# Deep Autoencoding Gaussian Mixture Model for Unsupervised Anomaly Detection

```
import numpy as np
import pandas as pd
import torch
from data_loader import *
from main import *
from tqdm import tqdm
```

### KDD Cup 1999 Data (10% subset)

This is the data set used for The Third International Knowledge Discovery and Data Mining Tools Competition, which was held in conjunction with KDD-99 The Fifth International Conference on Knowledge Discovery and Data Mining. The competition task was to build a network intrusion detector, a predictive model capable of distinguishing between "bad" connections, called intrusions or attacks, and "good" normal connections. This database contains a standard set of data to be audited, which includes a wide variety of intrusions simulated in a military network environment.

```
data = pd.read_csv("kddcup.data_10_percent_corrected", header=None,names=['duration', 'protocol_type', 'service', 'flag',
```

### **Pre-processing**

Following the paper, since the "normal" only comprises of approximately 20% of the entries, the "normal" data were considered as anomalies instead.

```
data.loc[data["type"] != "normal.", 'type'] = 0
data.loc[data["type"] == "normal.", 'type'] = 1
```

Next, the categorical variables are converted to a one hot encoding representation. My implementation is a bit different from the original paper in this aspect. Since I amonly using the 10% subset to generate the columns, I get 118 features instead of 120 as reported in the paper.

```
one_hot_protocol = pd.get_dummies(data["protocol_type"])
one_hot_service = pd.get_dummies(data["service"])
one_hot_flag = pd.get_dummies(data["flag"])

data = data.drop("protocol_type",axis=1)
data = data.drop("service",axis=1)
data = data.drop("flag",axis=1)

data = pd.concat([one_hot_protocol, one_hot_service, one_hot_flag, data],axis=1)
data.head()
```

	icmp	tcp	udp	IRC	X11	Z39_50	auth	bgp	courier	csnet_ns	
0	0	1	0	0	0	0	0	0	0	0	
1	0	1	0	0	0	0	0	0	0	0	

2	0	1	0	0	0	0	0	0	0	0	
3	0	1	0	0	0	0	0	0	0	0	
4	0	1	0	0	0	0	0	0	0	0	
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#### 5 rows × 119 columns

```
proportions = data["type"].value_counts()
print(proportions)
print("Anomaly Percentage",proportions[1] / proportions.sum())
```

```
0 396743
1 97278
Name: type, dtype: int64
Anomaly Percentage 0.19691065764410826
```

```
#proportions_alfa = data["type"].value_counts(normalize=True)
#print(proportions_alfa)
```

Normalize all the numeric variables.

I saved the preprocessed data into a numpy file format and load it using the pytorch data loader.

```
np.savez_compressed("kdd_cup",kdd=data.as_matrix())

C:\Users\cncluser\Anaconda3\lib\site-packages\ipykernel_launcher.py:1: FutureWarning: Method .as_matrix will be removed in """Entry point for launching an IPython kernel.
```

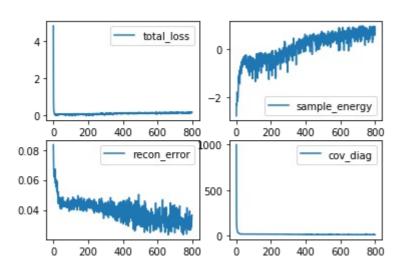
I initially implemented this to be ran in the command line and use argparse to get the hyperparameters. To make it runnable in a jupyter notebook, I had to create a dummy class for the hyperparameters.

```
class hyperparams():
    def __init__(self, config):
        self.__dict__.update(**config)
defaults = {
```

```
'lr' : 1e-4,
    'num_epochs' : 200,
    'batch_size' : 1024,
    'gmm_k' : 4,
    'lambda_energy' : 0.1,
    'lambda_cov_diag' : 0.005,
    'pretrained_model' : None,
    'mode' : 'train',
    'use_tensorboard' : False,
    'data_path' : 'kdd_cup.npz',
    'log_path' : './dagmm/logs',
    'model_save_path' : './dagmm/models',
    'sample_path' : './dagmm/samples',
    'test_sample_path' : './dagmm/test_samples',
    'result_path' : './dagmm/results',
    'log_step' : 194//4,
    'sample_step' : 194,
    'model_save_step' : 194,
}
```

```
solver = main(hyperparams(defaults))
accuracy, precision, recall, f_score = solver.test()
```

```
Elapsed 1:30:00.990700/0:00:00.417624 -- 0:00:00.417624 , Epoch [200/200], Iter [192/194], lr 0.0001, total_loss: 0.1735,
```



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```
[ 3.2677, -0.4000, 0.7402]]])
100%
                                                                           194/194 [00:10<00:00, 18.23it/s]
tensor([6.6445e-04, 2.1110e-03, 8.6473e-04, 9.9635e-01],
      grad_fn=<DivBackward0>)
tensor([[-0.2979, 2.3099, 0.0913],
       [-0.3379, 2.8321, 0.1679],
       [-0.2949, 2.1895, 0.0803],
       [ 2.0863, 0.3287, 0.4579]], grad_fn=<DivBackward0>)
tensor([[[10.4720, -0.1645, 2.2967],
        [-0.1645, 5.2059, 0.2242],
[ 2.2967, 0.2242, 0.5920]],
       [[ 7.5691, 0.5594, 1.5819],
        [ 0.5594, 5.3244, 0.2878],
[ 1.5819, 0.2878, 0.4418]],
       [[10.9752, -0.3034, 2.4145],
        [-0.3034, 5.0558, 0.1932],
[ 2.4145, 0.1932, 0.6160]],
       [[14.5482, -1.7952, 3.2824],
        [-1.7952, 0.3579, -0.3949],
        [ 3.2824, -0.3949, 0.7454]]], grad_fn=<DivBackward0>)
Threshold : 6.946145534515381
Accuracy : 0.9746, Precision : 0.9681, Recall : 0.9542, F-score : 0.9611
```

I copy pasted the testing code here in the notebook so we could play around the results.

Incrementally compute for the GMM parameters across all training data for a better estimate

```
solver.data_loader.dataset.mode="train"
solver.dagmm.eval()
N = 0
mu sum = 0
cov_sum = 0
gamma sum = 0
for it, (input_data, labels) in enumerate(solver.data_loader):
   input_data = solver.to_var(input_data)
   enc, dec, z, gamma = solver.dagmm(input_data)
   phi, mu, cov = solver.dagmm.compute_gmm_params(z, gamma)
   batch_gamma_sum = torch.sum(gamma, dim=0)
   gamma_sum += batch_gamma_sum
   mu_sum += mu * batch_gamma_sum.unsqueeze(-1) # keep sums of the numerator only
   N += input_data.size(0)
train_phi = gamma_sum / N
train_mu = mu_sum / gamma_sum.unsqueeze(-1)
train_cov = cov_sum / gamma_sum.unsqueeze(-1).unsqueeze(-1)
print("N:",N)
```

```
print("phi :\n",train_phi)
print("mu :\n",train_mu)
print("cov :\n",train_cov)
```

```
N: 198371
phi :
tensor([6.6445e-04, 2.1110e-03, 8.6473e-04, 9.9636e-01],
      grad_fn=<DivBackward0>)
tensor([[-0.2979, 2.3099, 0.0913],
       [-0.3379, 2.8321, 0.1679],
       [-0.2949, 2.1895, 0.0803],
       [ 2.0863, 0.3287, 0.4579]], grad_fn=<DivBackward0>)
cov:
 tensor([[[10.4748, -0.1583, 2.2963],
       [-0.1583, 5.2358, 0.2256],
       [ 2.2963, 0.2256, 0.5921]],
       [[ 7.5787, 0.5723, 1.5813],
        [ 0.5723, 5.3620, 0.2891],
        [ 1.5813, 0.2891, 0.4420]],
       [[10.9772, -0.2983, 2.4141],
        [-0.2983, 5.0827, 0.1944],
        [ 2.4141, 0.1944, 0.6160]],
       [[14.5466, -1.7950, 3.2821],
        [-1.7950, 0.3579, -0.3949],
        [ 3.2821, -0.3949, 0.7453]]], grad_fn=<DivBackward0>)
```

```
train_energy = []
train_labels = []
train_z = []
for it, (input_data, labels) in enumerate(solver.data_loader):
    input_data = solver.to_var(input_data)
    enc, dec, z, gamma = solver.dagmm(input_data)
    sample_energy, cov_diag = solver.dagmm.compute_energy(z, phi=train_phi, mu=train_mu, cov=train_cov, size_average=False
    train_energy.append(sample_energy.data.cpu().numpy())
    train_z.append(z.data.cpu().numpy())
    train_labels.append(labels.numpy())

train_labels.append(labels.numpy())

train_z = np.concatenate(train_energy,axis=0)
train_z = np.concatenate(train_labels,axis=0)
```

### Compute the energy of every sample in the test data

```
solver.data_loader.dataset.mode="test"
test_energy = []
test_labels = []
test_z = []
for it, (input_data, labels) in enumerate(solver.data_loader):
    input_data = solver.to_var(input_data)
    enc, dec, z, gamma = solver.dagmm(input_data)
    sample_energy, cov_diag = solver.dagmm.compute_energy(z, size_average=False)
    test_energy.append(sample_energy.data.cpu().numpy())
    test_z.append(z.data.cpu().numpy())
    test_labels.append(labels.numpy())

test_energy = np.concatenate(test_energy,axis=0)
test_z = np.concatenate(test_z,axis=0)
```

```
test_labels = np.concatenate(test_labels,axis=0)

combined_energy = np.concatenate([train_energy, test_energy], axis=0)
combined_z = np.concatenate([train_z, test_z], axis=0)
combined_labels = np.concatenate([train_labels, test_labels], axis=0)
```

Compute for the threshold energy. Following the paper I just get the highest 20% and treat it as an anomaly. That corresponds to setting the threshold at the 80th percentile.

```
thresh = np.percentile(combined_energy, 100 - 20)
print("Threshold:", thresh)

Threshold: 5.002402305603027

pred = (test_energy>thresh).astype(int)
gt = test_labels.astype(int)

from sklearn.metrics import precision_recall_fscore_support as prf, accuracy_score

accuracy = accuracy_score(gt,pred)
precision, recall, f_score, support = prf(gt, pred, average='binary')

print("Accuracy: {:0.4f}, Precision: {:0.4f}, Recall: {:0.4f}, F-score: {:0.4f}".format(accuracy,precision, recall, f_score)

Accuracy: 0.9746, Precision: 0.9687, Recall: 0.9537, F-score: 0.9611
```

## Visualizing the z space

It's a little different from the paper's figure but I assume that's because of the small changes in my implementation.

```
from mpl_toolkits.mplot3d import Axes3D
import matplotlib.pyplot as plt
%matplotlib notebook
fig = plt.figure()
ax = fig.add_subplot(111, projection='3d')
ax.scatter(test_z[:,1],test_z[:,0], test_z[:,2], c=test_labels.astype(int))
ax.set_xlabel('Encoded')
ax.set_ylabel('Euclidean')
ax.set_zlabel('Cosine')
plt.show()
```

Python.core.display.Javascript object									