Lecture Notes 4: Sampling, Simulation, Low-Level Extensions

Sampling from Distributions

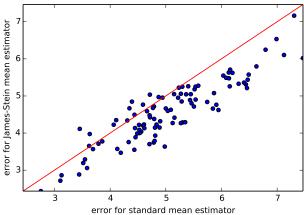
```
In [1]: import numpy,numpy.random
    Sampling from a uniform distribution between 0 and 1
In [2]: print numpy.random.uniform(0,1,[10])
[ 0.72611138     0.24067817     0.71938139     0.45283454     0.70368496     0.94023131     0.77130642     0.04720969     0.29922044     0.66994318]
    Sampling from normal distibution of mean 10 and standard deviation 0.01
In [3]: print numpy.random.normal(10,0.01,[10])
[ 10.00598152     10.00735394     10.00212889     9.9986185     10.00672975     9.98147864     10.00094005     9.995714     10.00350466     10.00550996]
```

Demo 1: Testing A Statistical Paradox (the James-Stein Paradox)

When estimating the mean of a data distribution, moving the estimator from the empirical mean, and closer to some arbitrary point (e.g. the origin) makes the estimator more accurate. Let's verify it:

```
In [6]: # Testing many times:
       errstd = []
       errstd2 = []
       n = 10
       d = 50
       for i in range(100):
            # sample from a distribution of mean vector 1 and standard deviation 1
            m = 1.0
            X = numpy.random.normal(m,1,[n,d])
            # empirical mean
            m_emp = X.mean(axis=0)
            # some coefficient
            c = (1-(d-2)*1.0/n/((m_emp)**2).sum())
            # james-stein estimator
            m_js = c*m_emp
            # the error between the true mean and the standard estimator
            errstd += [((m - m_emp)**2).sum()]
            # the error between the true mean and the said better estimator
            errstd2 += [((m - m_js)**2).sum()]
```

```
In [7]: import matplotlib
        from matplotlib import pyplot as plt
        %matplotlib inline
        from IPython.display import set_matplotlib_formats
        set_matplotlib_formats('pdf','png')
        plt.rcParams['savefig.dpi'] = 90
In [9]: plt.scatter(errstd,errstd2)
        1,h = min(errstd+errstd2),max(errstd+errstd2)
       plt.plot([1,h],[1,h],color='red')
        plt.xlabel('error for standard mean estimator')
        plt.ylabel('error for James-Stein mean estimator')
       plt.axis([1,h,1,h])
Out[9]: [2.4355343709629809,
        7.4646057017558283,
         2.4355343709629809,
         7.4646057017558283]
```



Observation: most points are below the curve (i.e. James-Stein estimator is better).

Making discrete choices

Let us suppose we have 7 fruits to choose from:



```
In [10]: fruits = ['watermelon', 'apple', 'grape', 'grapefruit', 'lemon', 'banana', 'cherry']
We use the function random.choice to randomly choose from that list
```

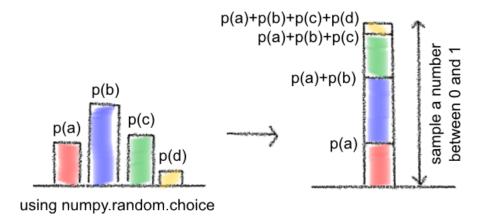
```
grapefruit
```

Another way to make discrete choices

'watermelon' 'apple' 'cherry' 'apple']

print numpy.random.choice(fruits,[20],p=p)

['grapefruit' 'apple' 'apple' 'apple' 'apple' 'apple' 'apple' 'banana' 'apple' 'watermelon' 'cherry' 'apple' 'apple' 'apple' 'apple'



```
In [13]: # Define fruits probabilities
    p = [0.05,0.70,0.05,0.05,0.05,0.05,0.05]

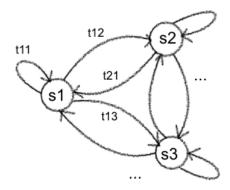
# Cumulate them
    1 = numpy.cumsum([0]+p[:-1]) # lower-bounds
    h = numpy.cumsum(p) # upper-bounds

# Draw a number between 0 and 1
    u = numpy.random.uniform(0,1)
```

```
# Find which basket it belongs to
         s = (u>1)*(u<h)
         # retrieve the label
         fruits[numpy.argmax(s)]
Out[13]: 'apple'
  The code can be parallelized to choose multiple fruits at the same time
In [14]: # generate many numbers between 0 and 1
         u = numpy.random.uniform(0,1,[20])
         # find the basket to which they belong
         na = numpy.newaxis
         s = (u[:,na]>1[na,:])*(u[:,na]<h[na,:])
         print([fruits[i] for i in numpy.argmax(s,axis=1)])
['watermelon', 'apple', 'apple', 'apple', 'apple', 'apple', 'apple', 'grapefruit', 'apple', 'banana', '
   More importantly, the code can be parallelized to choose from multiple distributions at the same time. Let us
create a matrix of probability distributions.
In [15]: P = [
             [0.0,0.05,0.05,0.05,0.05,0.05,0.70,0.05],
             [0.0,0.05,0.05,0.70,0.05,0.05,0.05,0.05]
         # (note that we have added a zero-probability state
         # at the beginning for ease of implementation)
  We stack them, and compute the bounds L and H
In [16]: C = numpy.cumsum(P,axis=1) # lower-bounds
         L,H = C[:,:-1],C[:,1:]
  Draw 8 fruits from each distribution, and test the bounds
In [19]: R = numpy.random.uniform(0,1,[3,8])
         S = (R[:,:,na] > L[:,na,:])*(R[:,:,na] < H[:,na,:])
In [20]: for s in S:
             print([fruits[i] for i in numpy.argmax(s,axis=1)])
['lemon', 'cherry', 'cherry', 'cherry', 'apple', 'cherry', 'cherry']
['banana', 'banana', 'banana', 'banana', 'grapefruit', 'cherry', 'grapefruit', 'banana']
['grape', 'grape', 'grape', 'grape', 'grape', 'grape', 'lemon', 'grape']
```

Demo 2: Monte Carlo Simulation of a Markov Chain

A Markov chain transits between a set of states, where the transition between pairs of states is associated with a fixed probability. The set of probabilities can be stored in a transition matrix.



Thanks to our parallel implementation, we can choose from multiple probability distributions, and thus, simulate multiple markov chains in parallel, each of them being in a different state.

```
In [24]: # Add empty state to transition matrix
         P = numpy.pad(T,1,mode='constant')[1:-1,:-1]
         print(P)
[[ 0.
       0.9 0.1 0. ]
ΓО.
       0.
            0.9 0.17
 Γ0.
                  0.]]
In [25]: # Implementing transition for each particle
         def mcstep(X):
             Xp = numpy.dot(X,P)
             Xc = numpy.cumsum(Xp,axis=1)
             L,H = Xc[:,:-1],Xc[:,1:]
             R = numpy.random.uniform(0,1,[len(Xp),1])
             return (R > L)*(R < H)*1.0
In [26]: # Initialize all particles to state 1
         A = numpy.outer(numpy.ones([30]),[1.0,0,0])
         print 'initial distribution of particles states: [%.3f %.3f %.3f]'%tuple(A.mean(axis=0))
         for i in range(20):
                 print '(iter %2d)'%i,numpy.argmax(A,axis=1)
                 A = mcstep(A)
         # Print distribution of sampled states
         print 'final distribution of particles states: [%.3f %.3f %.3f]' %tuple(A.mean(axis=0))
```

```
initial distribution of particles states: [1.000 0.000 0.000]
(iter 7) [0 1 0 0 1 0 0 0 2 0 0 0 0 0 1 0 1 0 0 1 1 0 2 0 1 1 0 0 1 1]
(iter 8) [0 1 0 0 1 0 0 0 0 0 0 0 0 0 1 0 1 0 0 1 1 0 0 1 2 1 0 0 1 2]
(iter 9) [0 1 0 0 1 0 0 0 0 1 0 0 0 0 1 0 1 0 0 1 1 0 0 1 0 1 0 0 1 0]
(iter 11) [0 1 0 1 0 1 0 0 0 1 0 0 0 0 1 0 2 0 0 0 1 0 0 2 0 1 0 0 1 0]
(iter 14) [0 1 0 1 1 2 0 0 0 1 0 0 1 0 1 0 0 0 1 0 2 0 1 0 0 1 0 0 1 0]
(iter 16) [0 1 0 1 1 1 0 0 0 1 0 0 0 0 1 0 0 1 0 0 0 1 0 1 0 1 0 1 0 0 1 0]
(iter 19) [0 1 0 0 1 0 0 0 0 1 0 0 0 0 1 0 0 1 0 0 1 0 1 0 1 0 1 0 0 0 0 1]
final distribution of particles states: [0.633 0.367 0.000]
```

Low-level extensions

Idea: Write highly optimized functions directly in C/C++, Fortran, or Cuda, and make them accessible to the Python user. Numpy is also based on this principle: making very efficient functions accessible in Python in a user-friendly manner.

Example: F2PY (Fortran to Python)

Main steps:

- Install gfortran (fortran compiler) and f2py (binding between Fortran and Python)
- Create file (e.g. convolution.f90) containing some optimized code for convolutions.
- Compile convolution.f90 into a Python module using f2py
- Import module convolution directly from Python.

Applying batch convolutions

We would like to use the following Fortran code that computes a batch of convolutions in the same way as the convolutional layer of a neural network.

Figure 1: file: convolution.f90

Demo 3: Filtering images to detect interesting features



```
# horizontal edge detector in reverse direction
         W[:2,:,:,1] = -1.0
         W[2:,:,:,1] = 1.0
         # vertical edge detector
         W[:,:2,:,2] = 1.0
         W[:,2:,:,2] = -1.0
         # vertical edge detector in reverse direction
         W[:,:2,:,3] = -1.0
         W[:,2:,:,3] = 1.0
         # more red than yellow detector
         W[:,:,0,4] = 1.0
         W[:,:,1,4] = -1.0
In [36]: Z = convolution.conv(X,W)
         Y = numpy.maximum(0,Z)
In [37]: f = plt.figure(figsize=(10,7))
         for i in range(4):
             p = f.add_subplot(2,2,i+1)
             p.imshow(Y[:,:,i])
                                               100
                                               150
                                               250
                        150
                                               150
                        200
                                               200
                        250
                                               250
```

In [38]: plt.imshow(Y[:,:,4])

Out[38]: <matplotlib.image.AxesImage at 0x7f9d8e9ef690>

