June 30, 2016

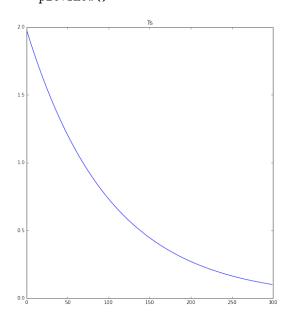
In [1]: import numpy as np

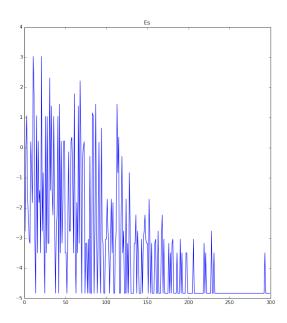
```
import matplotlib.pyplot as plt
        \% matplotlib inline
    9.1 Simulated Annealing
1
In [8]: #set initial state s randomly
       N = 6
        s = np.sign(np.random.rand(N)-.5)
        #w arbitrary, but symmetrically and with zero diagonal
        w = np.random.rand(N,N)*2-1
        for i in range(N):
            w[i,i] = 0
            for j in range(i):
                w[j,i] = w[i,j]
       print(s)
[-1. -1. 1. 1. 1. 1.]
In [3]: #the cost (\energy")
       def E(s, w):
            total = 0
            for i in range(len(s)):
                for j in range(len(s)):
                    total = total + w[i,j]*s[i]*s[j]
            return -.5 * total
        def Ei(s, w, i):
            total = 0
            for j in range(len(s)):
                total = total + w[i,j]*s[i]*s[j]
            return -.5 * total
        #The probability that the network is in a state s with energy E(s) is given by
        def P(s, w, b):
            \#print("P = ", (1.0/Z(s, w, b)), "*", np.exp(-b * E(s, w)))
            \#print("Z = ", Z(s, w, b), "->", (1.0/Z(s, w, b)))
            \#print("E = ", E(s, w))
            return (1.0/Z(s,w,b)) * np.exp(-b * E(s,w))
        #the partition function Z guarantees P (s) to be a valid probability mass function
        #is given as the sum over all possible configurations
```

```
def Z(s, w, b):
            total = 0
            for i in range(len(s)):
                total = total + np.exp( -b * E(s, w))
            return total
In [4]: def SimulatedAnnealing(s, w):
            # Initialization
            mys = s.copy()
            myw = w.copy()
            #set \beta0,\tau>1
            b0 = .5
            tau = 1.01
            #tmax
            tmax = 300
            #number of times to repeat
            \#try both M=1 and M=500
            M = 500
            # Optimization
            Ts = np.zeros(tmax)
            Es = np.zeros(tmax)
            bt = b0
            #for each iteration t = 0, \ldots, tmax
            for t in range(tmax):
                #repeat the following M times (state update loop):
                for m in range(M):
                    #select node i randomly
                    i = np.random.randint(0, N)
                    #determine the energies for the two options of state si and compute their difference
                    sci = mys.copy()
                    sfi = mys.copy()
                    sfi[i] = sfi[i] * -1
                    Ec = Ei(sci, myw, i)
                    Ef = -Ec
                    \#Ef = E(sfi, myw)
                    dE = Ef - Ec
                    #Change to new state with some probability
                    #flip state si with probability
                    Pf = 1.0 / (1 + np.exp(bt*dE))
                    if np.random.rand() < Pf:</pre>
                         mys[i] = mys[i] * -1
                     \#print(Pf)
                     #print(Pf*bt)
                #increase \beta
                bt = tau * bt
```

```
#For plotting
Ts[t] = 1.0/bt
Es[t] = E(mys, myw)
return s, Ts, Es
snew, Ts, Es = SimulatedAnnealing(s, w)
```

2 Plotting





```
In [7]: def MakeAllS(curs, depth):
    if depth == 0:
        return curs
    else:
        reso = MakeAllS(curs, depth-1)
        resn = []
        for i in range(len(reso)):
            resn.append(reso[i]+[0])
            resn.append(reso[i]+[1])
        return resn

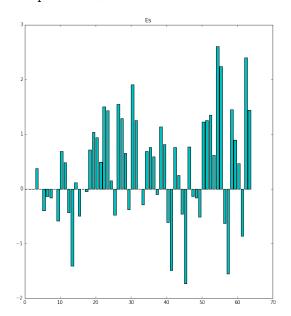
allEs = [E(sc,w) for sc in MakeAllS([[0],[1]], N-1)]

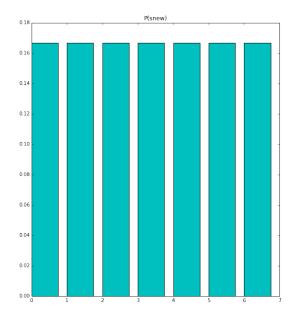
ind = np.arange(len(np.arange(len(allEs))))  # the x locations for the groups
width = .75  # the width of the bars

bs = [110, 11, 1.1, 1.01, 1.001, 1.0001, 1.00001]
```

```
cb = bs[4]
Pbs = [P(s, w, myb) for myb in bs]

f, axarr = plt.subplots(1, 2, figsize=(20,10))
axarr[0].bar(ind, allEs, width, color='c')
axarr[0].set_title('Es')
axarr[1].bar(np.arange(len(np.arange(len(Pbs)))), Pbs, width, color='c')
axarr[1].set_title('P(snew)')
plt.show()
```





3 9.2 Mean-Field Annealing

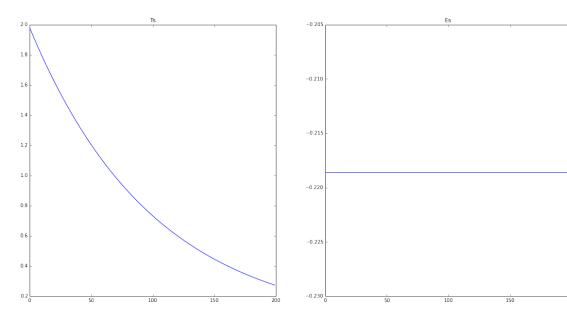
```
In [9]: #set initial state s randomly
        s = np.sign(np.random.rand(N)-.5)
        #w from above
In [10]: #The cost (energy) function remains the same
         def E(s, w):
             total = 0
             for i in range(len(s)):
                 for j in range(len(s)):
                     total = total + w[i,j]*s[i]*s[j]
             return -.5 * total
         \#The\ approximated\ probability\ of\ a\ state\ s\ is\ now\ given\ by
         def Q(s, e, b):
             total=0
             for j in range(len(s)):
                 total += e[j]*s[j]
             return (1.0/Z(s,w,b)) * np.exp(-b * total)
```

```
#is given as the sum over all possible configurations
         def Z(s, w, b):
             total = 0
             for i in range(len(s)):
                 total = total + np.exp( -b * E(s, w))
             return total
         def e(s, w):
             mye = np.zeros(len(s))
             for i in range(len(s)):
                 mye[i] = ei(s, w, i)
             return mye
         def ei(s, w, i):
             myei = 0
             for j in range(len(s)):
                 myei = myei + w[i,j]*s[j]
             return -1 * myei
In [11]: def MeanFieldAnnealing(s, w):
             # Initialization
             mys = s.copy()
             myw = w.copy()
             #\beta0 small enough, \tau > 1
             b0 = .5
             tau = 1.01
             #set tmax, \varepsilon small enough
             tmax = 200
             epsilon = .00001
             #number of times to repeat
             #try both M=1 and M=500
             M = 500
             # Optimization
             Ts = np.zeros(tmax)
             Es = np.zeros(tmax)
             bt = b0
             eold = np.zeros(len(s))
             enew = eold
             iteration = 0
             maxiterations = 500
             #for each iteration t = 0, \ldots, tmax
             for t in range(tmax):
                 eold = enew
                 enew = e(s,w)
                  #repeat the following until convergence, i.e., /enew - eold/ < \varepsilon
```

#the partition function Z guarantees P (s) to be a valid probability mass function

```
while np.absolute( np.min(enew - eold) ) > epsilon and iteration < maxiterations:
            #for i=1,...,N
            for i in range(N):
                \#compute\ mean-field
                enew[i] = ei(s, w, i)
                #update the state
                s[i] = np.tanh(-bt * enew[i])
            iteration = iteration + 1
        #if iteration < maxiterations:</pre>
             print("converged at", iteration)
        #increase \beta
        bt = tau * bt
        #For plotting
        Ts[t] = 1.0/bt
        Es[t] = E(mys, myw)
    return s, Ts, Es
snew, Ts, Es = MeanFieldAnnealing(s, w)
```

4 Plotting



In []: