STA663_Final_Project

April 30, 2015

1 1) Background

1.1 Problem:

If a researcher is working with longitudinal or panel data, he or she may want to detect latent groupings or clusters in these data. Specifically, for life-course approaches to research, you may want to identify latent trajectories of a behavior or outcome across time.

1.2 Outline of Algorithm:

The paper I have chosen describes a type of finite mixture algorithm, one for modeling patterns of an outcome of interest as clusters across time. The method is termed by the authors as Group-based Latent Trajectory Modeling, and was developed by Daniel Nagin in 1999.

The general specification for the group based latent trajectory models is provided in equations (1) and (2). Equation (1) describes the basic form of the finite mixture model—i.e., summing a finite number of latent groupings believed to compose the underlying population. Since group membership is not observed, the proportion of the j underlying population belonging to each of the latent trajectory groups π_j , must be estimated. This requires the aggregation of the J conditional likelihood functions, forming the unconditional probability of the data, Y_i (Nagin 2005).

$$P(Y_i) = \sum_{j=1}^{J} \pi_j P^j(Y_i) \tag{1}$$

Here, $P(Y_i)$ estimates the unconditional probability of seeing the trajectory of the outcome measure for individual i.

The key to the "group-based" approach is an underlying idea called "conditional independence." The likelihood function this produces is denoted in equation (2). For a give group j, conditional independence makes the assumption that the distribution of y_{it} is independent of the observed value of the outcome in prior periods, $y_{it-1}, y_{it-2},...$ This assumption helps to reduce the complexity of the model, and when combined with the EM algorithm, allows for identification of latent trajectory groupings despite the complex nature of the model and data.

$$L = \prod_{i=1}^{N} P(Y_i) \tag{2}$$

1.3 Readme for Python Module:

1.3.1 Group-Based Finite Mixture Models: A Latent Trajectory Approach

Code for course project use only at this time.

Final project for STA 663 * Takes time variable and continuous outcome * Generates latent trajectories (time-dependent clusters) of outcome behavior over time

Dependencies

Required

- Python (created using 2.7.x)
- Numpy
- Scipy
- Pymix (https://github.com/klahnakoski/pymix.git)

Recommended

- ggplot (for trajectory plotting function included in the repository)
- pandas (for trajectory plotting function included in the repository)
- STAN (for Bayesian alternative specification)
- pystan (for Bayesian alternative specification)
- rpy2 (for loading in test data from R)

2 2) Implementation

Note: I had started to build these distribution functions myself (see py scripts in GitHub repo) but then I discovered a python library with mixture distributions that could be combined and reshaped to estimate these models ('Pymix' package)!

```
In [78]: from __future__ import division
         import os
         import sys
         import glob
         from mixture import *
         import mixture
         import numpy as np
         import pandas as pd
         from ggplot import *
         import pandas as pd
         import pystan
         %matplotlib inline
         %precision 4
         %load_ext rpy2.ipython
         np.random.seed(34)
         # ---- #
         # Normal Mixture #
         # ---- #
         # Normal Distribution Functions
         # Generate Means of Normal Mixture
         def norm_means(data):
             cols = data.shape[1]
             return np.repeat(0.0,cols)
         # Generate Covariance Structure of Normal Mixture based on conditional independence assumption
         # (See Equation (2) in outline of algorithm)
         def norm_cov(data):
             m = data.shape[1]
             covs = np.repeat(1.0,m)
             return np.diag(covs)
```

Single Multivariate Normal distribution function

```
def norm_singdist(ng, data):
   m = data.shape[1]
   dist = mixture.MultiNormalDistribution(m,
        norm_means(data),norm_cov(data))
   return dist
# Generalize to multiple clusters of multivariate distribution function
def norm_multdist(ng, data):
   temp = list()
   for x in range(ng):
        temp.append(norm_singdist(2, data))
   return temp
# Data functions
def build_mix_dat(data):
   mixdat = mixture.DataSet()
   mixdat.fromArray(data)
   return mixdat
# Normal Mixture Model
def intialize_normal_model(ng, data):
   mod_ps = np.repeat(1.0/ng, ng)
   if ng ==2:
        n1,n2 = norm_multdist(ng, data)
       mix_ = mixture.MixtureModel(ng, mod_ps,[n1,n2])
   elif ng == 3:
        n1,n2,n3 = norm_multdist(ng, data)
        mix_ = mixture.MixtureModel(ng, mod_ps,[n1,n2,n3])
        n1,n2,n3,n4 = norm_multdist(ng, data)
        mix_ = mixture.MixtureModel(ng, mod_ps,[n1,n2,n3,n4])
   elif ng == 5:
        n1,n2,n3,n4,n5 = norm_multdist(ng, data)
        mix_ = mixture.MixtureModel(ng, mod_ps,[n1,n2,n3,n4,n5])
   elif ng == 6:
       n1,n2,n3,n4,n5,n6 = norm_multdist(ng, data)
        mix_ = mixture.MixtureModel(ng, mod_ps,[n1,n2,n3,n4,n5,n6])
   elif ng == 7:
        n1,n2,n3,n4,n5,n6,n7 = norm_multdist(ng, data)
       mix_ = mixture.MixtureModel(ng, mod_ps,[n1,n2,n3,n4,n5,n6,n7])
   elif ng == 8:
        n1,n2,n3,n4,n5,n6,n7,n8 = norm_multdist(ng, data)
        mix_ = mixture.MixtureModel(ng, mod_ps,[n1,n2,n3,n4,n5,n6,n7,n8])
   return mix_
# Run Normal Mixture Model
def normal_mixmod(ng, data, s=0):
   mix_mod = intialize_normal_model(ng, data)
   mix_dat = build_mix_dat(data)
   mix_mod.modelInitialization(mix_dat)
   mix_mod.EM(mix_dat, 40,.2, silent=s)
   mix_mod.classify(mix_dat, silent=s)
   return mix_mod
```

```
# Results
def GBFMM(ng,data, s=0):
    out = normal_mixmod(ng, data, s=s)
    traj = list()
    for i in range(ng):
        traj.append(out.components[i].distList[0].mu)
    return traj
# Trajectory Plotting Function
def plot_traj(outs):
    if len(outs) == 2:
        d = {'one' : outs[0],
        'two' : outs[1],
        'Time': range(1,len(outs[0])+1)}
    elif len(outs) == 3:
        d = {'one' : outs[0],
        'two' : outs[1],
        'three' : outs[2],
        'Time': range(1,len(outs[0])+1)}
    elif len(outs) == 4:
        d = {'one' : outs[0],
        'two' : outs[1],
        'three' : outs[2],
        'four' : outs[3],
        'Time': range(1,len(outs[0])+1)}
    elif len(outs) == 5:
        d = {'one' : outs[0],
        'two' : outs[1],
        'three' : outs[2],
        'four' : outs[3],
        'five' : outs[4],
        'Time': range(1,len(outs[0])+1)}
    elif len(outs) == 6:
        d = {'one' : outs[0],
        'two' : outs[1],
        'three' : outs[2],
        'four' : outs[3],
        'five' : outs[4],
        'six' : outs[5],
        'Time': range(1,len(outs[0])+1)}
    elif len(outs) == 7:
        d = {'one' : outs[0],
        'two' : outs[1],
        'three' : outs[2],
        'four' : outs[3],
        'five' : outs[4],
        'six' : outs[5],
        'seven' : outs[6],
        'Time': range(1,len(outs[0])+1)}
    elif len(outs) == 8:
        d = {'one' : outs[0],
        'two' : outs[1],
        'three' : outs[2],
        'four' : outs[3],
```

```
'five' : outs[4],
    'six' : outs[5],
    'seven' : outs[6],
    'eight' : outs[7],
    'Time': range(1,len(outs[0])+1)}
    dat = pd.DataFrame(d)
    dm = pd.melt(dat, id_vars=['Time'], var_name='Latent Class', value_name='Value')
    print ggplot(dm, aes('Time', 'Value', color='Latent Class')) + \
        geom_point(size=80) + \
        geom_line(size=3) + \
        theme_bw()
The rpy2.ipython extension is already loaded. To reload it, use:
    %reload_ext rpy2.ipython
```

3 3) Testing

```
In [25]: # ----- #
         # Tests for convergence given mixture distributions with given number of groupings
         # (these also test that the number of latent groups produced equal the number of groups specif
         # test_numgroups_3
         # ----- #
         dist3 = MultiNormalDistribution(3, [0.0, 1.0, 3.0],
                              [[1.0, 0.1, 0.14],
                              [0.2, 1.0, 0.2],
                              [0.3, 0.2, 1.0]
         data3 = dist3.sampleSet(100)
         assert len(GBFMM(ng=3, data=data3, s=1)) == 3
         # test_numgroups_4
         dist4 = MultiNormalDistribution(4, [0.0, 1.0, 3.0,2.0],
                             [[1.0, 0.1, 0.14, 0.2],
                              [0.2, 1.0, 0.2, 0.4],
                              [0.3, 0.2, 1.0, 0.3],
                              [0.3, 0.2, 0.3, 1.0])
         data4 = dist4.sampleSet(100)
         assert len(GBFMM(ng=4, data=data4, s=1)) == 4
         # test_numgroups_5
         dist5 = MultiNormalDistribution(5, [0.0, 1.0, 3.0,2.0,0.5],
                              [[1.0, 0.1, 0.14, 0.2, 0.1],
                              [0.2, 1.0, 0.2, 0.4, 0.35],
                              [0.3, 0.2, 1.0, 0.3, 0.05],
                              [0.3, 0.2, 0.3, 1.0, 0.32],
                               [0.1, 0.22, 0.08, .15, 1.0])
         data5 = dist5.sampleSet(100)
         assert len(GBFMM(ng=5, data=data5, s=1)) == 5
         # test_numgroups_6
         dist6 = MultiNormalDistribution(6, [0.0, 1.0, 3.0,2.0,0.5,1.5],
                              [[1.0, 0.1, 0.14, 0.2, 0.1, 0.6],
                              [0.2, 1.0, 0.2, 0.4, 0.35, 0.2],
                              [0.3, 0.2, 1.0, 0.3, 0.05, 0.0],
```

[0.3, 0.2, 0.3, 1.0, 0.32,0.1],

```
[0.1, 0.22, 0.08, 0.15, 1.0, 0.25],
                      [0.2, 0.1, 0.2, 0.4, 0.35, 1.0]])
data6 = dist6.sampleSet(1000)
assert len(GBFMM(ng=6, data=data6, s=1)) == 6
# test_numgroups_7
dist7 = MultiNormalDistribution(7, [0.0, 1.0, 3.0, 2.0, 0.5, 1.5, -1.0],
                    [[1.0, 0.1, 0.14, 0.2, 0.1, 0.6, 0.25],
                      [0.2, 1.0, 0.2, 0.4, 0.35, 0.2, 0.2],
                      [0.3, 0.2, 1.0, 0.3, 0.05, 0.0, 0.3],
                     [0.3, 0.2, 0.3, 1.0, 0.32, 0.1, 0.2],
                      [0.1, 0.22, 0.08, 0.15, 1.0, 0.25, 0.1],
                      [0.2, 0.1, 0.2, 0.4, 0.35,1.0, 0.0],
                      [0.1, 0.22, 0.08, 0.15, 0.2, 0.15, 1.0]])
data7 = dist7.sampleSet(1000)
assert len(GBFMM(ng=7, data=data7, s=1)) == 7
# test_numgroups_8
dist8 = MultiNormalDistribution(8, [0.0, 1.0, 3.0,2.0,0.5,1.5,-1.0,1.0],
                    [[1.0, 0.1, 0.14, 0.2, 0.1, 0.6, 0.25, 0.2],
                     [0.2, 1.0, 0.2, 0.4, 0.35, 0.2, 0.2, 0.1],
                     [0.3, 0.2, 1.0, 0.3, 0.05, 0.0, 0.3, 0.2],
                     [0.3, 0.2, 0.3, 1.0, 0.32, 0.1, 0.2, 0.35],
                     [0.1, 0.22, 0.08, 0.15, 1.0, 0.25, 0.1, 0.05],
                     [0.2, 0.1, 0.2, 0.4, 0.35, 1.0, 0.0, 0.1],
                     [0.1, 0.22, 0.08, 0.15, 0.2, 0.15, 1.0, 0.3],
                     [0.15, 0.2, 0.1, 0.2, 0.2, 0.35, 0.0, 1.0]])
data8 = dist8.sampleSet(1000)
assert len(GBFMM(ng=8, data=data8, s=1)) == 8
# ----- #
# Tests for convergence for small N
# ----- #
dist_small_1 = MultiNormalDistribution(3, [0.0, 1.0, 3.0],
                    [[1.0, 0.1, 0.14],
                     [0.2, 1.0, 0.2],
                     [0.3, 0.2, 1.0]
data_small_1 = dist_small_1.sampleSet(30)
assert len(GBFMM(ng=3, data=data_small_1, s=1)) == 3
dist_small_2 = MultiNormalDistribution(4, [0.0, 1.0, 3.0,2.0],
                    [[1.0, 0.1, 0.14, 0.2],
                     [0.2, 1.0, 0.2, 0.4],
                     [0.3, 0.2, 1.0, 0.3],
                     [0.3, 0.2, 0.3, 1.0])
data_small_2 = dist_small_2.sampleSet(50)
assert len(GBFMM(ng=4, data=data_small_2, s=1)) == 4
# ----- #
# Tests for convergence for large N
# ---- #
dist_large_1 = MultiNormalDistribution(3, [0.0, 1.0, 3.0],
                    [[1.0, 0.1, 0.14],
                     [0.2, 1.0, 0.2],
```

```
[0.3, 0.2, 1.0]
         data_large_1 = dist_large_1.sampleSet(2000)
         assert len(GBFMM(ng=3, data=data_large_1, s=1)) == 3
         dist_large_2 = MultiNormalDistribution(4, [0.0, 1.0, 3.0,2.0],
                             [[1.0, 0.1, 0.14, 0.2],
                              [0.2, 1.0, 0.5, 0.4],
                              [0.3, 0.2, 1.0,0.3],
                              [0.0, 0.4, 0.2, 1.0]
         data_large_2 = dist_large_2.sampleSet(3000)
         assert len(GBFMM(ng=4, data=data_large_2, s=1)) == 4
In [69]: # ----- #
         # Test Trajectory Plotting Function & Visual Check of Trajectories
         # ----- #
         d1 = MultiNormalDistribution(4, [0.0, 1.0, 3.0,2.0],
                             [[1.0, 0.1, 0.14, 0.2],
                              [0.8, 1.0, 0.2, 0.4],
                              [0.5, 0.2, 1.0, 0.3],
                              [0.1, 0.3, 0.3, 1.0]])
         d2 = MultiNormalDistribution(4, [1.0,0.2, 2.0,4.5],
                             [[1.0, 0.3, 0.5, 0.0],
                              [0.5, 1.0, 0.2, 0.5],
                              [0.3, 0.4, 1.0, 0.5],
                             [0.3, 0.4, 0.1, 1.0]])
         d3 = MultiNormalDistribution(4, [3.0,2.0, 1.0,0.0],
                             [[1.0, 0.3, 0.5, 0.0],
                              [0.5, 1.0, 0.2, 0.5],
                              [0.3, 0.4, 1.0, 0.5],
                             [0.3, 0.4, 0.1, 1.0])
         data1 = d1.sampleSet(300)
         data2 = d2.sampleSet(100)
         data3 = d3.sampleSet(100)
         # print np.array(data1)
         data3= np.concatenate([data1, data2, data3], axis=0)
         outs_1 = GBFMM(ng=4, data=data3, s=0)
         plot_traj(outs_1)
Step 1: log likelihood: -3312.4960746
                                         (diff=-3311.4960746)
Step 2: log likelihood: -3298.60641088
                                          (diff=13.8896637209)
Step 3: log likelihood: -3263.23596942
                                          (diff=35.3704414607)
Step 4: log likelihood: -3187.20997306
                                          (diff=76.0259963601)
Step 5: log likelihood: -3094.92279202
                                          (diff=92.2871810369)
Step 6: log likelihood: -3027.36520871
                                          (diff=67.5575833093)
Step 7: log likelihood: -2996.13875147
                                          (diff=31.2264572371)
Step 8: log likelihood: -2980.69605671
                                          (diff=15.4426947604)
Step 9: log likelihood: -2969.45658442
                                          (diff=11.2394722896)
Step 10: log likelihood: -2958.9513708
                                          (diff=10.505213623)
Step 11: log likelihood: -2949.23252172
                                          (diff=9.718849085)
Step 12: log likelihood: -2941.73097139
                                           (diff=7.50155032951)
                                           (diff=4.37318547048)
Step 13: log likelihood: -2937.35778592
```

```
Step 14: log likelihood: -2935.01655564
                                          (diff=2.34123027403)
Step 15: log likelihood: -2933.56541109
                                          (diff=1.45114455759)
Step 16: log likelihood: -2932.50329834
                                          (diff=1.06211274104)
Step 17: log likelihood: -2931.6131419
                                          (diff=0.890156442188)
Step 18: log likelihood: -2930.84316869
                                          (diff=0.769973207313)
Step 19: log likelihood: -2930.20869633
                                          (diff=0.634472362919)
Step 20: log likelihood: -2929.71403334
                                          (diff=0.494662993972)
Step 21: log likelihood: -2929.30273173
                                          (diff=0.411301606659)
Step 22: log likelihood: -2928.83515997
                                          (diff=0.467571759812)
Step 23: log likelihood: -2928.14892766
                                          (diff=0.686232310017)
Step 24: log likelihood: -2927.26750858
                                          (diff=0.881419084902)
Step 25: log likelihood: -2926.17741575
                                          (diff=1.09009283025)
Step 26: log likelihood: -2924.90417299
                                          (diff=1.27324275228)
Step 27: log likelihood: -2923.94720685
                                          (diff=0.95696614857)
Step 28: log likelihood: -2923.4871062
                                          (diff=0.460100644041)
Step 28: log likelihood: -2923.28954442
                                          (diff=0.197561781967)
Convergence reached with log_p -2923.28954442 after 28 steps.
classify loglikelihood: -2923.28954442.
```

** Clustering **

Cluster 0, size 30

[60, 75, 148, 158, 160, 200, 217, 407, 408, 409, 412, 414, 422, 424, 433, 438, 448, 454, 460, 464, 479,

Cluster 1, size 76

[400, 401, 402, 403, 404, 405, 406, 410, 411, 413, 415, 416, 417, 418, 419, 420, 421, 423, 425, 426, 42

Cluster 2 , size 99

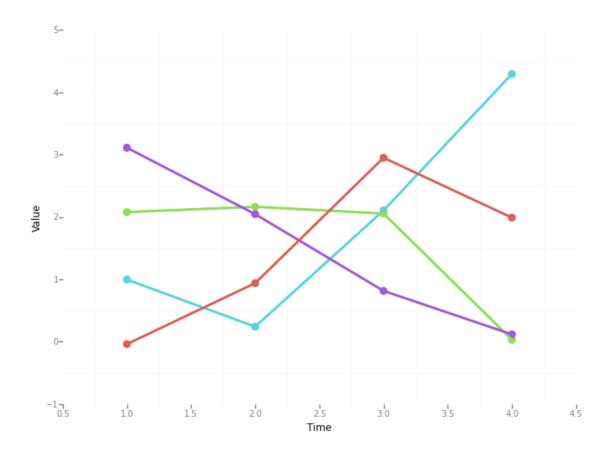
[135, 141, 186, 300, 301, 302, 303, 304, 305, 306, 307, 308, 309, 310, 311, 313, 314, 315, 316, 317, 31

Cluster 3, size 295

[0, 1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14, 15, 16, 17, 18, 19, 20, 21, 22, 23, 24, 25, 26, 27,

Unassigend due to entropy cutoff:

[]



<ggplot: (303148237)>

4 4) Application and Comparison

4.0.2 Data

To apply this algorithm to a real problem on real data, I examine the Adjudicated Toronto Youth Data. The data include the number of criminal court appearances by 378 youth in Toronto. Each individual has 5 years of data, where each row contains the number of unique court contacts per year for an individual from the age of 18 to 23.

4.0.3 Comparison

I compare the EM algorithm approach to a Bayesian finite mixture written in the probabilistic programming language STAN, and called by the python module pystan. This approach using MCMC instead of the MLE approach based on the EM algorithm.

```
Step 5: log likelihood: -5721.56812432
                                         (diff=4.02643363104)
Step 6: log likelihood: -5709.27463866
                                         (diff=12.2934856551)
Step 7: log likelihood: -5677.63503708
                                         (diff=31.6396015824)
Step 8: log likelihood: -5577.26677048
                                         (diff=100.368266603)
Step 9: log likelihood: -5360.78622798
                                         (diff=216.480542493)
Step 10: log likelihood: -5161.11428695
                                          (diff=199.671941031)
Step 11: log likelihood: -4949.69647743
                                          (diff=211.417809524)
Step 12: log likelihood: -4864.92774278
                                          (diff=84.7687346443)
Step 13: log likelihood: -4837.29260321
                                          (diff=27.63513957)
Step 14: log likelihood: -4831.85916386
                                          (diff=5.43343935681)
Step 15: log likelihood: -4830.13973582
                                          (diff=1.71942804021)
Step 16: log likelihood: -4829.07981934
                                          (diff=1.05991648075)
Step 17: log likelihood: -4828.08993193
                                          (diff=0.989887406076)
Step 18: log likelihood: -4827.06546163
                                          (diff=1.02447029583)
Step 19: log likelihood: -4826.4053168
                                         (diff=0.660144833062)
Step 19: log likelihood: -4826.21801076
                                          (diff=0.187306036481)
Convergence reached with log_p -4826.21801076 after 19 steps.
classify loglikelihood: -4826.21801076.
** Clustering **
Cluster 0, size 134
[3, 4, 6, 7, 14, 19, 25, 32, 35, 37, 40, 51, 53, 55, 56, 58, 61, 64, 65, 66, 68, 72, 75, 78, 80, 85, 88
Cluster 1, size 244
[0, 1, 2, 5, 8, 9, 10, 11, 12, 13, 15, 16, 17, 18, 20, 21, 22, 23, 24, 26, 27, 28, 29, 30, 31, 33, 34,
Unassigend due to entropy cutoff:
In [77]: plot_traj(outs)
```

(diff=-5873.69620147)

(diff=92.408446164)

(diff=46.9149555604)

(diff=9.77824179596)

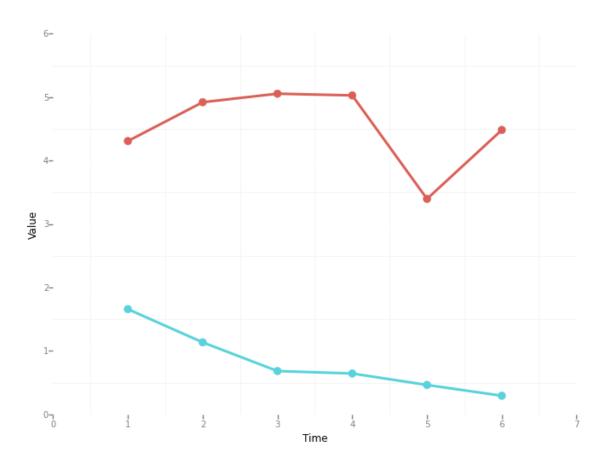
In [76]: outs = GBFMM(2,TO1adj)

Step 1: log likelihood: -5874.69620147

Step 2: log likelihood: -5782.28775531

Step 3: log likelihood: -5735.37279974

Step 4: log likelihood: -5725.59455795



<ggplot: (308464657)>

4.0.4 Results

The EM based finite mixture algorithm has identified two clustered trajectories. The red shows a cluster of youth who are coming into contact with criminal courts frequently over this five year period. The blue line shows a low risk trajectory.

Now let's try the MCMC approach:

```
In [97]: model_code = """
         data {
           int<lower=1> K;
                                                                              # K components
           int<lower=1> N;
                                                                              # N observations
           real y[N];
                                                                              # variable of interest
         parameters {
           simplex[K] theta;
                                                                              # mixing proportions
           simplex[K] mu_prop;
           real mu_loc;
           real<lower=0> mu_scale;
           real<lower=0> sigma[K];
                                                                              # sds of the components
         transformed parameters {
           ordered[K] mu;
```

```
model {
           // prior
           mu_loc ~ cauchy(0,5);
           mu_scale ~ cauchy(0,5);
           sigma ~ cauchy(0,5);
           // likelihood
             real ps[K];
             vector[K] log_theta;
             log_theta <- log(theta);</pre>
             for (n in 1:N) {
               for (k in 1:K) {
                 ps[k] <- log_theta[k]</pre>
                          + normal_log(y[n],mu[k],sigma[k]);
               increment_log_prob(log_sum_exp(ps));
           }
         }
         0.00
         dat = {
             'K': 2,
             'N': len(np.array(T01stan)),
             'y': np.array(T01stan),
         fit = pystan.stan(model_code=model_code, data=dat, iter=1000, chains=2)
/Users/Josh/anaconda/lib/python2.7/multiprocessing/queues.py:390: UserWarning: Pickling fit objects is
The relevant StanModel instance must be pickled along with this fit object.
When unpickling the StanModel must be unpickled first.
  return send(obj)
/Users/Josh/anaconda/lib/python2.7/multiprocessing/queues.py:390: UserWarning: Pickling fit objects is
The relevant StanModel instance must be pickled along with this fit object.
When unpickling the StanModel must be unpickled first.
  return send(obj)
In [98]: print fit
Inference for Stan model: anon_model_0ea983c07018240174b20728e647dd40.
2 chains, each with iter=1000; warmup=500; thin=1;
post-warmup draws per chain=500, total post-warmup draws=1000.
                                   2.5%
                                            25%
                                                          75% 97.5% n_eff
             mean se_mean
                              sd
                                                   50%
                                                                              Rhat
theta[0]
             0.35 1.9e-3
                            0.03
                                   0.29
                                           0.33
                                                  0.35
                                                         0.37
                                                                0.41 257.0
                                                                                1.0
theta[1]
             0.65 1.9e-3
                            0.03
                                   0.59
                                           0.63
                                                  0.65
                                                         0.67
                                                                0.71 257.0
                                                                                1.0
             0.38
                     0.02
                             0.2
                                   0.02
                                          0.21
                                                  0.38
                                                         0.54
                                                                0.72
                                                                       79.0
                                                                                1.0
mu_prop[0]
             0.62
                     0.02
                                  0.28
                                          0.46
                                                         0.79
                                                                0.98
                                                                       79.0
mu_prop[1]
                             0.2
                                                  0.62
                                                                                1.0
mu\_loc
             0.2
                     0.18
                            1.53 -3.61 -0.67
                                                  0.61
                                                         1.43
                                                                1.98
                                                                       74.0
                                                                                1.0
```

means of the components

mu <- mu_loc + mu_scale * cumulative_sum(mu_prop);</pre>

}

 mu_scale

sigma[0]

sigma[1]

4.08

0.18

0.25 1.4e-3

0.44 1.7e-3

1.53

0.02

0.03 0.39

2.29

0.2

2.86

0.23

0.42

3.68

0.24

0.44

4.95

0.26

7.86

0.45 0.49 244.0

0.29 268.0

74.0

1.0

1.0

1.0

```
mu[0]
             2.02
                   1.5e-3
                             0.03
                                    1.97
                                            2.01
                                                   2.02
                                                           2.04
                                                                  2.08
                                                                        306.0
                                                                                  1.0
mu[1]
                                                           4.3
             4.28
                   1.9e-3
                             0.03
                                     4.2
                                            4.25
                                                   4.28
                                                                  4.34
                                                                       319.0
                                                                                 1.0
           -283.6
                     0.13
                             1.75 -288.0 -284.5 -283.2 -282.3 -281.2 182.0
lp__
                                                                                 1.0
```

Samples were drawn using NUTS(diag_e) at Thu Apr 30 15:29:21 2015. For each parameter, n_eff is a crude measure of effective sample size, and Rhat is the potential scale reduction factor on split chains (at convergence, Rhat=1).

4.0.5 Results

While similar inferences can be drawn from this MCMC method, and I would argue in many cases the Bayesian results are more robust, convergence time is much slower. For mixture model problems where the underlying structure is not incredibly complex, the EM algorithm employed by the Group-Based Finite Mixture Model may be more practical.