

Python for Scientists

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Me

- Researcher at the CSIC
- Computational biochemistry and biophysics
- Started with Fortran. Learnt some Perl, R, Mathematica.
 - But now mainly use Python
- Analysis of simulations
- Developing new methods

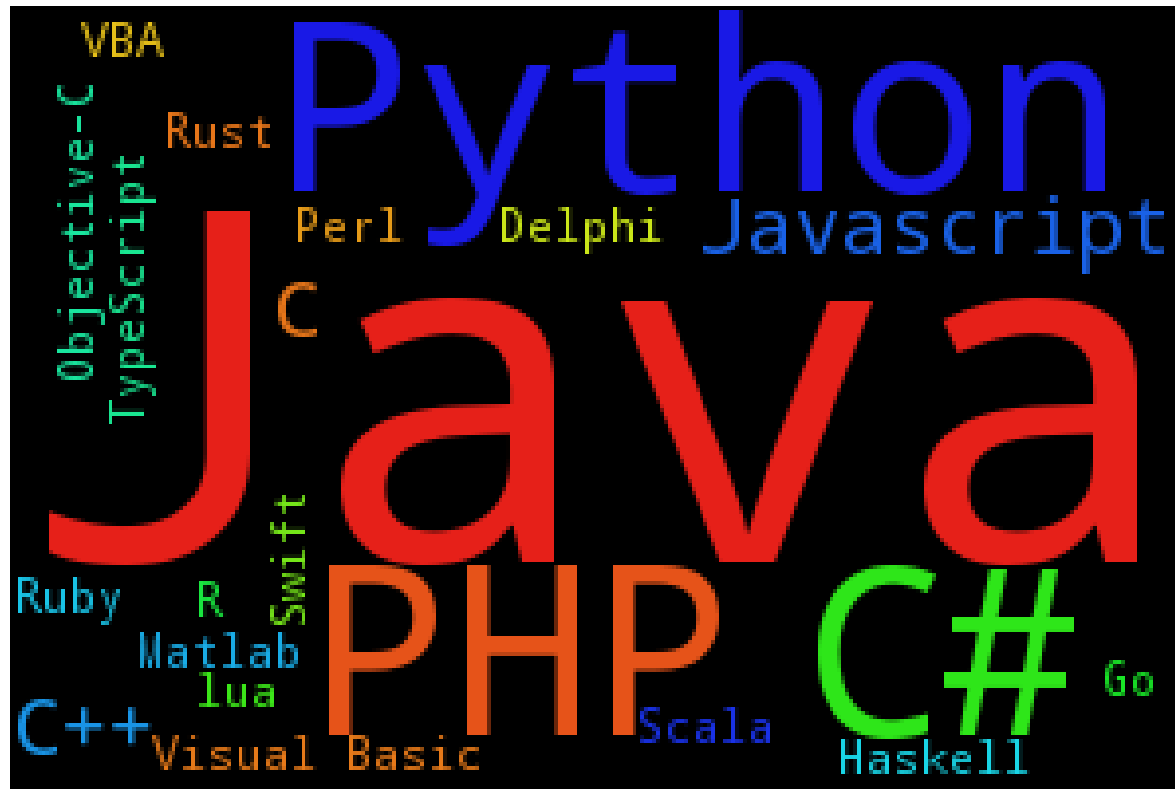
You?

- Research institute?
- Research field?
- Known languages?
- Kind of data?
- Expectations?

Overview

- Why Python
- Language basics
- Working with files
- Working with arrays: Numpy
- Data visualization: matplotlib
- Functions and modules
- Scientific modules. Scipy
- Classes and objects (bare minimum!)
- Other scientific modules: scikit-learn, biopython...
- Profiling and optimization and beyond Python

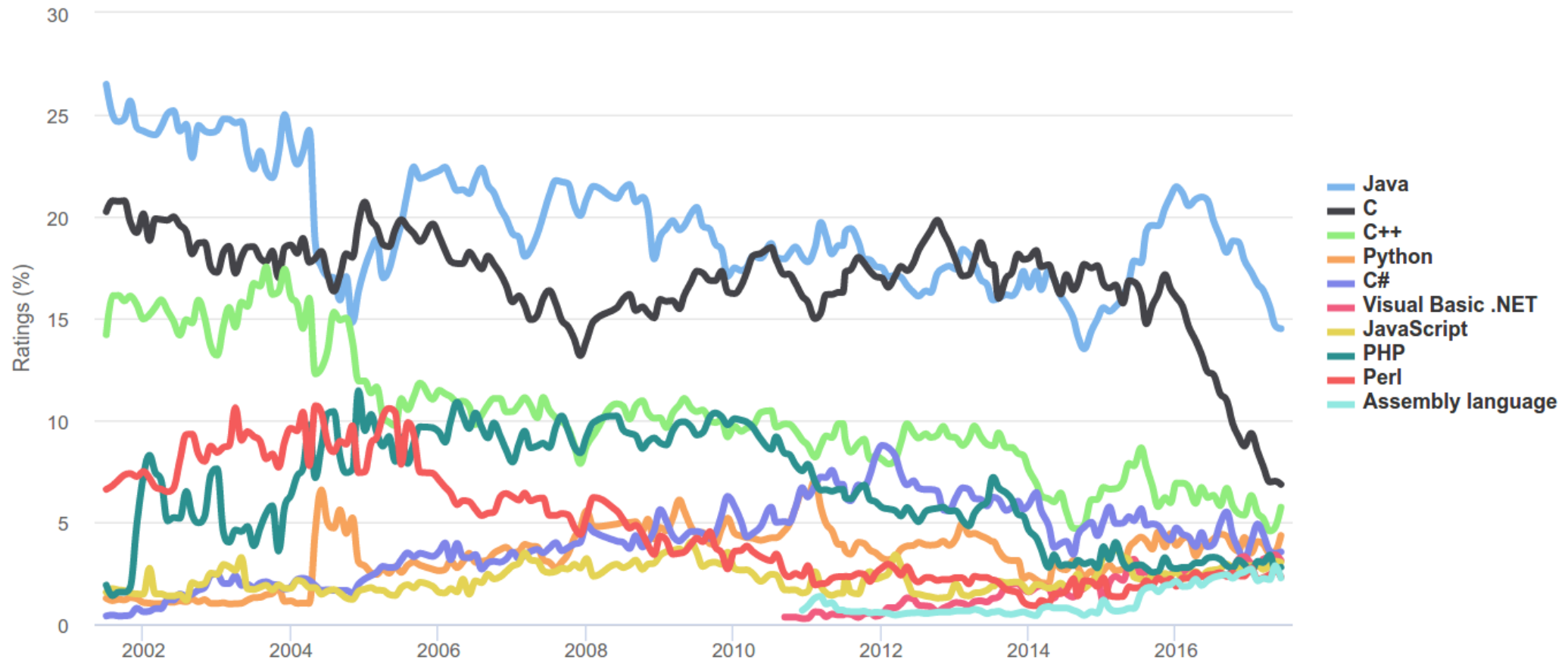
Language popularity



<http://pypl.github.io/PYPL.html>

Language popularity

Source: www.tiobe.com



<http://www.tiobe.com/index.php/content/paperinfo/tpci/index.html>

Hammerprinciple.com

PYTHON



<http://hmrp.pl/zuT1C8>

Based on 141739 responses from 12230 people, this is the picture we've built up of Python.

DOES WELL AT...

- ↑ I enjoy using this language
- ↑ I find this language easy to prototype in
- ↑ I would use this language for casual scripting
- ↑ I know this language well
- ↑ I regularly use this language
- ↑ When I write code in this language I can be very sure it is correct
- ↑ I rarely have difficulty abstracting patterns I find in my code
- ↑ I usually use this language on solo projects
- ↑ I would use this language for a web project
- ↑ I would use this language for a desktop GUI project

DOES POORLY AT...

- ↓ I often get angry when writing code in this language
- ↓ I often feel like I am not smart enough to write this language
- ↓ This language has an annoying syntax
- ↓ I learned this language early in my career as a programmer
- ↓ This language makes it easy to shoot yourself in the foot
- ↓ Developers who primarily use this language often burn out after a few years
- ↓ This language is unusually bad for beginners
- ↓ I am reluctant to admit to knowing this language
- ↓ The thought that I may still be using this language in twenty years time fills me with dread
- ↓ Writing code in this language is a lot of work

Python for science

- A high level language gives more time to more complex problems
 - At the expense of hiding important details
- Example:
 - A reaction mechanism
 - Optimisation of an energy function
 - Steepest descent, conjugate gradients, quasi-Newton
 - Implementation of BFGS quasi-Newton
 - Memory issues, diagonalization, matrix inversion...
 - Calculation of numerical gradients or hessians:
 - machine precision, central differences, etc.

http://fperez.org/py4science/why_python.html

Python for science

“We then generated 1000 random sequences with randomly specified $\langle H \rangle$ and $\langle Q \rangle$ values, and conducted molecular-dynamics simulations to calculate the $\langle R_g \rangle$ for each chain. The obtained $\langle R_g \rangle$ values were combined with the two-state formalism to determine whether the chain was ordered (globule) or disordered (coil)”

Biophysical Journal, **104**, 2013, 488–495



OS calls

Numerical analysis

File parsing

Data visualization

Python for science

Compiled languages

Fast
Difficult
non-interactive

Matlab, Mathematica, Octave

Slow
Rich libraries
Nice development environment
Restricted base language
Expensive (some)

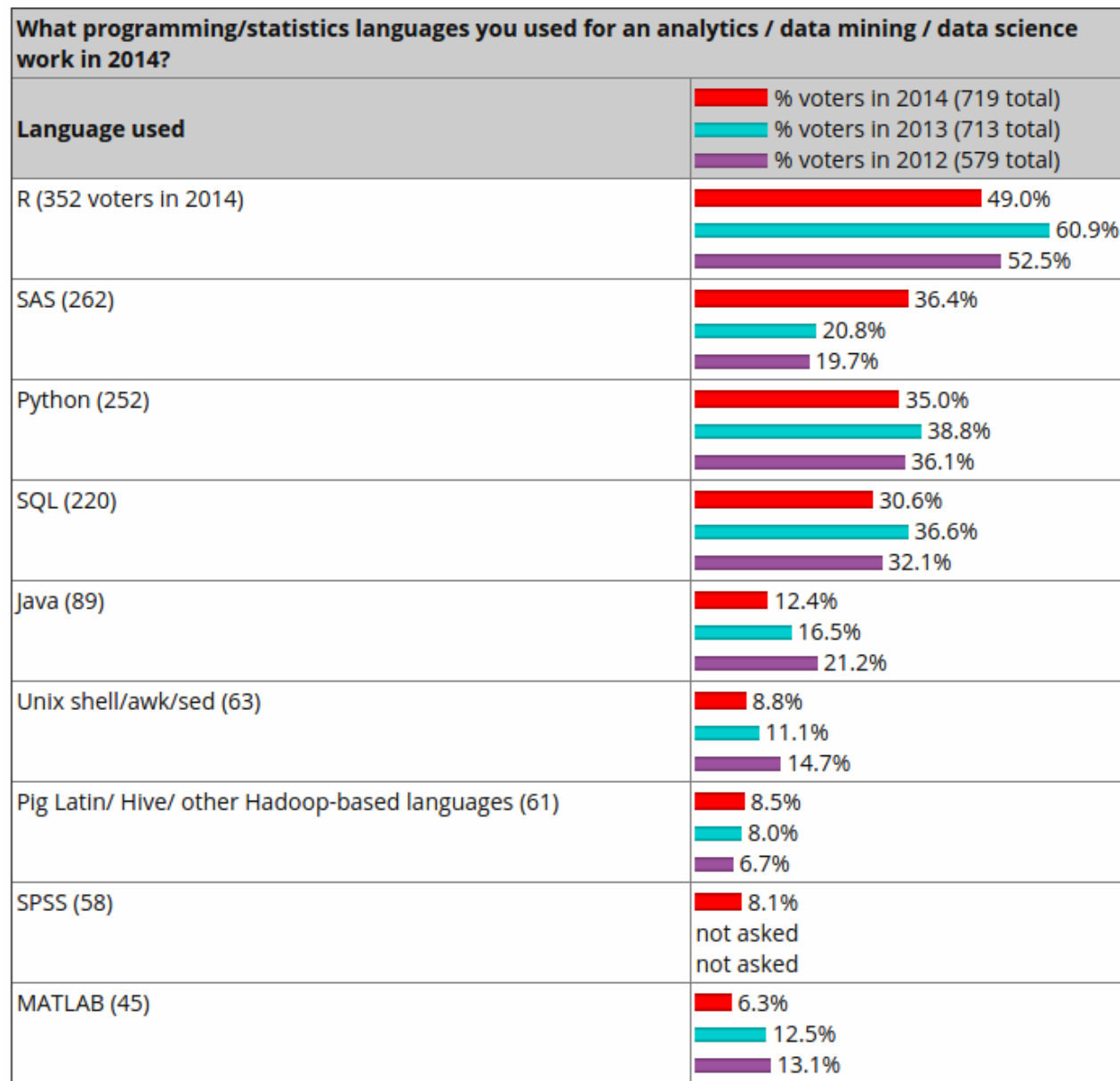
Python

Rich libraries (less than matlab)
Other libraries
Free
Active community
Harder than Matlab

Matlab, Mathematica?

- Scientific computing:
 - ipython + scipy + matplotlib
- Free
- Open source
- Extensible
- Bioinformatics
 - Biopython
- Molecular Dynamics
 - MMTK
- Efficiency
 - Numba, Cython, Fortran, C
- Server control
- XML parser

Python for data science



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THIS LANGUAGE ENCOURAGES WRITING CODE THAT IS EASY TO MAINTAIN.



THIS IS A MAINSTREAM LANGUAGE



THIS LANGUAGE IS GOOD FOR BEGINNERS



I WOULD USE THIS LANGUAGE AS A SCRIPTING LANGUAGE EMBEDDED INSIDE A LARGER APPLICATION



I WOULD USE THIS LANGUAGE FOR WRITING SERVER PROGRAMS



I WOULD USE THIS LANGUAGE FOR MOBILE APPLICATIONS



I CAN IMAGINE THIS WILL BE A POPULAR LANGUAGE IN TWENTY YEARS TIME



<http://hmrp.pl/x79RTk#73>

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DEVELOPERS WHO PRIMARILY USE THIS LANGUAGE OFTEN BURN OUT AFTER A FEW YEARS



THE SEMANTICS OF THIS LANGUAGE ARE MUCH DIFFERENT THAN OTHER LANGUAGES I KNOW.



WRITING CODE IN THIS LANGUAGE IS A LOT OF WORK



THIS LANGUAGE HAS A NICHE IN WHICH IT IS GREAT



THIS LANGUAGE HAS AN ANNOYING SYNTAX



THIS LANGUAGE IS UNUSUALLY BAD FOR BEGINNERS



I OFTEN FEEL LIKE I AM NOT SMART ENOUGH TO WRITE THIS LANGUAGE



THIS LANGUAGE HAS A NICHE OUTSIDE OF WHICH I WOULD NOT USE IT



<http://hmrp.pl/x79RTk#27>

Python for data science

- Which is better for data analysis: R or Python?
<http://www.quora.com/Which-is-better-for-data-analysis-R-or-Python>
- SAS vs. R (vs. Python) – which tool should I learn?
<http://www.analyticsvidhya.com/blog/2014/03/sas-vs-vs-python-tool-learn/>
- Python Vs R Machine learning
<http://datascience.stackexchange.com/questions/326/python-vs-r-machine-learning>
- How to Choose Between Learning Python or R First
<http://blog.udacity.com/2015/01/python-vs-r-learn-first.html>
- Python, Machine Learning, and Language Wars
<http://sebastianraschka.com/blog/2015/why-python.html>

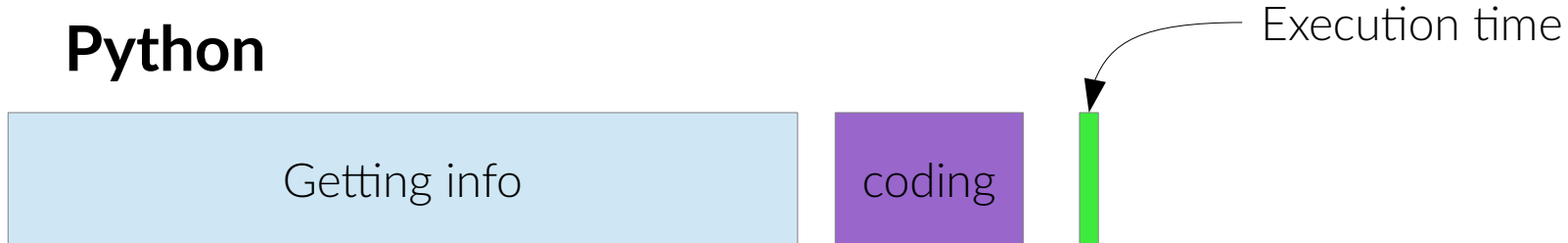
Low level vs. high level

- Python is a high level language
- You can focus on:
 - Low level issues
 - Higher complexity of problems
- Low level issues
 - Variable types
 - Machine precision
- But also
 - Extend
 - Maintain
 - Document code

Python vs. Fortran/C

Different time distribution to get a task done

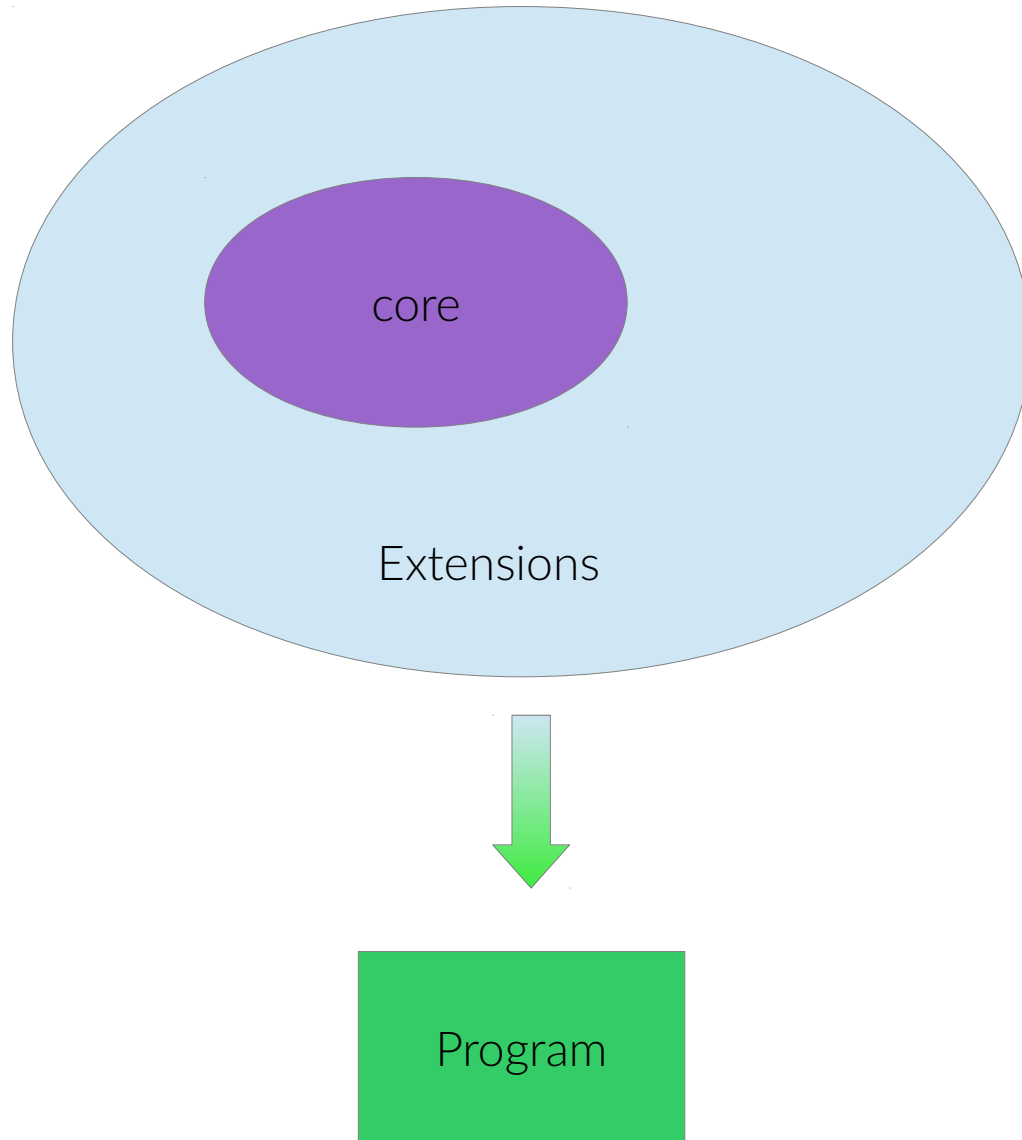
Python



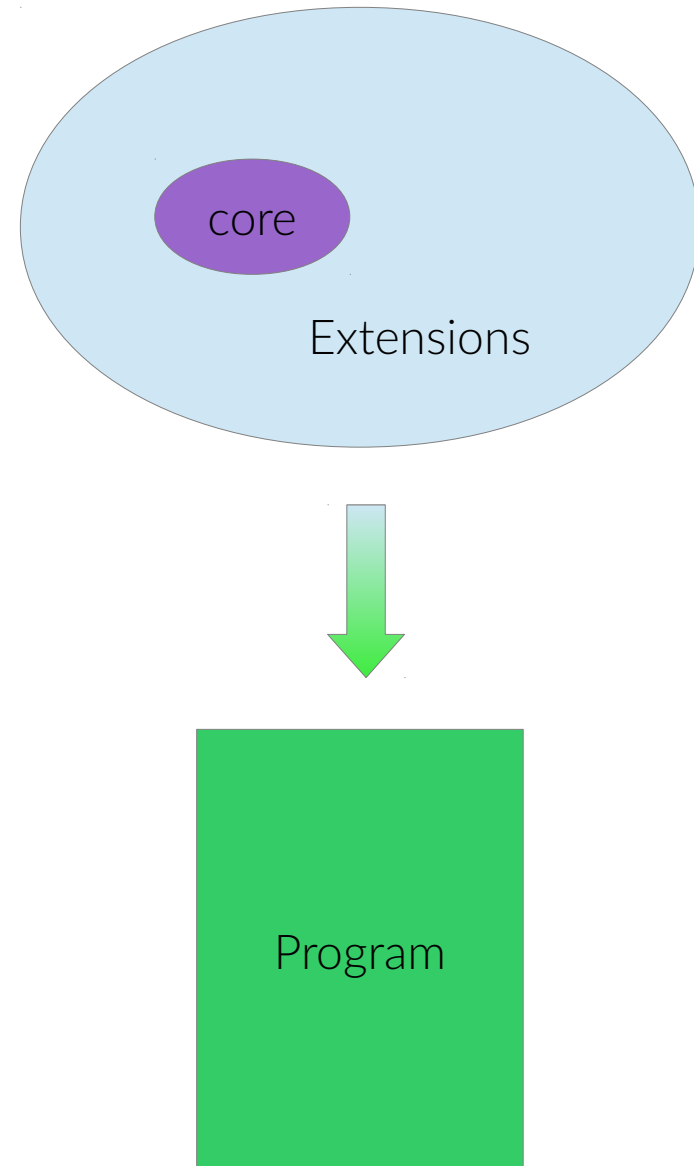
Fortran / C



Python



Fortran/C



Python for science

- The homogenization of scientific computing, or why Python is steadily eating other languages' lunch

<http://www.talyarkoni.org/blog/2013/11/18/the-homogenization-of-scientific-computing-or-why-python-is-steadily-eating-other-languages-lunch/>

- 10 Reasons Python Rocks for Research (And a Few Reasons it Doesn't)

<http://www.stat.washington.edu/~hoytak/blog/whypython.html>



Hello World program

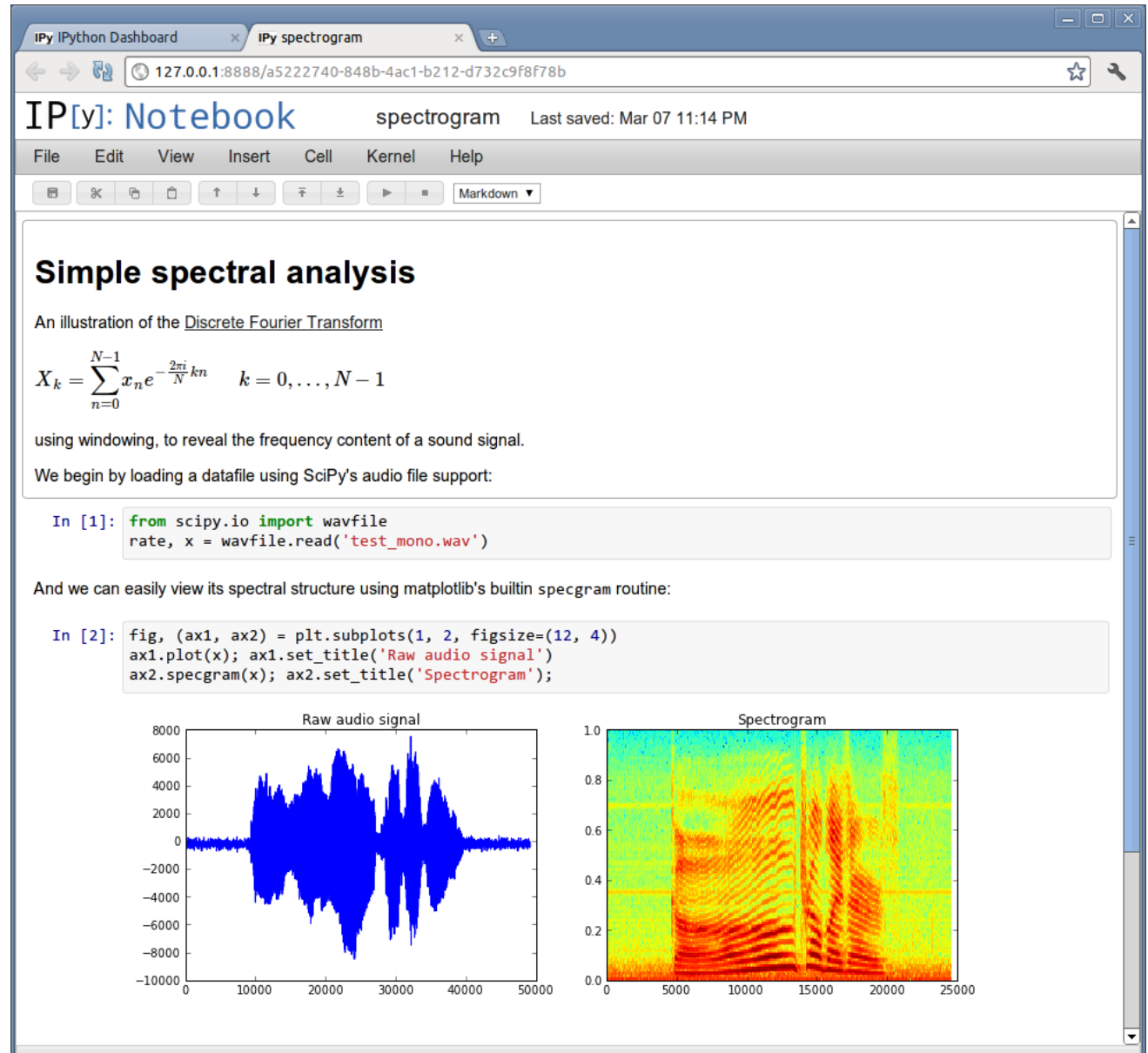
```
print("Hello World!")
```

```
print("Hello World!")
```

```
$ python3 hello.py
```

Interactive shells

- python
- IDLE
- Jupyter
(previously called ipython)
 - shell
 - notebook
- spyder
- eric
- PIDA
- Sage



Python distributions

- Anaconda
 - <https://www.continuum.io/downloads>
- Enthought Canopy
 - <https://www.enthought.com/products/canopy/>

Dynamically typed

```
>>> a = 4
>>> type(a)
<class 'int'>
>>> b = 7.6
>>> type(b)
<class 'float'>
>>> type(a+b)
<class 'float'>
>>> c = 'Hola'
>>> c + ' Que tal?'
'Hola Que tal?'
>>> c + a
Traceback (most recent call last):
  File "<stdin>", line 1, in <module>
TypeError: Can't convert 'int' object to str implicitly
```

Which python version?

- Language is fast evolving
- 2 versions now coexist: 3.x and 2.x
- These versions are not completely compatible
- 3.x is better and continued
- 2.x has some software still not ported
- Both can safely coexist
 - Packages and shells are for a specific version
- **2to3 -w hello.py**

Language elements

Numbers

Integers:

```
> i = 5  
> j = i**i**i
```

Limited by amount of memory:

```
>>> i.bit_length()  
3  
  
>>> j = i**i**i  
>>> j.bit_length()  
7257  
  
>>> 9 % 5 #modulo  
4
```

Floating point:

```
>>> x = 5.  
>>> y = x**x**x  
Traceback (most recent call last):  
  File "<stdin>", line 1, in <module>  
OverflowError: (34, 'Numerical result  
out of range')
```

Division vs integer division (Python 3):

```
>>> 3/2  
1.5  
  
>>> 3//2  
1  
  
>>> j/i #Returns a Float
```

Assignments

Explicit notation:

```
> i = i+1  
> j = j / 10.
```

Short notation:

```
> i += 1  
> j /= 10.
```

Floating point:

```
>>> x = 5.  
>>> y = x**x**x
```

Traceback (most recent call
last):

File "<stdin>", line 1, in
<module>

OverflowError: (34, 'Numerical
result out of range')

strings

Strings:

```
> str(6.7)
> c = 'Hola'
```

Operations:

```
> s = 'numeric ' + 'python'
> len(s)
> s[5]
'i'
> s.split()
['numeric', 'python']
```

```
> print('Result: %5.3f' % (11./3.))
3.667
```

Non mutable:

```
> s[6]
> s[6]='7' #Error!
```

Regular expressions

```
import re
```

Lists, sets and tuples

- Heterogeneous containers:

```
> l=[6, 'a', [5,[9,8,7,6]],  
    -6.5, (True, True)]  
> [1,2]+[3,4]  
> l.append(6)
```

- sets:


```
s=set([4,3,2,3])  
> 4 in s  
True  
> s  
set([2, 3, 4])
```

- Tuples are immutable lists
t=(1,2,3)

Lists, sets and tuples

List indexing and methods:

```
> l = list(range(10))
> l[4] = 20
> l[4:]
> l[-4]
> l[:]
> l[::-1] #reverse
> l.reverse()
> l.pop()
> l.extend([3,4,5])
> l.sort()
```



First index
is 0

Set methods:

```
> s1=set([1,2,3,4])
> s2=set([3,4,5,6,7])
> s1.union(s2)
> s1.intersection(s2)
> s1.difference(s2)
> s2.difference(s1)
> s1.intersection(s2) == s2 & s1
True
> s1 - s2 == s2 - s1
False
```

Uses of lists, sets and tuples

- Calculate and keep all the primes < 1000
- Given a coordinate file, calculate for each atom a list of all the atoms that are at less than 0.2nm.
- Get the solutions of a quadratic equation (0,1,2) or (real vs. complex).
- <http://docs.python.org/3/tutorial/datastructures.html>

Copying and looping over lists

lists are treated as pointers:

copying lists, makes a copy of the pointer.

```
> l=[1,2,3,4]
```

```
> l2=l
```

```
> l[2]=1000
```

```
l1
```

```
[1, 2 , 1000, 4]
```

Looping over lists:

Fortran/C style:

```
num=[2,3,2,3,4,5,5]
```

```
for i in range(len(num)):
```

```
    print(num[i])
```

Pythonic style:

```
for item in num:
```

```
    print(item)
```

This can be used for sets, dictionaries, and tuples.

Dictionaries

Setting elements:

```
> phone={}
> phone['Ramon']='1242'
> phone['Joan']='1323'
> phone['Quique']='1242'
> phone.keys()
['Quique', 'Joan', 'Ramon']
> d2 = dict(Ramon=1242, Joan=1323,
    Quique=1242)
```

Dictionaries are not ordered

Getting elements:

```
> for key in phone:
...     print(key, phone[key])
Quique 1242
Joan 1323
Ramon 1242
```

Removing elements:

```
> del(phone['Ramon'])
```

The beauty of Python blocks

We are usually told to indent blocks for clarity.

Python makes this the syntax rule to identify blocks.

The code has to be nice!

Convention:

- Use 4 spaces
- Use spaces, not tabs.

```
while iter < maxIter:
    x = f(x)
    iter = iter + 1

if i>0:
    print("i is positive")
elif i==0:
    print("i is zero")
else:
    print("i is negative")
```

Execution control: if

if... elif... else

if <condition>:

<block>

elif <condition>:

<block>

else:

<block>

4==4 #True

5!=4 #True

4>=5 #False

4 in [4,5] #True

result=True

if result: print('yes')

Execution control

Conditions can be combined with:

`and or not ()`

Object identity:

```
> a=[1,2,3]
```

```
> b=a
```

```
> b is a
```

```
True
```

Any non-zero number or non-empty string is True:

```
> if []: print ('yes')  
      else: print('no')
```

```
no
```

```
> if 5 and 'result':  
      print('yes')
```

```
else:
```

```
      print('no')
```

```
yes
```

```
> if 5 or 1/0: print('yes')
```

```
yes
```

for and while loops

For loops

```
> dict={4:'a',3:'b', 2:'c',  
1:'d'}  
> for i in dict:  
    print(i, dict[i])
```

While

```
while <condition>:  
    <block>
```

Break continue pass

```
> pass # does nothing
```

break: Exit loop

```
if x>0:  
    pass  
else:  
    break
```

cycle: Continue with the next iteration

list comprehensions

simple way to create lists:

```
> l=[x**2 for x in range(8)]  
[0, 1, 4, 9, 16, 25, 36, 49]
```

with conditionals:

```
l2= [(i, -2*i+3) for i in l if i % 3 == 0]  
[(0, 3), (9, -15), (36, -69)]
```

Nested lists:

```
> [(x, y) for x in [1,2,3] for y in [3,1,4] if x != y]  
[(1, 3), (1, 4), (2, 3), (2, 1), (2, 4), (3, 1), (3, 4)]
```

list comprehension and enumerate

Enumerate indexes lists:

```
line='how do you do?'  
line=line.split()  
for i, word in enumerate(line):  
    print(i, word.upper())  
0 HOW  
1 DO  
2 YOU  
3 DO?
```

Enumerate returns an iterator

```
> enumerate(['a', 'b', 'c'])  
<enumerate object at 0x1ebeaa50>
```

Be pythonic

Convert the negative elements of a list to positive

```
>>> x = [1, 2, -4, -5, 3, -5]
```

```
j = 0
```

```
while j < len(x):
```

```
    x[j] = abs(x[j])
```

```
    j += 1
```

```
for j in range(len(x)):
```

```
    x[j] = abs(x[j])
```

Or with list comprehensions

```
x = [abs(j) for j in x]
```

Or with functional programming

```
x= map(abs, x) #returns an iterator
```

<http://docs.python-guide.org/en/latest/writing/style/>

More python functions

```
print(3,4,5, sep='o', end='<<<<\n')
```

```
zip([1,2,3], ['a', 'b', 'c', 'd'])
```

```
a = input('Write a number: ')
```

```
len([1,2,3])
```

```
list(range(5))
```

```
range(20,10,-1)
```

```
sorted([5,4,3,5])
```

```
sum([5,4,3,5])
```

<http://docs.python.org/3.3/library/functions.html>

Mutable and immutable

- Mutable objects can be mutated.
 - Their identity remains the same
- Immutable objects are “mutated” by creating a new object

```
>>> a = 4
>>> id(a)
9157088
>>> a += 2
>>> id(a)
9157152
>>> s = 'Hola'
>>> id(s)
140165884365656
>>> s = s + ' que tal?'
>>> id(s)
140165884365712
>>> ll = [3,4,5]
>>> id(ll)
140165884674416
>>> ll.append(6)
>>> id(ll)
140165884674416
```

Identity and equality

```
>>> 1.0 is 1.0
```

```
True
```

```
>>> 1.0 == 1.0
```

```
True
```

```
>>> 1 == 1.0
```

```
True
```

```
>>> 1 is 1.0
```

```
False
```

```
>>> a = 4
```

```
>>> b = a
```

```
>>> a is b
```

```
True
```

```
>>> id(a)
```

```
9157088
```

```
>>> id(b)
```

```
9157088
```

```
>>> l1 = [1,2,3,]
```

```
>>> l2 = l1
```

```
>>> l2 = l1[:]
```

```
>>> l2 is l1
```

```
False
```

```
>>> l2 == l1
```

```
True
```

Objects: everything

```
>>> a = 5
>>> isinstance(a, int)
True
>>> object
<class 'object'>
>>> int
<class 'int'>
>>> isinstance(a, object)
True
>>> issubclass(int, object)
True
```

Objects have variables:

```
> c = 4+5j
> c.real
```

Objects have methods:

```
> c.conjugate #the method
> c.conjugate() #its call
```

And we can apply functions to objects:

```
> abs(c)
```

Python flow with pythontutor

The screenshot displays the Python Tutor interface with the following components:

- Code Editor:** Contains Python code for list manipulation and function calls.
- Frames:** A table showing the current execution frame (Global frame) and its variables.
- Objects:** A table showing the objects created in memory, including lists and functions.
- Execution Flow:** A progress bar and navigation buttons indicating the current line of execution.
- Program Output:** A box showing the output of the program.

Code:

```
1 x = [1, 2, 3]
2 y = [4, 5, 6]
3 z = y
4 y = x
5 x = z
6
7 x = [1, 2, 3] # a different [1, 2, 3] list!
8 y = x
9 x.append(4)
10 y.append(5)
11 z = [1, 2, 3, 4, 5] # a different list!
12 x.append(6)
13 y.append(7)
14 y = "hello"
15
16
17 def foo(lst):
18     lst.append("hello")
19
```

Frames:

Variable	Value
x	[1, 2, 3]
y	"hello"
z	[1, 2, 3, 4, 5]
foo	function foo(lst)
bar	function bar(myLst)

Objects:

Object	Value
list	[1, 2, 3, 4, 5, 6, 7, "hello"]
list	[1, 2, 3, 4, 5, "hello"]

Program output:

```
[1, 2, 3, 4, 5, 6, 7, 'hello']
[1, 2, 3, 4, 5, 'hello']
```

try... except

“Look before you leap”:

```
def safe_divide_1(x, y):  
    if y==0:  
        print("Divide-by-0 attempt  
detected")  
        return None  
    else:  
        return x/y
```

“It's easier to ask forgiveness than permission”:

```
def safe_divide_2(x, y):  
    try:  
        return x/y  
    except ZeroDivisionError:  
        print("Divide-by-0 attempt  
detected")  
        return None
```

Short ipython tutorial

beyond python

TAB autocomplete:

- functions
- methods
- files
- ...

`reload` command

cursor keys get history (for console only):

- even previous sessions!
- text + keys: previous match

`?:` help

`%quickref`

`Ctrl-r` : previous commands

Without ipython:

`python3 -u script.py` enters interactive mode

Magic functions

`%timeit x=10` : time the 'x=10' statement with high precision.

`%%timeit x=2**100`

`x*100` : time 'x*100' with a setup of 'x=2**100'; setup code is not counted. This is an example of a cell magic.

`%cpaste, %paste`: Paste & execute a pre-formatted code block from clipboard.

`%history`

`%load_ext`

`%run`

`%pdb`: Control the automatic calling of the pdb interactive debugger.

`%pylab`

`%timeit`

`%pwd`

`%cd`

`%%bash` <http://ipython.org/ipython-doc/dev/interactive/tutorial.html>

running scripts

`%run script.py`

`import script.py`

are not the same!

`%run script.py` is like `python3 script.py`

Imports are only “impoted” once in a session (see later `%autoreload` magic function)

jupyter notebook

- Nice presentation
- Allows parallel execution
- Combines text and code
- Executable or exportable to:
 - html
 - LaTeX
 - python
- Start with: `ipython3 notebook`
- Examples:

<https://github.com/jrjohansson/scientific-python-lectures>

Files

Files

- Files can be *text* or *binary*
- Files can be opened for read, write or append
 - 'r', 'w', 'a+'
- **with open('name') as filein:**
 - Allows automatic file closure
 - Explanation of the **with** statement:
<http://effbot.org/zone/python-with-statement.htm>

Reading / Writing Files

```
file_in=open('indata.txt','r')
file_out=open('outdata.txt','w')
for line in file_in:
    # Take some information (split() method is very useful!)
    x = float(line.split()[0])
    # Apply a given function (fact)
    fx = fact(x)
    # Write the result in an output file with a defined format
    file_out.write('{:010.3f}\n'.format(fx))
```

But for loading numerical data **Numpy** is more efficient. And **pandas** even more.

File parsing

- The basic:

```
for line in filein:  
    do something
```

- Common things:

```
if 'optimized' in line:  
    do something
```

```
line = line.split()
```

```
if line.upper().startswith('GEOM'): ...  
energy = float(line[2])
```

skipping lines

- Lines can be skipped by calling **next()** to a file:
for line in filein:
 if 'Optimized' in line:
 next(filein); next(filein) #skip two lines
 do something...

Formatting

- There are several function:

```
'12'.rjust(5), '12'.zfill(5)
```

- But format is more general:

```
print('{0:2d} {1:3d}'.format(x, x*x))
```

```
print("{:10.3f} {:10.3f} {:10.3f}".format(x,y,z))
```

- List of unkown length (use argument unpacking):

```
vals = np.linspace(0,1,11)
```

```
print((len(vals)*"{:10.2e} ").format(*vals))
```

<http://docs.python.org/3/library/string.html#formatspec>

Useful modules

- Similar to `ls`:

```
import glob
files = glob.glob(pattern)
```

- Working with shell-like commands:

```
import os
os.rename(src, dst)
os.mkdir(path)
os.chown(path, uid, gid)
os.getenv(key)
os.walk(directory)
```

<http://docs.python.org/3/library/os.htm>
|

Useful modules

- Reading Excel files <http://www.python-excel.org/>:

```
import xlrd
```

- Pandas uses this library

- Working with image files

- <http://scikit-image.org/>
- <http://pillow.readthedocs.io/en/latest/>

Numpy

Why Numpy / Scipy?

- Python (alone) is not efficient for numerical calculations
- Python (alone) is not practical for array manipulation
- Numpy provides the data types and methods for arrays
- Scipy provides more elaborate numerical methods
 - Optimization
 - Fast Fourier Transform
 - Linear algebra, etc

```
import numpy as np
```

```
import scipy.optimize
```

```
import scipy.stats as stats
```

numpy arrays

- without numpy:

```
> a=[[1,2],[3,4]]
```

```
> b=[[10,20], [30,40]]
```

```
> a+b
```

```
[[1, 2], [3, 4], [10, 20], [30, 40]]
```

- with numpy:

```
> a=np.array(a)
```

```
> b=np.array(b)
```

```
> a+b
```

```
array([[11, 22],[33, 44]])
```

- Array creation

```
a=np.array([1,2,3,4]).reshape([2,2])
```

```
a=np.array([[1,2], [3,4]])
```

```
a=np.zeros([2,2], dtype=int)
```

```
a[0,0]=1.
```

```
a=np.ones((4,4))
```

```
a=np.arange(10)
```

```
a=np.diag([1,2,3,4])
```

```
a=np.tile(a, (10,2))
```

```
a=np.identity(3)
```

```
a=np.linspace(-5,5, 20)
```

Ufuncs

- Unary:
`a.min()`
`a.sum()`
`a.cumsum()`
`a.mean()`
`np.argmin(a)`
`np.exp(-a)`
`np.cov(a)`
`a.tolist()`
- Binary:
`a + b`
`np.dot(a, b)` same as `a@b`
- Applying to parts of an array:
`> a=np.array([[1,2], [3,4]])`
`> a.min(axis=0)`
`array([1, 2])`
`a.sum(axis=1)`
`array([3, 7])`
- Python functions are less efficient than numpy functions:
`a.sum()` faster than `sum(a)`
`np.min(a)` faster than `min(a)`

many implemented as methods and functions

Accessing array elements

- Slicing:

```
> a[2:5]
```

```
> b[:, ::5]
```

```
> a[1:4, ...]
```

- Fancy indexing:

- Boolean arrays (masks):

```
> a = np.arange(10,15)
> indices = (a**2 > 115) & (a < 14)
> a[indices]
array([11, 12, 13])
```

- With lists:

```
> a = np.arange(10,15)
> y=a[[4,4,1]]
> y
array([14, 14, 11])
> a[[4,4,1]] = [-2, -4, 5]
> a
array([10,  5, 12, 13, -4])
```


Accessing array elements

```
>>> a[0,3:5]  
array([3,4])
```

```
>>> a[4:,4:]  
array([[44, 45],  
       [54, 55]])
```

```
>>> a[:,2]  
array([2,12,22,32,42,52])
```

```
>>> a[2::2,::2]  
array([[20,22,24]  
       [40,42,44]])
```

0	1	2	3	4	5
10	11	12	13	14	15
20	21	22	23	24	25
30	31	32	33	34	35
40	41	42	43	44	45
50	51	52	53	54	55

Accessing array elements

```
>>> a[(0,1,2,3,4),(1,2,3,4,5)]  
array([ 1, 12, 23, 34, 45])
```

```
>>> a[3:,[0, 2, 5]]  
array([[30, 32, 35],  
       [40, 42, 45]],  
      [50, 52, 55])
```

```
>>> mask = array([1,0,1,0,0,1],  
                 dtype=bool)
```

```
>>> a[mask,2]  
array([2,22,52])
```

0	1	2	3	4	5
10	11	12	13	14	15
20	21	22	23	24	25
30	31	32	33	34	35
40	41	42	43	44	45
50	51	52	53	54	55

Accessing array elements

- Slices return views

```
> a = np.arange(5)
> y=a[2:5]
> y *= -1
> a
array([ 0,  1, -2, -3, -4])
> y.flags.owndata
False
```

- np.where

```
> np.where((a>=2)&(a<4), a**2, -1)
```

```
Array([-1, -1,  4,  9, -1])
```

- np.choose

- Powerful, but complex!

- np.nonzero

- Boolean arrays return copies

```
> a = np.arange(5)
> y = a[a>1]
> y *= -1
> a
array([0, 1, 2, 3, 4])
> y.flags.owndata
True
```

- Fancy indexing returns copies:

```
> a = np.arange(5)
> y=a[[2,3,4]]
> y *= -1
> a
array([0, 1, 2, 3, 4])
> y.flags.owndata
True
```

Broadcasting

```
> a = 4.  
> b = np.array([1,2,3])  
> c = np.array([[1,2,3], [4,5,6]])  
> b+a, c+a  
(array([ 5.,  6.,  7.]), array([[ 5.,  6.,  7.],  
                                [ 8.,  9., 10.])))  
> b+c  
array([[2, 4, 6],  
       [5, 7, 9]])  
> c.dot(b)  
> b.dot(c)  
ValueError: objects are not aligned  
> b[1:]*c  
ValueError: operands could not be broadcast together with shapes (2) (2,3)  
> b[1:]*c.T  
• Use matrix if you want more algebra-like behaviour  
• Python 3.5 introduces a@b as a dot product abbreviation
```

Broadcasting rules

- When operating on two arrays, NumPy compares their shapes element-wise. It **starts with the trailing dimensions**, and works its way forward. Two dimensions are compatible when

- ✓ they are equal, or
- ✓ one of them is 1

- <https://docs.scipy.org/doc/numpy/user/basics.broadcasting.html>

- Examples OK

✓ A (3d array): 15 x 3 x 5
B (2d array): 3 x 1
Result (3d array): 15 x 3 x 5

✓ A (3d array): 15 x 3 x 5
B (3d array): 15 x 1 x 5
Result (3d array): 15 x 3 x 5

✓ A (4d array): 8 x 1 x 6 x 1
B (3d array): 7 x 1 x 5
Result (4d array): 8 x 7 x 6 x 5

- Example not OK

A (2d array): 2 x 1
B (3d array): 8 x 4 x 3

Broadcasting

- Change the shape to allow for broadcasting:

```
> c = np.array([[1,2,3], [4,5,6]])  
> b = c.mean(axis=1)  
> c+b[:,np.newaxis] #or c+b[:,None]  
> c+b.reshape((-1,1))
```

- Or keep the shape:

```
> b = c.mean(axis=1, keepdims=True)  
> c+b
```

- See also:

- `np.atleast_2d`, `np.atleast_1d` and `np.atleast_3d`

np.einsum

- Complex but powerful function to avoid the use of loops
 - Dot product, outer product, and others can be written as **einsum**

```
> c = np.array([[1,2,3], [4,5,6]])
```

```
> np.allclose(c.dot(c.T), np.einsum('ij, kj->ik',c,c))
```

True

- See numpy documentation and the following blog entry:
 - <http://docs.scipy.org/doc/numpy-1.10.0/reference/generated/numpy.einsum.html>
 - <http://ajcr.net/Basic-guide-to-einsum/>

array functions and methods

- Array reduction and logical operations:

```
> a=np.arange(5)
> np.all(a>3)
False
> np.any(a>3)
True
> a > 3
array([False, False, False, False,
       True], dtype=bool)
> (a > 3) & (a < 5)
array([False, False, False, False,
       True], dtype=bool)
```

- Some details of memory use:

- `> a.flags`

```
C_CONTIGUOUS : True
F_CONTIGUOUS : True
OWNDATA : True
WRITEABLE : True
ALIGNED : True
UPDATEIFCOPY : False
```


Loading and saving data

- Pickle is the usual way to save and restore data in Python
- We often have data file in text format:

```
#Dist Energy  
1.0 34.  
1.2 38.  
2.4 42.
```
- ```
> f=np.loadtxt("energies.dat")

> f
array([[1. , 34.],
 [1.2, 38.],
 [2.4, 42.]])
```
- Pandas adds more flexibility
- Save single arrays with:  

```
> np.save('result_y', y)
```
- Save in text mode with:  

```
> np.savetxt('result_y', y)
```
- and multiple arrays with (saves a dictionary):  

```
> np.savez('results', x, y)
```
- Recover them with load:  

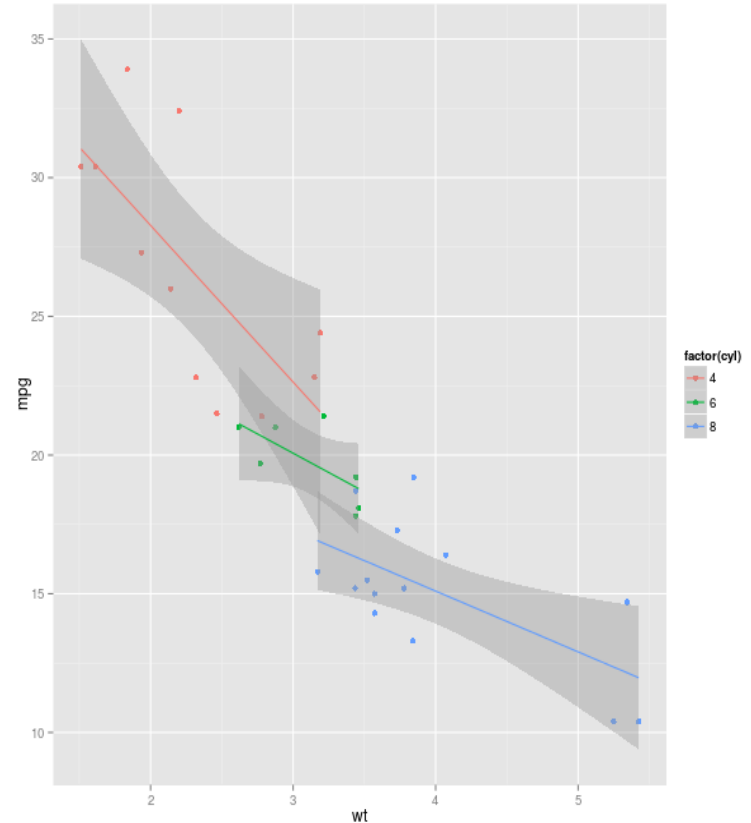
```
> y=np.load('results_y.npy')
> npz=np.load('results.npz')
```

# Acess R from python

- Use the rpy2 module.
- From the documentation:

```
import math, datetime
import rpy2.robjects.lib.ggplot2 as ggplot2
import rpy2.robjects as ro
from rpy2.robjects.packages import importr
base = importr('base')
datasets = importr('datasets')
```

```
mtcars = datasets.data.fetch('mtcars')['mtcars']
pp = ggplot2.ggplot(mtcars) + \
 ggplot2.aes_string(x='wt', y='mpg', col='factor(cyl)') + \
 ggplot2.geom_point() + \
 ggplot2.geom_smooth(ggplot2.aes_string(group = 'cyl'),
 method = 'lm')
pp.plot()
```



# Other tutorials

- Take a look at these tutorials:
  - [http://wiki.scipy.org/Tentative\\_NumPy\\_Tutorial](http://wiki.scipy.org/Tentative_NumPy_Tutorial)
  - From: <http://jrjohansson.github.io/>
    - Lecture-2-Numpy.ipynb
    - Lecture-3-Scipy.ipynb

# Matplotlib

# Matplotlib

- A module for plotting 2D and 3D data
- Combines well with numpy
- Starts with

```
import matplotlib.pyplot as plt
%matplotlib inline
```

```
import pylab or similar is deprecated.
```

# Matplotlib

Simplest plots:

```
> plt.plot([1,2,3], [1,4,9])
> plt.plot(x, sin(x), '--') #where x is a numpy array
> plt.figure() # creates new figure
> plt.clf() # Clears current figure
> plt.matshow(m) # m is a 2D array
> plt.imshow(m) # m is a 2D array. Similar to matshow.
> d = np.loadtxt('data.txt')
> plt.plot(d[:,0], d[:,1], 's') #just slightly longer than
gnuplot
```

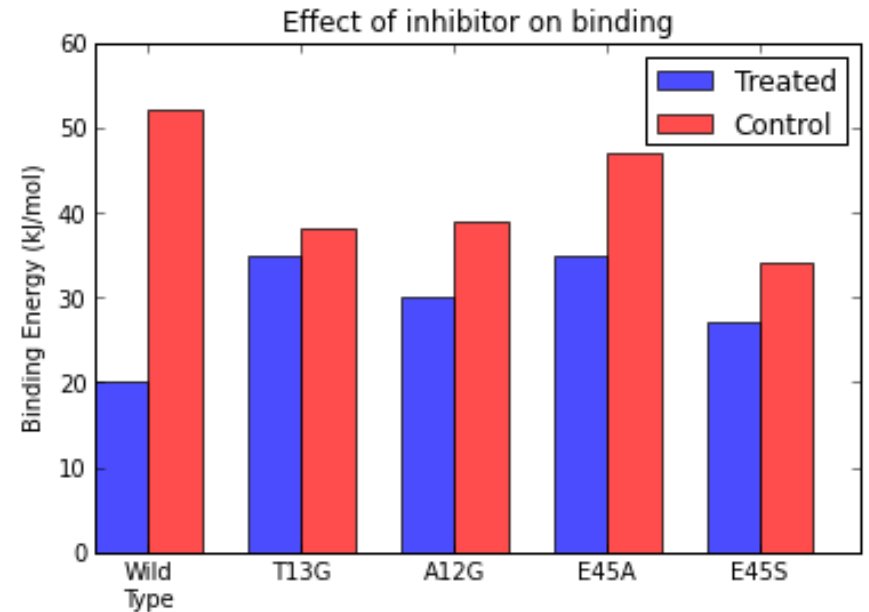
# Matplotlib

Totally reproducible  
figures

```
N = 5
treated = (20, 35, 30, 35, 27)
control = (52, 38, 39, 47, 34)
ind = np.arange(N) # the x locations for the groups
width = 0.35 # the width of the bars

fig, ax = plt.subplots()
rects1 = ax.bar(ind, treated, width, color='b', alpha=0.7, label='Treated')
rects2 = ax.bar(ind+width, control, width, color='r', alpha=0.7, label='Control')

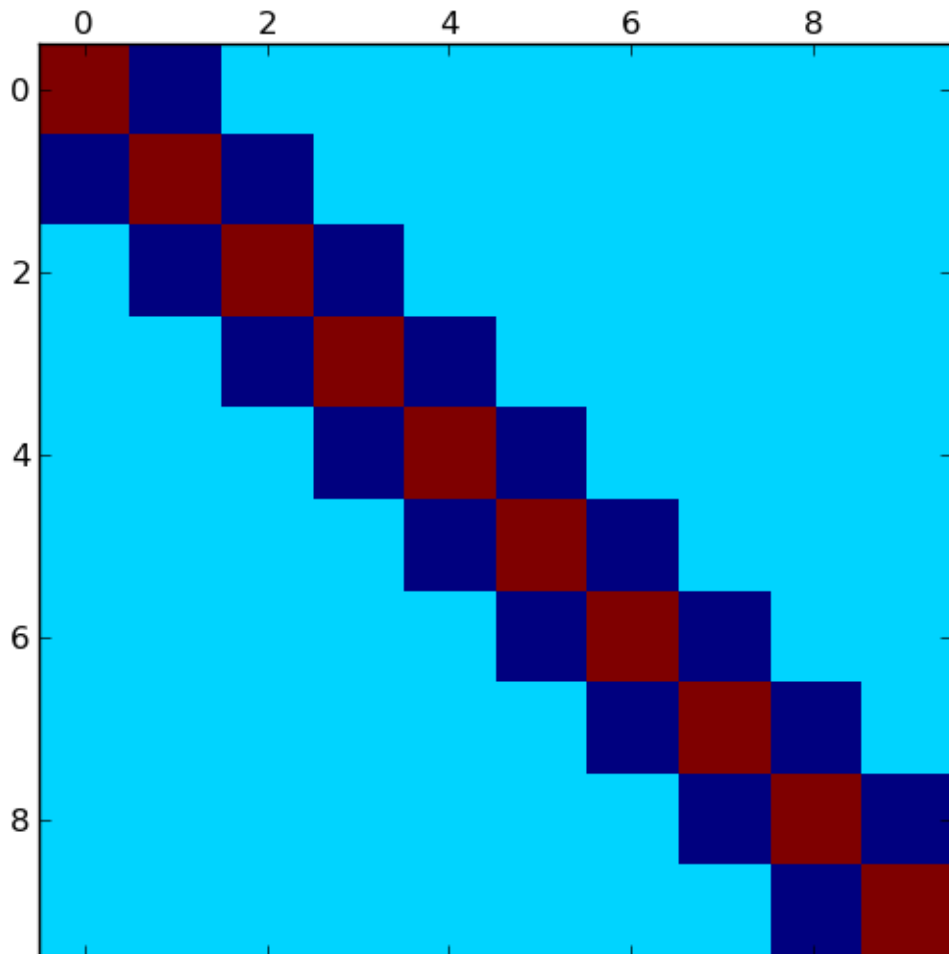
add some
ax.set_ylabel('Binding Energy (kJ/mol)')
ax.set_title('Effect of inhibitor on binding')
ax.set_xticks(ind+width)
ax.set_xticklabels(('Wild\nType', 'T13G', 'A12G', 'E45A', 'E45S'))
ax.legend()
```



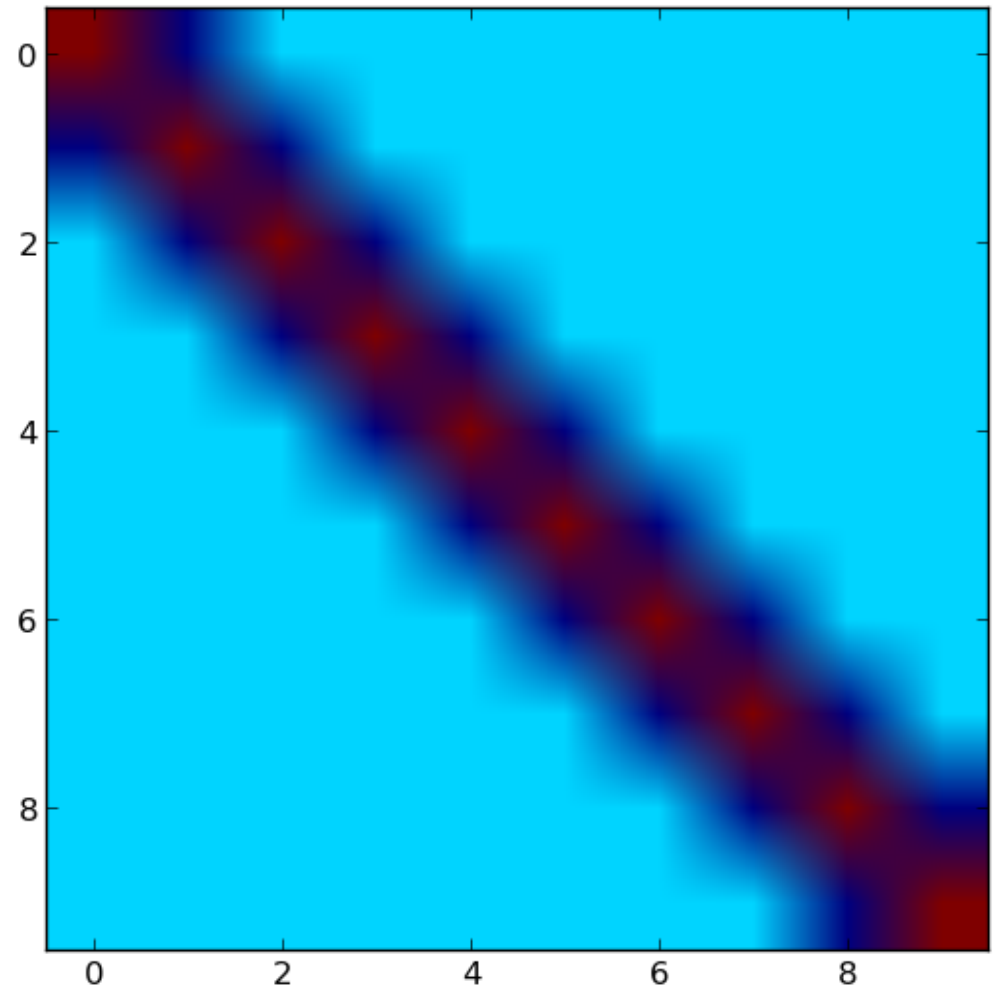
# Plotting matrices

```
m=np.diag(2*np.ones(10))+np.diag(-1*np.ones(9),1)+np.diag(-1*np.ones(9), -1)
```

matshow(m)



imshow(m)



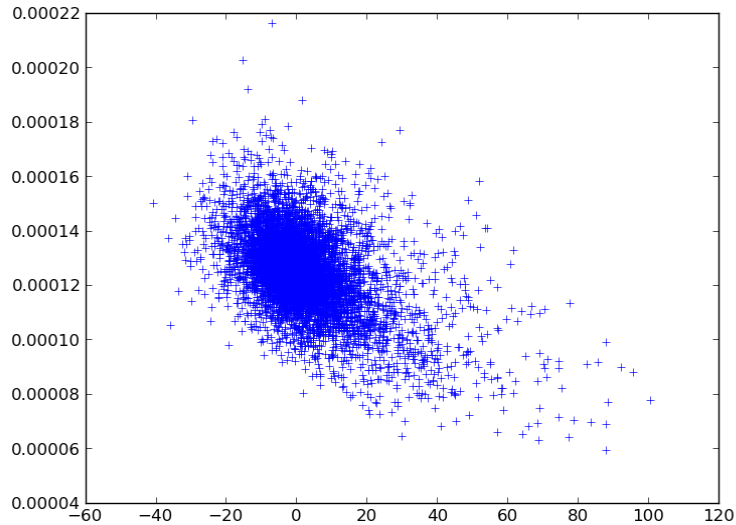


# Matplotlib styles

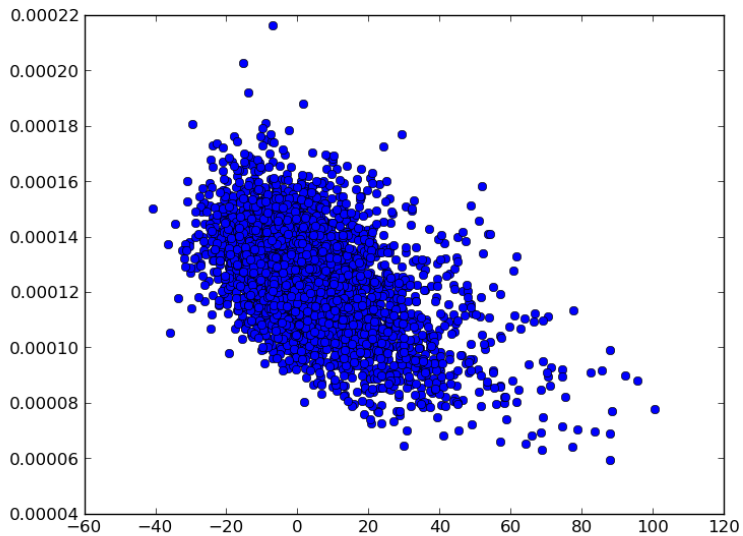
- Since version 1.5 several default styles.
- Try the following code

```
x= np.linspace(0, np.pi, 100)
for s in plt.style.available:
 with plt.style.context(s):
 plt.figure()
 plt.title(s)
 plt.plot(x,np.sin(x)*np.cos(x**2), label='A')
 plt.plot(x,np.sin(x)*np.cos(x**2)*np.cos(x), label='B')
 plt.plot(x,np.sin(x)-np.cos(x)*np.sin(x), label='C')
 plt.legend(loc='best')
```

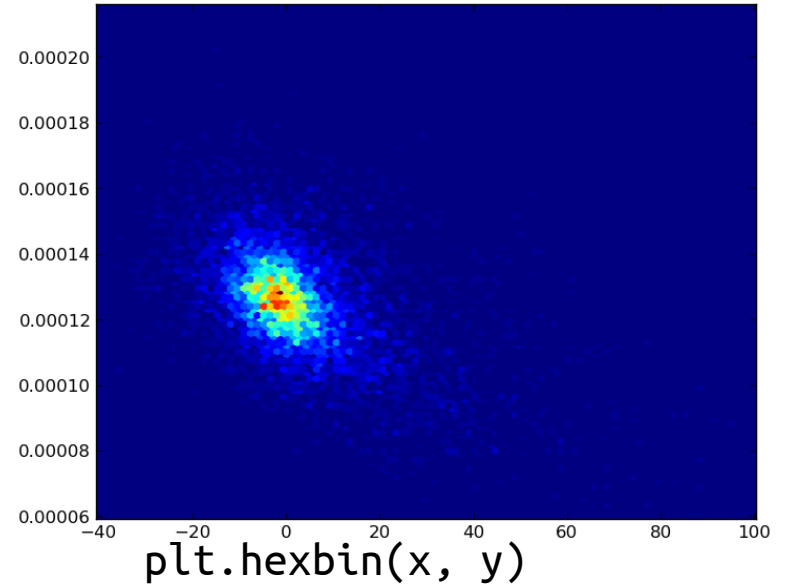
# Plotting lots of points:hexbin



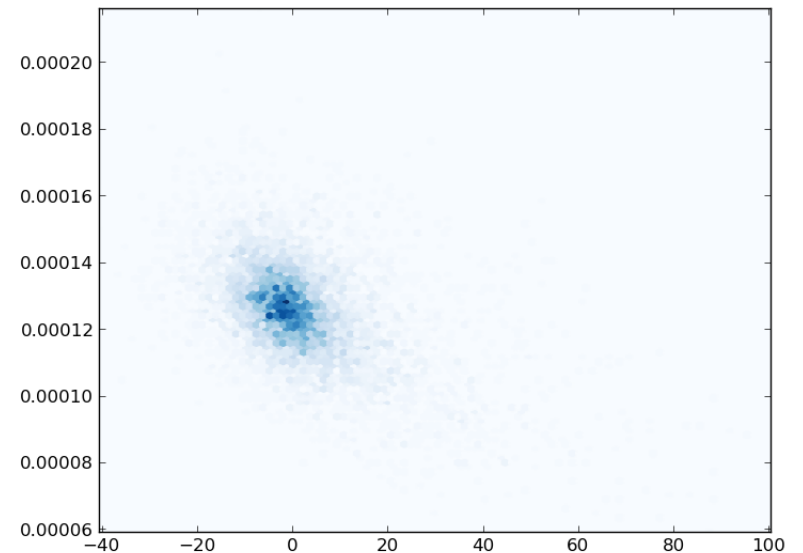
`plt.plot(x, y, '+')`



`plt.plot(x, y, 'o')`

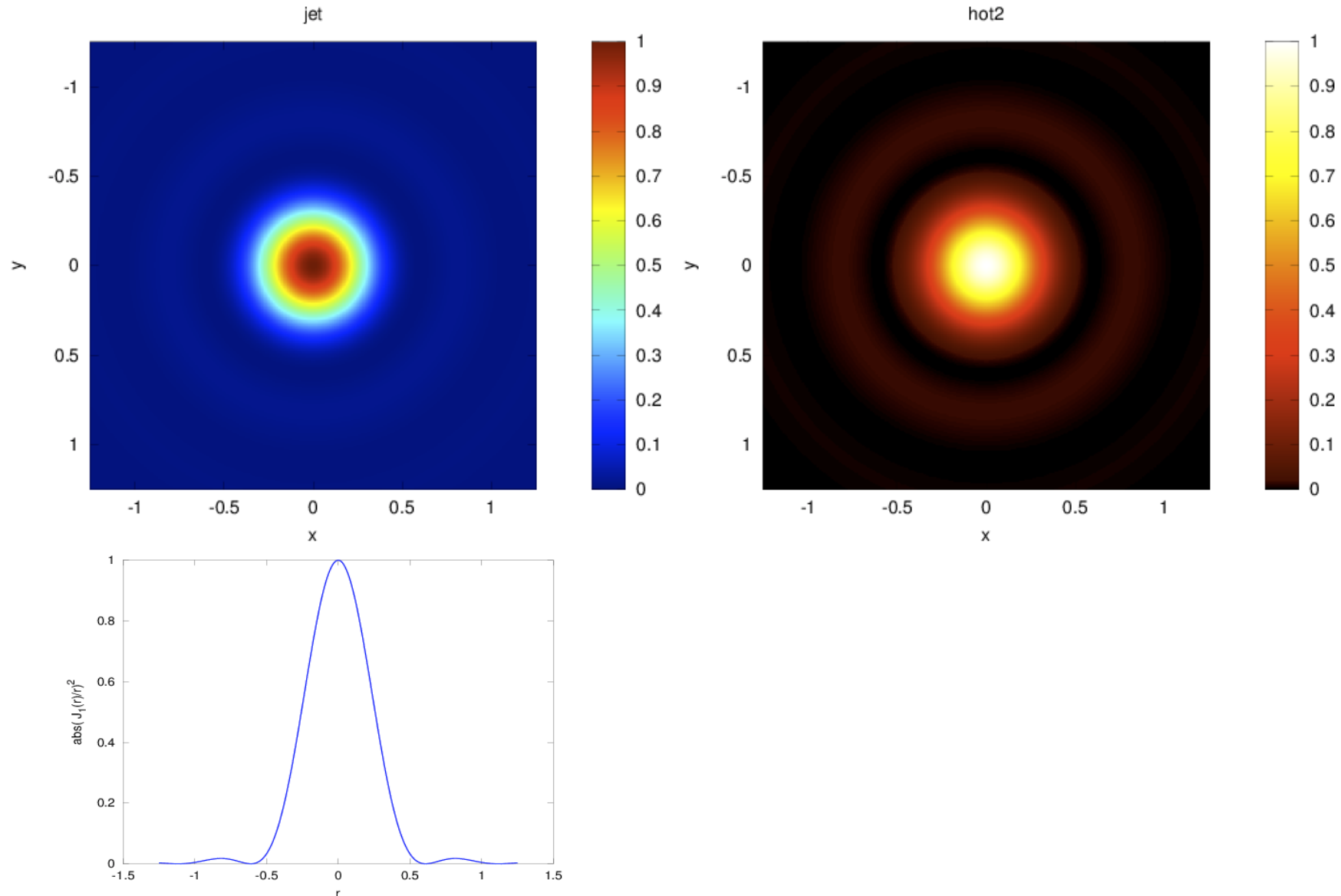


`plt.hexbin(x, y)`



`plt.hexbin(x, y, cmap=pylab.cm.Blues)`

# Jet is not a good colormap



<http://cresspahl.blogspot.com.es/2012/03/expanded-control-of-octaves-colormap.html>  
<https://jakevdp.github.io/blog/2014/10/16/how-bad-is-your-colormap/>

# Matplotlib

- Do Lecture-4-Matplotlib.ipynb from <http://jrjohansson.github.io/>
  - Other interesting material there...
- Check matplotlib gallery
  - <http://matplotlib.org/gallery.html>
- Quick reference of symbols and colours:
  - <http://www.loria.fr/~rougier/teaching/matplotlib/#quick-references>  
(part of a larger tutorial)
- Some more tricks and examples:
  - <http://wiki.scipy.org/Cookbook/Matplotlib>

# Extensions

- Seaborn
  - Data visualization
  - Nice color palettes from <http://colorbrewer2.org/>
  - <https://stanford.edu/~mwaskom/software/seaborn/>
- Bokeh:
  - <http://bokeh.pydata.org/en/latest/>
- Plotly:
  - <https://plot.ly/>
- <http://pbpython.com/visualization-tools-1.html>

# Plotting structured data

- If your data has structure: pandas dataframe
  - matplotlib is too imperative
  - Use: pandas, seaborn, altair, ggplot
  - Read:
    - <https://dsaber.com/2016/10/02/a-dramatic-tour-through-pythons-data-visualization-landscape-including-ggplot-and-altair/>
- View:
  - <https://speakerdeck.com/jakevdp/pythons-visualization-landscape-pycon-2017>

# Functions and modules

# Functions

defined by **def** and a colon:

```
def add(x,y):
 return x+y
```

Remember indentation!

Automatic (and recommended)  
documentation:

```
def add(x,y):
 """ Returns the
 sum of 2 numbers """
 return x+y
```

- Arguments are passed by reference
- there is access to global variables:

```
> def x_val(): print(x)
> x=60
> x_val()
60
```



# Functions II

Function variables are local :

```
> def x_val():
```

```
... x=40
```

```
... print(x)
```

```
> x=60
```

```
> x_val()
```

```
40
```

```
> x
```

```
60
```

to assign variables, use **return**

```
def x_val():
```

```
... x=40
```

```
... print(x)
```

```
... return x
```

```
> x = x_val()
```

```
40
```

```
> x
```

```
40
```

# Functions III

Mutable objects are passed by reference:

```
> def square_0(lst):
... lst[0]*=lst[0]

> a=[3,2,1]

> square_0(a)

> a
[9,2,1]
```

Copy variables that need to be preserved:

```
> a_copy=a[:]
> square_0(a)
> import copy
> a_copy=copy.deepcopy(a)
```

# Functions IV

Functions can have default arguments :

```
> def submit(job, priority=10, nprocs=1):
... pass
> submit('job1.sh')
```

Function arguments do not have explicit types.

```
> add('Python ', 'summerschool')
Python summerschool
```

Functions can be recursive

```
def fact(n):
 if n == 1:
 return 1
 else:
 return n * fact(n-1)
```

# Argument unpacking

Starred arguments are tuples that collect positional arguments :

```
> def prod(*args): ...
> prod(2,3,4)
> x = (4, 5, 6)
> prod(*x)
```

In prod, args=(2,3,4)

Keyword arguments can be passed as a dictionary:

```
> options = dict(paper='A4', color =
 True)
print_setup(options)
```

Unpacking can be a convenient way to print a list:

```
> vals = [1,2,3,4,5]
> print((4*'{:03d} ').format(*vals))
001 002 003 004
```

<https://docs.python.org/3/tutorial/controlflow.html#unpacking-argument-lists>

# Python optional arguments

Def statements are executed  
when function is defined:

```
> def f(arg=[]):
...

```

Therefore `arg` is associated to a  
unique list, and not to a  
different list for each function  
call.

Solution:

```
> def f(arg=None):
 if arg is None:
 arg = []
 ...

```

**arg** is now a local variable that is  
created for each function call

Explained here:

<http://docs.python-guide.org/en/latest/writing/gotchas/>

And here with more gory details:

<http://effbot.org/zone/default-values.htm>

# Python Functions III

Functions can have default arguments:

```
> def submit(job, priority=10, nprocs=1):
... pass

> submit('job1.sh')
```

These arguments become **optional** and take the default value when absent

Function arguments do not have explicit types.

```
> add('Python ', 'summerschool')
Python summerschool
```

# Lists or iterators?

- Lists are iterable objects
- Iterators generate objects on-the-fly
- Iterators can be created with a generator function
  - Uses **yield** statement
- Relevant for efficiency

```
def rang_llista(n):
 result = []
 i = 0
 while i < n:
 result.append(i)
 i += 1
 return result
```

```
def rang_gen(n):
 i = 0
 while i < n:
 yield i
 i += 1
```

# Modules

- Modules allow packing libraries or extensions
- There are built-in and external modules
- When imported modules are executed
- Modules can be written in C or Fortran!

```
> import math
```

```
> m = math
```

```
> import math as m
```

```
> from math import cos, sin
```

```
> from math import * #dangerous. All into the same namespace
```



# Modules

- Python checks if a module is already loaded.
  - The interpreter does not reload a module already imported
  - This can cause unexpected behaviour interactively
- Ipython has a more versatile module loading

```
%load_ext autoreload
```

```
autoreload 2 #Will reload a module if it changes
```

# Some useful modules

- `sys` — System-specific parameters and functions
- `os` — Miscellaneous operating system interfaces
- `os.path` — Common pathname manipulations
- `glob` — Unix style pathname pattern expansion
- `re` — regular expressions
- `copy` — Shallow and deep copy operations
- `argparse` — Parser for command-line options, arguments and sub-commands
- `subprocess` — Subprocess management
- `inspect` — Inspect live objects

# Some useful modules

```
if len(sys.argv)!=3):
 print('Error: Use two arguments.') sys.exit()

method = sys.argv[1]
filelist = glob.glob('/home/ramon/*')
for fileName in filelist:
 if os.path.isfile(fileName): print(fileName)
```

# Modules: too many...

From the python documentation:

It is also possible to use a list as a queue, where the first element added is the first element retrieved (“first-in, first-out”); however, lists are not efficient for this purpose. While appends and pops from the end of list are fast, doing inserts or pops from the beginning of a list is slow (because all of the other elements have to be shifted by one).

To implement a queue, use **collections.deque** which was designed to have fast appends and pops from both ends.

# Modules: too many...

```
>>> import math
>>> import cmath
>>> import numpy.lib.scimath as scimath
>>> math.sqrt(4)
2.0
>>> math.sqrt(-4)
Traceback (most recent call last):
 File "<stdin>", line 1, in <module>
ValueError: math domain error
>>> cmath.sqrt(4)
(2+0j)
>>> cmath.sqrt(-4)
2j
>>> scimath.sqrt(4)
2.0
>>> scimath.sqrt(-4)
2j
```

# Working with your modules

- `import` reads from local directory and from the directories in `sys.path` (`import sys` first)
- Put your modules in a directory and add it to the environment variable `$PYTHONPATH`.
- Python will add the directories in `$PYTHONPATH` to `sys.path`
- Document your modules and the functions therein.
- Use `if __name__ == '__main__':` to execute code only if Python is running the module, and not if it is imported.
  - <http://stackoverflow.com/questions/419163/what-does-if-name-main-do>

# Installing external Modules

- Use conda distribution. Then `$ conda install module`
- Many come as part of the linux distributions (usually older versions than those in conda or PyPI)
  - ipython, numpy, biopython...
- For modules in the PyPI repository (most of them)  
<https://pypi.python.org/pypi>
  - `(sudo) pip3 install module`
- Manual installation (dependencies have to be also manually installed):
  - `$ python setup.py build`
  - `$ (sudo) python setup.py install`

# Updating external Modules

- With conda `$ conda update module`
- For modules in the PyPI repository(most of them)  
<https://pypi.python.org/pypi>
  - `pip3 install -U module`
- **pip** can also be used in the conda installation.
- Remember that modules are installed for a given version of python. If you have python 2.x and 3.x you need to check for which version you are installing. For example using **pip3** or  
`$ which pip`



Scipy

# Linear algebra

- Support for LAPACK, BLAS and ATLAS
  - Can make Scipy compilation more involved

```
> A=matrix(random.rand(5,5))
```

```
> A.I
```

```
> linalg.det(A)
```

```
> linalg.eigvals(A)
```

```
> linalg.eig(A)
```

```
> linalg.svd(A)
```

```
> linalg.cholesky(A)
```

- Solving linear systems:

- $A.x=b$

```
> b=matrix(random.rand(5)).\
reshape((5,1))
```

```
> linalg.solve(A,b)
```

- LAPACK, BLAS wrappers

```
> from scipy.lib import lapack
```

```
> from scipy.lib import blas
```

```
blas.fblas.sdot?
```

# Optimization

- There are different optimization methods:

- > `import scipy.optimize as so`

- Some only need the function value:

- > `fmin, fmin_powell`

- Some need the gradient or the hessian:

- > `fmin_cg, fmin_bfgs, fmin_ncg`

- Some look for global minima:

- > `anneal`

- Remember:

- > `scipy.info('optimize')`

- 2 sources of official documentation:

- <http://docs.scipy.org/doc/scipy/reference/tutorial/optimize.html>

- <http://docs.scipy.org/doc/scipy/reference/optimize.html>

# f2py

- Many things are fast with Numpy
- Iterative algorithms over **array values** are slow
- You can import Fortran functions and subroutines with f2py
- You could also call external fortran programs with
  - > `subprocess.call(<program>, shell=True)`
    - but data exchange has to be through files (slower)
- f2py finds your fortran compiler. Works with gfortran, ifort,...
- f2py creates a module you can import in python
- As simple as:
  - `$ f2py -c <file> -m <module>`
    - Tip: first compile it to check it works

# f2py II

```
module funcs
implicit none
contains
function f1(x,y)
 real,intent(in):: x,y
 real:: f1
 f1=x+y**2
end function f1

function f2(x,y)
 real,intent(in):: x,y
 real, dimension(3):: f2
 f2(1)=x+y**2
 f2(2)=sin(x*y)
 f2(3)=2*x-y
end function f2
end module
```

```
$ f2py -c test.f90 -m test
```

- go to python:

```
> import test
> test.funcs.f1(1,2)
5.0
> test.funcs.f2(1,2)
array([5., 0.90929741, 0.],
 dtype=float32)
```

# f2py III

Using ipython magicfunctions:

```
sudo pip3 install -U fortran-magic
```

Useful for performing long array operations

```
In [5]: %load_ext fortranmagic
```

```
In [6]: %%fortran
 subroutine f1(x, y, z)
 real, intent(in) :: x,y
 real, intent(out) :: z

 z = sin(x+y)

 end subroutine f1
```

```
In [7]: f1(1.0, 2.1415)
```

```
Out[7]: 9.26574066397734e-05
```

# Big data, big memory

- Numpy arrays are meant to live in memory
- If that is not possible:
  - Use op= operations (they use half the memory):
    - `p *=alpha` is better than `p = p*alpha`
  - Use `scipy.sparse` matrices
    - <http://docs.scipy.org/doc/scipy/reference/sparse.html>
  - Use `PyTable` to store (compressed) matrices on disk
    - <http://www.pytables.org/>
  - Modify your algorithm to work with submatrices

# Sympy: Symbolic math

- Symbolic algebra
- Analytic solution of equations
- Integration, derivation
- Polynomials
- Limits

Alternate forms:

`(cos(x + y)).expand(trig=True)`

$$-\sin(x)\sin(y) + \cos(x)\cos(y)$$

`trigsimp(cos(x + y))`

$$\cos(x + y)$$

`(cos(x + y)).rewrite(csc, sin, sec, cos, cot, tan)`

$$\frac{-\tan^2\left(\frac{x}{2} + \frac{y}{2}\right) + 1}{\tan^2\left(\frac{x}{2} + \frac{y}{2}\right) + 1}$$

`(cos(x + y)).rewrite(sin, exp, cos, exp, tan, exp)`

$$\frac{1}{2}e^{i(-x-y)} + \frac{1}{2}e^{i(x+y)}$$

<http://sympy.org/en/index.html>

```
>>> integ = Integral(sin(x**2), x)
```

```
>>> integ
```

$$\int \sin(x^2) dx$$

```
>>> integ.doit()
```

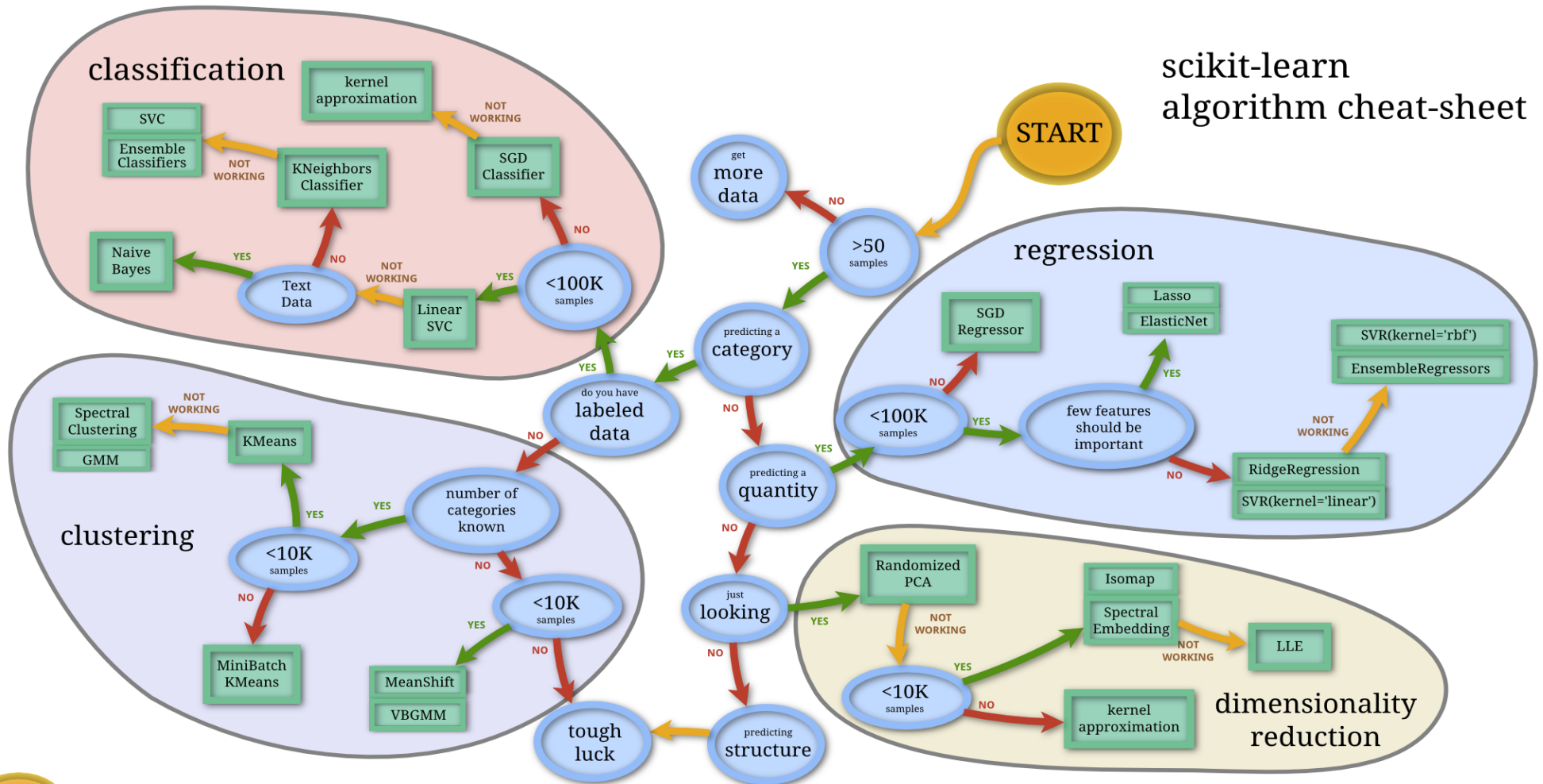
$$\frac{3 \cdot \sqrt{2} \cdot \sqrt{\pi} \cdot \text{fresnels}\left(\frac{\sqrt{2} \cdot x}{\sqrt{\pi}}\right) \cdot \Gamma(3/4)}{8 \cdot \Gamma(7/4)}$$



# Machine Learning

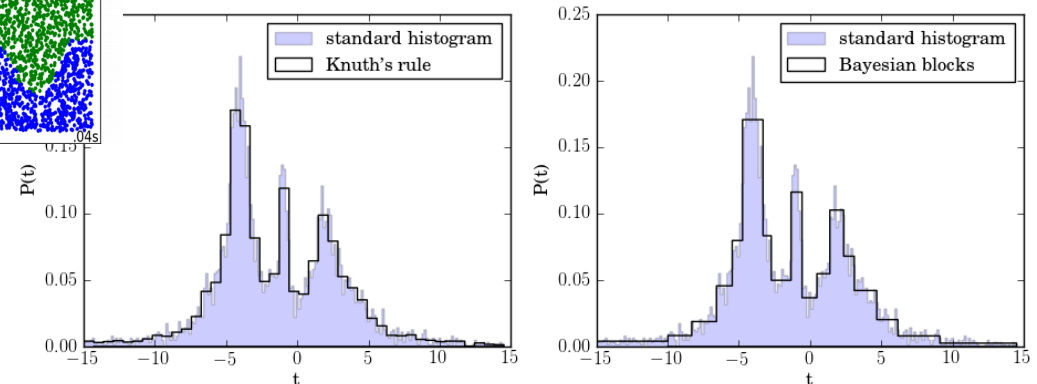
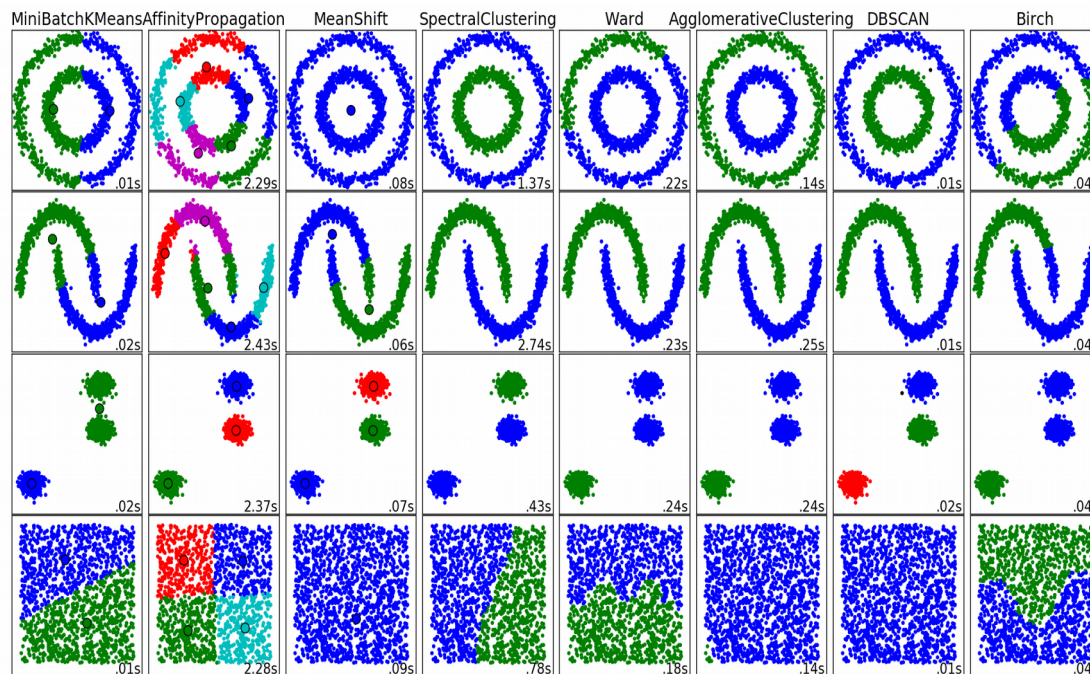
# scikit-learn

scikit-learn  
algorithm cheat-sheet



# scikit-learn and AstroML

- scikit-learn
  - Very well documented
- AstroML
  - Designed for astronomy
  - Builds on top of scikit-learn



# Deep learning & Bayesian Modelling

- Deep Learning
  - Caffe2
  - Theano
    - Lassagne
  - Tensorflow
  - Keras
- Bayesian Modelling
  - pyMC3

# Optimization and debugging

# Optimization

- “Premature optimization is the root of all evil”  
Knuth
- `%timeit a=np.random.random(100000)`
- `a=np.random.random(100000)`  
`n_dim=3`  
`%%timeit`  
`x=np.zeros(shape=(100000,n_dim),order='F')`  
`for j in range(0,n_dim):`  
`x[:,j]=a*j`
- Evaluated in a separate environment

# Profiling

- Use line-profiler:
- `pip3 install line_profiler`  
`%load_ext line_profiler`
- `lprun -f func2 func1()`
- Memory profiling with:
- `pip3 install memory_profiler`  
`%load_ext memory_profiler`
- `%mprun -f func2 func1()`

# Exceptions and errors

Although the language is interpreted there are some syntax errors that prevent execution:

```
def safe_divide_1(x, y)
```

```
File"/home/ramon/python/prova.py",
line 1
```

```
def safe_divide_1(x, y)
```

^

```
SyntaxError: invalid syntax
```

Exceptions leave a trace easy to follow.

Easy debugging with

```
%pdb
```

```
%debug
```



# pdb: python debugger

```
In [1]: pdb
```

```
Automatic pdb calling has been turned ON
```

```
In [4]: run foo.py
```

```
NameError: name 'b' is not defined
```

```
> /home/ramon/python/foo.py(2)<module>()
```

```
1 a = 3
```

```
----> 2 print(b)
```

```
ipdb> ?
```

# pdb: python debugger

```
In [9]: run foo.py
```

```

NameError Traceback (most recent call last)
/home/ramon/python/foo.py in <module>()
 1 a = 3
----> 2 print(b)
```

```
NameError: name 'b' is not defined
```

```
In [10]: %debug
> /home/ramon/python/foo.py(2)<module>()
 1 a = 3
----> 2 print(b)
ipdb>
```

# Numba

- Numba compiles in a virtual machine.
- Developed by Continuum analytics, so easiest install from conda.
- **\$ conda install numba**

# Cython

- An extension to python that generates C code that can be compiled
- Available in most linux distributions
- See also:
- <https://jakevdp.github.io/blog/2013/06/15/numba-vs-cython-take-2/>

# Other alternatives

- Use Julia
  - A different language
  - Close in syntax to Python
- Theano: “define, optimize, and evaluate mathematical expressions involving multi-dimensional arrays efficiently”
  - <https://theano.readthedocs.org/en/latest/>
- Parakeet: a runtime compiler for scientific computing in Python
  - <http://www.parakeetpython.com/http://www.parakeetpython.com/>
- Just-in-time compilers for number crunching in Python
  - <http://www.phi-node.com/2013/01/just-in-time-compilers-for-number.html>
- See also the notebooks here:
  - [http://nbviewer.jupyter.org/github/rasbt/One-Python-benchmark-per-day/tree/master/ipython\\_nbs/](http://nbviewer.jupyter.org/github/rasbt/One-Python-benchmark-per-day/tree/master/ipython_nbs/)

# Add-ons

# Add ons: Biopython

## Biopython

```
from Bio.PDB import *
```

```
p=PDBParser(PERMISSIVE=1)
```

```
s=p.get_structure('1OJR', filename)
```

Print out the coordinates of all CA atoms with B factor > 50:

```
for model in s.get_list():
```

```
 for chain in model.get_list():
```

```
 for residue in chain.get_list():
```

```
 if residue.has_id("CA"):
```

```
 ca=residue["CA"]
```

```
 if ca.get_bfactor()>50.0:
```

```
 print ca.get_coord()
```

<http://biopython.org>

# Add ons: Machine learning and statistics

- Basic statistics in scipy.stats
  - Tutorial:  
<http://docs.scipy.org/doc/scipy/reference/tutorial/stats.html>
  - Reference: <http://docs.scipy.org/doc/scipy/reference/stats.html>
- Machine learning with sklearn
  - <http://scikit-learn.org/stable/>
  - Choosing the method:  
[http://scikit-learn.org/stable/tutorial/machine\\_learning\\_map/](http://scikit-learn.org/stable/tutorial/machine_learning_map/)
- More algorithms (and a textbook) with AstroML
  - <http://www.astroml.org/>



# Add ons: itertools

```
> import itertools
> perms = itertools.permutations('ABC', 3)
> list(perms)
[('A', 'B', 'C'),
 ('A', 'C', 'B'),
 ('B', 'A', 'C'),
 ('B', 'C', 'A'),
 ('C', 'A', 'B'),
 ('C', 'B', 'A')]
> list(itertools.combinations('ABC',2))
[('A', 'B'), ('A', 'C'), ('B', 'C')]
```

# Add ons: active papers

ActivePapers is a framework for doing and publishing reproducible research. An ActivePaper is a file that contains code (Python modules and scripts) and data (HDF5 datasets), plus the dependency information between all these pieces. You can change a script and re-run all the computations that depend on it, for example. Once your project is finished, you can publish the ActivePaper as supplementary material to your standard paper.

<http://khinsen.wordpress.com/2013/09/27/activepapers-for-python/>

# Resources

# Resources

On-line Official documentation (contains Tutorial in PDF or HTML):

<http://www.python.org/doc>

General introductory books (also in paper):

<http://diveintopython.org/> (This one is simpler!)

<http://www.greenteapress.com/thinkpython/thinkpython.html>

Comparison of codes in different languages:

<http://rosetacode.org>

<http://www.codecodex.com>

Python package index: where to find modules

<http://pypi.python.org/pypi>

# Resources

- Interactive tutorial
  - <http://pythonmonk.com/>
- A Crash Course in Python for Scientists (with applications in Quantum chemistry)
  - <http://nbviewer.ipython.org/5920182>
  - Written in an ipython notebook
- Python Scientific Lecture notes
  - <http://scipy-lectures.github.io/>
- Python flow with Pythontutor
  - <http://www.pythontutor.com>

# Resources: Books

- Rossant, C, *Learning Ipython for Interactive Computing and Data Visualization*.
  - Basic level. Covers several subjects, including matplotlib and parallelism. Recipes book.
- Stewart, J.M., *Python for Scientists*
  - Basic level. Unfortunately in Python 2. Covers a lot on differential equations.
- DeCaria A. J. *Python Programming and Visualization for Scientists*
  - DeCaria teaches Python programming and visualization for meteorology and ocean sciences majors.
- Packt Publishing. Wide variety, lots on GIS and Python.
- <https://wiki.python.org/moin/AdvancedBooks>

# Resources: Video Tutorials

- Check:  
<https://www.youtube.com/user/EnthoughtMedia>
- Check youtube videos of PyCon
- Check Scipy Conference and Euroscipy:
  - <http://conference.scipy.org/proceedings/scipy2015/>

# Resources: MOOCs

- General Python programming:
  - <https://www.coursera.org/course/programming1>
  - <https://www.coursera.org/course/programming2>
- Advanced scientific programming with Fortran, Python, OMP, OpenMPI...
  - <https://www.coursera.org/course/scicomp>



# Resources: Teaching

- On teaching programming with Python 3

<http://www.comp.leeds.ac.uk/nde/papers/teachpy3.html>

- Online Syntax Highlighting

<http://tohtml.com/python/>

- Style Guide for Python Code:

- [www.python.org/dev/peps/pep-0008/](http://www.python.org/dev/peps/pep-0008/)

# K. Hinsen views

- “NumPy has introduced incompatible changes with almost every new version over the last years”
- “Given the importance of NumPy in the scientific Python ecosystem, I consider its lack of stability alarming”.
- “What makes me hesitate to recommend not using Python is that there is no better alternative”.
- <https://khinsen.wordpress.com/2014/09/12/the-state-of-numpy/>

# Jake VanderPlas

- Great blog about python with applications in
  - <https://speakerdeck.com/jakevdp/the-unexpected-effectiveness-of-python-in-science>
  - Science
  - Statistics
  - Cycling...
  - All entries are jupyter notebooks.
  - <https://jakevdp.github.io/>
  - See also his book and library on machine learning:
  - <http://www.astroml.org/>
  - <http://press.princeton.edu/titles/10159.html>

# Software in python

- QM/MM with pDynamo: <http://www.pdynamo.org>
- MM with MMTK: <http://dirac.cnrs-orleans.fr/MMTK/>
- Molecular visualization:
  - VMD: <http://www.ks.uiuc.edu/Research/vmd/>
  - pymol: <http://www.pymol.org/>
- QM calculation with
  - pyQuante: <http://pyquante.sourceforge.net/>
  - NWChem: <http://www.nwchem-sw.org/index.php/Python>
- Protein structure with pyRosetta: <http://pyrosetta.org/>
- Bioinformatics with BioPython: <http://biopython.org/>

# Python for modelling

- cclib: <http://cclib.github.io/>
- ORBKIT: <http://orbkit.github.io/>
- Nglview, chemical structures in jupyter:  
<https://github.com/arose/nglview>
- Trajectory analysis:
  - MDtraj : <http://mdtraj.org>
  - MDAnalysis: <http://www.mdanalysis.org/>
  - Pytraj: <https://github.com/Amber-MD/pytraj>