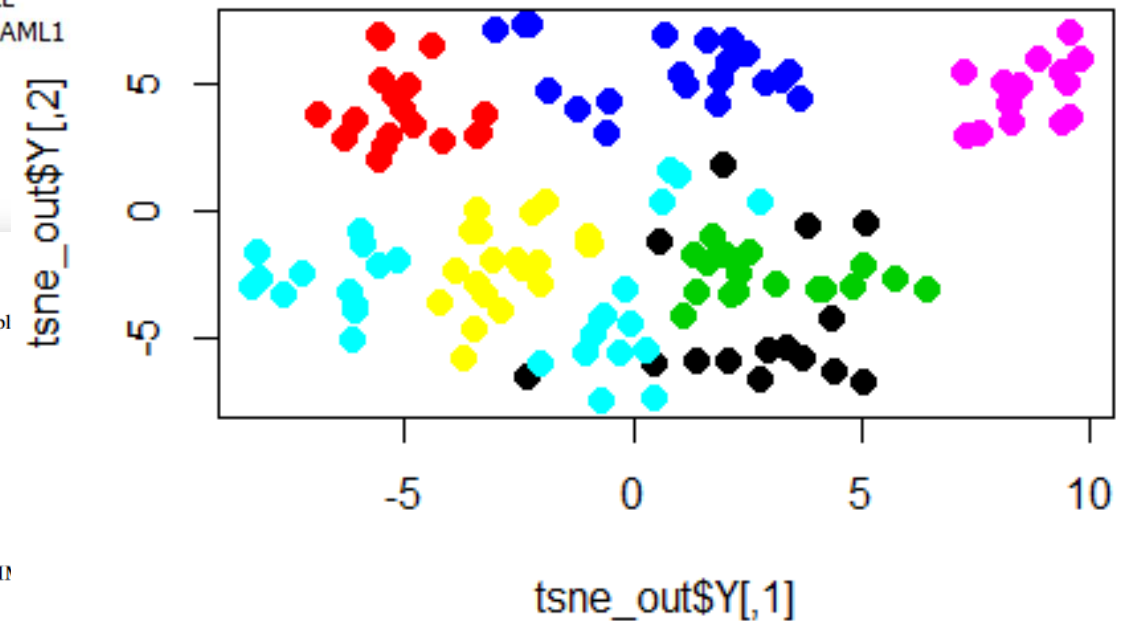
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Stochastic Neighbor Embedding (SNE) and tSNE

BCR-ABL
E2A-PBX1
Hyperdiploid (>50)
MLL
Other
T-ALL
TEL-AML1



Copenhagen 4th-8th of December 2017

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