## Kagle\_Submission

December 13, 2020

#### 1 Mechanisms of Action (MoA) Prediction

Predicting multiple targets of the Mechanism of Action (MoA) response(s) of different samples (sig\_id), given various inputs such as gene expression data and cell viability data.

# 1.1 Some of the important terms used in the headings of the tables are presented here:

```
g - : signifies gene expression data
    c - : signifies cell expression data
    cp_type : indicates samples treated with a compound (cp_vehicle) or with a control perturbation
    NOTE: (samples with control perturbations don't have MoAs)
    cp_time - treatment duration (24,48,72) Hours
    cp_dose - Dosage - HIGH or LOW
[1]: # Importing the multi label stratified k-fold
    # cross validator
    import sys
    sys.path.append('../input/iterative-stratification/
     from iterstrat.ml_stratifiers import MultilabelStratifiedKFold
    # Initial random imports
    import random
    import os
    import copy
    import warnings
    # warnings.filterwarnings('ignore')
    # Importing numpy
    import numpy as np
    # Importing pandas
    import pandas as pd
    # Importing matplotlib
    import matplotlib.pyplot as plt
```

```
# Importing seaborn
import seaborn as sns

# Importing sklearn
from sklearn import preprocessing
from sklearn.metrics import log_loss
from sklearn.preprocessing import StandardScaler
from sklearn.decomposition import PCA
from sklearn.feature_selection import VarianceThreshold

# Importing pytorch
import torch
import torch.nn as nn
import torch.nn.functional as F
import torch.optim as optim
```

#### 2 Using GPU if available

```
[2]: # using GPU if available
if torch.cuda.is_available():
    device_code = 'cuda'
else:
    device_code = 'cpu'
```

```
[3]: # setting the seed, so that every time the seed is started from the same number

def set_seed_characteristics(seed=55):
    # Setting a random seed value

    random.seed(seed)

# for guaranteering the reproducability of numbers by setting seed for NumPy

np.random.seed(seed)

# for setting the seed for cuda or cpu

torch.manual_seed(seed)

# To ensure that Pytorch doesnt just switch to the fastest possible_ualgorithm but
    # ensures that it selects a deterministic algorithm

torch.backends.cudnn.deterministic = True
```

#### 3 Reading the CSV Files

```
[4]: training_features = pd.read_csv('input/lish-moa/train_features.csv')
     # Reading the head rows and columns of train features
    training_features_head = training_features.head()
    training_targets_scored = pd.read_csv('input/lish-moa/train_targets_scored.csv')
     # Reading the head rows and columns of train targets scored
    training_targets_scored_head = training_targets_scored.head()
    testing features = pd.read csv('input/lish-moa/test features.csv')
     # Reading the head rows and columns of train targets non-scored
    testing_features_head = testing_features.head()
[5]: # Printing the head - training features
    training_features_head
[5]:
             sig_id cp_type cp_time cp_dose
                                                 g-0
                                                         g-1
                                                                 g-2
                                                                         g-3 \
    0 id_000644bb2 trt_cp
                                  24
                                          D1 1.0620 0.5577 -0.2479 -0.6208
    1 id_000779bfc trt_cp
                                          D1 0.0743 0.4087 0.2991 0.0604
                                  72
    2 id_000a6266a trt_cp
                                  48
                                          D1 0.6280 0.5817 1.5540 -0.0764
    3 id_0015fd391 trt_cp
                                  48
                                          D1 -0.5138 -0.2491 -0.2656 0.5288
    4 id_001626bd3 trt_cp
                                  72
                                          D2 -0.3254 -0.4009 0.9700 0.6919
          g-4
                  g-5 ...
                            c-90
                                    c-91
                                            c-92
                                                    c-93
                                                            c-94
                                                                    c-95
    0 -0.1944 -1.0120 ... 0.2862 0.2584 0.8076 0.5523 -0.1912 0.6584
    1 1.0190 0.5207 ... -0.4265 0.7543
                                          0.4708 0.0230 0.2957 0.4899
    2 -0.0323 1.2390 ... -0.7250 -0.6297 0.6103 0.0223 -1.3240 -0.3174
    3 4.0620 -0.8095 ... -2.0990 -0.6441 -5.6300 -1.3780 -0.8632 -1.2880
    4 1.4180 -0.8244 ... 0.0042 0.0048 0.6670 1.0690 0.5523 -0.3031
         c-96
                         c-98
                 c-97
                                 c-99
    0 -0.3981 0.2139 0.3801 0.4176
    1 0.1522 0.1241 0.6077 0.7371
    2 -0.6417 -0.2187 -1.4080 0.6931
    3 -1.6210 -0.8784 -0.3876 -0.8154
    4 0.1094 0.2885 -0.3786 0.7125
    [5 rows x 876 columns]
[6]: # Printing the head - train targets scored
    training_targets_scored_head
[6]:
             sig_id 5-alpha_reductase_inhibitor 11-beta-hsd1_inhibitor
    0 id_000644bb2
    1 id_000779bfc
                                               0
                                                                       0
                                               0
                                                                       0
    2 id_000a6266a
```

```
id_0015fd391
                                                0
                                                                          0
4 id_001626bd3
                                                0
                                                                          0
   acat_inhibitor
                     acetylcholine_receptor_agonist
0
                 0
                                                     0
1
2
                 0
                                                     0
3
                 0
                                                     0
   {\tt acetylcholine\_receptor\_antagonist} \quad {\tt acetylcholinesterase\_inhibitor}
0
                                       0
                                                                          0
1
                                       0
2
                                                                          0
3
                                       0
                                                                          0
4
                                       0
                                                                          0
   adenosine_receptor_agonist
                                 adenosine_receptor_antagonist
0
                               0
                                                                 0
1
2
                               0
                                                                 0
3
                               0
                                                                 0
4
                               0
   adenylyl_cyclase_activator
                                  ... tropomyosin_receptor_kinase_inhibitor
0
                                                                              0
                                                                              0
1
                               0
2
                               0
                                                                              0
3
                               0
                                                                              0
                               0
                                                                              0
                                     tubulin_inhibitor
   trpv_agonist trpv_antagonist
0
               0
                                  0
                                                       0
1
2
               0
                                  0
                                                       0
3
               0
                                                       0
   tyrosine_kinase_inhibitor
                                ubiquitin_specific_protease_inhibitor
0
                                                                         0
                              0
                              0
                                                                         0
1
                              0
                                                                         0
2
3
                              0
                                                                         0
   vegfr_inhibitor vitamin_b vitamin_d_receptor_agonist
                                                                wnt_inhibitor
0
```

```
      1
      0
      0
      0
      0

      2
      0
      0
      0
      0

      3
      0
      0
      0
      0

      4
      0
      0
      0
      0
```

[5 rows x 207 columns]

```
[7]: # Printing the head - test features
testing_features_head

[7]: sig_id cp_type cp_time cp_dose g-0 g-1 g-2 g-3 \
0 id_0004d9e33 trt_cp 24 D1 -0.5458 0.1306 -0.5135 0.4408
```

```
trt_cp
1 id_001897cda
                                  72
                                          D1 -0.1829
                                                     0.2320 1.2080 -0.4522
                                  24
                                          D1 0.1852 -0.1404 -0.3911 0.1310
2 id_002429b5b
               ctl_vehicle
3 id 00276f245
                                  24
                                         D2 0.4828 0.1955 0.3825 0.4244
                     trt_cp
4 id_0027f1083
                                          D1 -0.3979 -1.2680 1.9130 0.2057
                     trt_cp
                                  48
                       c-90
                               c-91
                                       c-92
                                               c-93
                                                      c-94
                                                              c-95
     g-4
             g-5 ...
0 1.5500 -0.1644 ... 0.0981 0.7978 -0.1430 -0.2067 -0.2303 -0.1193
1 -0.3652 -0.3319 ... -0.1190 -0.1852 -1.0310 -1.3670 -0.3690 -0.5382
2 -1.4380 0.2455 ... -0.2261 0.3370 -1.3840 0.8604 -1.9530 -1.0140
3 -0.5855 -1.2020 ... 0.1260 0.1570 -0.1784 -1.1200 -0.4325 -0.9005
4 -0.5864 -0.0166 ... 0.4965 0.7578 -0.1580 1.0510 0.5742 1.0900
     c-96
            c-97
                    c-98
                            c-99
0 0.0210 -0.0502 0.1510 -0.7750
1 0.0359 -0.4764 -1.3810 -0.7300
2 0.8662 1.0160 0.4924 -0.1942
3 0.8131 -0.1305 0.5645 -0.5809
4 -0.2962 -0.5313 0.9931 1.8380
```

[5 rows x 876 columns]

#### 4 Dataset classes, training and testing functions

```
[8]: # Pytorch data loader implementation of MoA dataset
class MoADataset:
    def __init__(self, features, targets):
        self.features = features
        self.targets = targets

def __len__(self):
        return (self.features.shape[0])

def __getitem__(self, idx):
```

```
[9]: # Pytorch model for the MoA determination
     class Model(nn.Module):
         # Instantiaing all the models before utilizing
         # them later in the forward function.
         def __init__(self, num_features, num_targets, hidden_size):
             # super keyword used to access data from the parent
             # pytorch.nn.Module class
             super(Model, self). init ()
             # Applying batch normalization. This is done to standardize
             # the input for each mini batches and will help reduce the
             # number of epochs for which the training is done. This limits
             # the covariate shift (this is the value by which the hidden
             # layer values shift around) and allows to learn from a more
             # stable set of data. Sometimes, it also allows for a
             # higher learning rate. This is also used for regularization
             # and helps reduce over fitting. Generally, if batch
             # normalization is used, you can use a smaller dropout,
             # which in turn means that lesser layers can be lost
             # in every step.
             self.batch_normalization_1 = nn.BatchNorm1d(num_features)
             # For regularization purposes the dropout is set
             # This is done by setting a probablity. Random
             # neural networks are picked at a probablity, say p
             # or dropped at a probablity of 1-p. This is essential
```

```
# to prevent overfitting of the model and also reduces
       # the computation time. A fully connected neural network, if
       # run without dropout will start forming dependancies between
       # each other and this can lead to over-fitting.
       self.dropoutlayer_1 = nn.Dropout(0.2)
       # nn.utils.weight_norm : This is weight normalization. Usually,
                                faster than batch normalization
       # nn.Linear : Applying linear transform to the incoming data
                     and creates a single layer feed forward network.
       # input size : num features
       # output size : hidden size
       self.denselayer_1 = nn.utils.weight_norm(nn.Linear(num_features,__
→hidden size))
       self.batch_normalization_2 = nn.BatchNorm1d(hidden_size)
       self.dropoutlayer_2 = nn.Dropout(0.2)
       # input size : hidden size
       # output size : hidden_size
       self.denselayer_2 = nn.utils.weight_norm(nn.Linear(hidden_size,__
→hidden_size))
       self.batch_normalization_3 = nn.BatchNorm1d(hidden_size)
       self.dropoutlayer_3 = nn.Dropout(0.1)
       # input size : hidden_size
       # output size : hidden_size
       self.denselayer_3 = nn.utils.weight_norm(nn.Linear(hidden_size,__
→hidden size))
       self.batch normalization 4 = nn.BatchNorm1d(hidden size)
       self.dropoutlayer_4 = nn.Dropout(0.1)
       # input size : hidden size
       # output size : hidden_size
       self.denselayer_4 = nn.utils.weight_norm(nn.Linear(hidden_size,__
→hidden_size))
       self.batch_normalization_5 = nn.BatchNorm1d(hidden_size)
       self.dropoutlayer_5 = nn.Dropout(0.1)
       # input size : hidden_size
       # output size : num_targets
       self.denselayer_5 = nn.utils.weight_norm(nn.Linear(hidden_size,__
→num_targets))
   # The forward function basically defines the model
   def forward(self, forward_x):
       forward_x = self.batch_normalization_1(forward_x)
```

```
forward_x = self.dropoutlayer_1(forward_x)
forward_x = F.relu(self.denselayer_1(forward_x))

forward_x = self.batch_normalization_2(forward_x)
forward_x = self.dropoutlayer_2(forward_x)
forward_x = F.relu(self.denselayer_2(forward_x))

forward_x = self.batch_normalization_3(forward_x)
forward_x = self.dropoutlayer_3(forward_x)
forward_x = self.denselayer_3(forward_x)

forward_x = self.denselayer_4(forward_x)

forward_x = self.dropoutlayer_4(forward_x)

forward_x = self.denselayer_4(forward_x)

forward_x = self.denselayer_5(forward_x)

forward_x = self.dropoutlayer_5(forward_x)

forward_x = self.denselayer_5(forward_x)

return forward_x
```

```
[10]: # Function to train the model
      def trainingFunction(model, optimizer, scheduler, lossFunction, trainloader,
       →device code):
          model.train()
          training loss = 0
          for training_data in trainloader:
              optimizer.zero_grad()
              inputs, targets = training_data['x'].to(device_code),__
       →training_data['y'].to(device_code)
              outputs = model(inputs)
              loss = lossFunction(outputs, targets)
              loss.backward()
              optimizer.step()
              scheduler.step()
              training_loss += loss.item()
          training_loss /= len(trainloader)
          return training_loss
```

```
outputs = model(inputs)
  loss = lossFunction(outputs, targets)
  validation_loss += loss.item()
  validation_predictions.append(outputs.sigmoid().detach().cpu().numpy())
  validation_loss /= len(validationloader)
  validation_predictions = np.concatenate(validation_predictions)
  return validation_loss, validation_predictions
```

```
[12]: # Adding the inference function
def inferenceFunction(model, inferenceloader, device_code):
    model.eval()
    inferences = []
    for data in inferenceloader:
        inputs = data['x'].to(device_code)
        with torch.no_grad():
            outputs = model(inputs)
        inferences.append(outputs.sigmoid().detach().cpu().numpy())
    inferences = np.concatenate(inferences)
    return inferences
```

```
[13]: # Adding dummy inserts to the cp_time and cp_dose columns
# Usually done to categorical variables
def addDummies(data):
    dummy_data = pd.get_dummies(data, columns=['cp_time','cp_dose'])
    return dummy_data
```

#### 5 Preparing the dataset

```
[14]: set_seed_characteristics(seed=55)

[15]: # Seperating out the Gene expression Column and Cell Viability Column

gene_expression = [g for g in training_features.columns if g.startswith('g-')]
    cell_viability = [c for c in training_features.columns if c.startswith('c-')]
```

```
[16]: # Since our dimensions are really high, we can resort to
# using PCA for dimensionality reduction, but is still able
# to capture the characteristics of the data.

# Now, this can be done by choosing a random dimension, and
# having the same random state as before. By doing this
# we observe that we do not encounter
# any 'nan' errors during training.

# Doing PCA for the Gene expression data
```

```
# can choose any random number here
      random_pca_dimension_genes = 20
      # Concatenating the training and test set
      data = pd.concat([pd.DataFrame(training_features[gene_expression]), pd.
       →DataFrame(testing_features[gene_expression])])
      # Performing PCA and converting to a random_pca_dimension_genes number of _{f U}
      → columns
      pca_genes = PCA(n_components = random_pca_dimension_genes, random_state=55)
      # Fitting the PCA transform
      data_pca = pca_genes.fit_transform(data[gene_expression])
      # Splitting the training and test columns
      train_pca_genes = data_pca[:training_features.shape[0]]
      test_pca_genes = data_pca[-testing_features.shape[0]:]
      # Converting training and testing into Pandas data frame shape
      train_pca_genes = pd.DataFrame(train_pca_genes, columns=[f'pca_G-{i}' for i in_
       →range(random_pca_dimension_genes)])
      test pca genes = pd.DataFrame(test pca genes, columns=[f'pca G-{i}' for i in_|
      →range(random_pca_dimension_genes)])
      # Concatenating these back to the original features
      training_features = pd.concat((training_features, train_pca_genes), axis=1)
      testing_features = pd.concat((testing_features, test_pca_genes), axis=1)
[17]: # Doing PCA for the Cell Viability Data
      # can choose any random number here
      random_pca_dimension_cells = 32
      # Concatenating the training and test set
      data = pd.concat([pd.DataFrame(training_features[cell_viability]), pd.
      →DataFrame(testing_features[cell_viability])])
      # Performing PCA and converting to a random pca dimension cells number of
      \rightarrow columns
      pca_cells = PCA(n_components = random_pca_dimension_cells, random_state=55)
      # Fitting the PCA transform
      data_pca = pca_cells.fit_transform(data[cell_viability])
      # Splitting the training and test columns
```

```
train_pca_cells = data_pca[:training_features.shape[0]]
      test_pca_cells = data_pca[-testing_features.shape[0]:]
      # Converting training and testing into Pandas data frame shape
      train_pca_cells = pd.DataFrame(train_pca_cells, columns=[f'pca_C-{i}' for i in_
      →range(random_pca_dimension_cells)])
      test_pca_cells = pd.DataFrame(test_pca_cells, columns=[f'pca_C-{i}' for i inu
       →range(random_pca_dimension_cells)])
      # Concatenating these back to the original features
      training_features = pd.concat((training_features, train_pca_cells), axis=1)
      testing features = pd.concat((testing features, test pca cells), axis=1)
[18]: # Setting a desired threshold to calculate the VarianceThreshold.
      # As per the math all the Features with a training-set variance
      # lower than this threshold will be removed.
      variancethreshold = VarianceThreshold(threshold=0.7)
      # Combining training and test features to create a single dataset
      combined_data = training_features.append(testing_features)
      # Fits to the data, before transforming it
      combined_data_transformed = variancethreshold.fit_transform(combined_data.iloc[:

→, 4:])

      # Extracting the training and the testing data out of the
      # transformed data
      training_features_transformed = combined_data_transformed[ : training_features.
       \rightarrowshape [0]]
      testing_features_transformed = combined_data_transformed[-testing_features.
       \rightarrowshape[0]:]
[19]: # Extracting the training features in a suitable
      # pandas dataset format and numbering the columns
      # after the labels of 'siq_id', 'cp_type', 'cp_time', 'cp_dose'.
      training_features = pd.
       →DataFrame(training_features[['sig_id','cp_type','cp_time','cp_dose']].values.
      →reshape(-1, 4), columns=['sig_id','cp_type','cp_time','cp_dose'])
      training_features = pd.concat([training_features, pd.
      →DataFrame(training_features_transformed)], axis=1)
      # Extracting the testing features in a suitable
      # pandas dataset format and numbering the columns
      # after the labels of 'sig_id', 'cp_type', 'cp_time', 'cp_dose'.
```

```
testing_features = pd.

→DataFrame(testing_features[['sig_id','cp_type','cp_time','cp_dose']].values.

→reshape(-1, 4), columns=['sig_id','cp_type','cp_time','cp_dose'])

testing_features = pd.concat([testing_features, pd.

→DataFrame(testing_features_transformed)], axis=1)
```

```
[21]: # Extracting the columns of the drugs that are sold from
# the train pandas dataframe

target = train[training_targets_scored.columns]

# Now that the ctl_vehicle drugs have been removed, we do not need
# cp_type. So we can go ahead and remove that columns as well.

train = train.drop('cp_type', axis=1)
test = test.drop('cp_type', axis=1)

# extracting the columns in the targets

target_columns = target.drop('sig_id', axis=1).columns.values.tolist()
```

#### 6 Dataset preparation complete

```
[22]: # multilabel stratified K Fold import causes a small warning and we do not want # to show that in the notebook.
warnings.filterwarnings('ignore')
```

```
folds = train.copy()
number_of_folds = 3
# creating a 3 fold multilabel stratified K Fold
multilabel_k fold = MultilabelStratifiedKFold(n_splits = number_of_folds)
# Standard k fold splitting. Here we are splitting into number_of_folds folds
for fol, (train_folds, validation_folds) in enumerate(multilabel_k_fold.
⇒split(X=train, y=target)):
   folds.loc[validation_folds, 'kfold'] = int(fol)
folds['kfold'] = folds['kfold'].astype(int)
# Isolating out the feature columns. This is done by first
# Isolating the columns that are not present in the target
# followed by extracting the columns except the sig id and
# kfold.
feature_columns = [c for c in addDummies(folds).columns if c not in_
→target columns]
feature_columns = [c for c in feature_columns if c not in ['kfold', 'sig_id']]
```

## 7 Declaring the HyperParameters

```
[23]: BATCH SIZE = 128
      max epochs = 16
      # When training neural networks, it is common to use
      # weight decay where after each update, the weights
      # are multiplied by a factor slightly less than 1
      weight_decay = 1e-5
      # deciding the initial learning rate
      # It controls how quickly or slowly a neural
      # network model can learn a model or a problem.
      lr = 1e-3
      # Boolean to decide on stopping early when the
      # validation loss > best loss
      bool early stop = True
      # steps to execute before early stopping
      steps_early_stopping= 10
      # number of features corresponding to the columns in the
      # targets
      num_features=len(feature_columns)
      # number of targets corresponding to the columns in the
      # features
```

```
num_targets=len(target_columns)
# in between neural netwrok size
hidden_size=1024
```

#### 8 Declaring the training functions and performing the training

```
[24]: # to plot validation loss
      valid_loss_list = []
      # to plot the training loss
      train_loss_list = []
      # to plot the best recorded loss
      best_loss_list = []
[25]: def run_training(fold, seed):
          # declaring the list as global to plot validation loss
          global valid_loss_list
          # declaring the training loss list as global to plot
          # the training loss
          global train_loss_list
          # declaring the best loss list as global to plot the
          # best losses recorded
          global best_loss_list
          # setting the seed to start from the same number as
          # explained previously
          set_seed_characteristics(seed)
          # adding dummy variables to the training set
          train = addDummies(folds)
          # extracting the validating rows numbers for the
          # respective k fold values
          val_idx = train[train['kfold'] == fold].index
          # Dropping all the rows from the training set
          # that does not belong to this kth fold
          train_necessary_rows = train[train['kfold'] != fold].reset_index(drop=True)
          # Dropping all the rows from the valiadtion set
          # that does not belong to this kth fold
          valid_necessary_rows = train[train['kfold'] == fold].reset_index(drop=True)
          # splitting the x and y values for training set
          train_features, train_targets = train_necessary_rows[feature_columns].
       →values, train_necessary_rows[target_columns].values
```

# splitting the x and y values for test set

```
validation_features, validation_targets = __
→valid_necessary_rows[feature_columns].values,
→valid_necessary_rows[target_columns].values
   # Converting the training data to standard pytorch
   # dataset class format
   train_dataset = MoADataset(train_features, train_targets)
   # Converting the validation data to standard pytorch
   # dataset class format
   valid_dataset = MoADataset(validation_features, validation_targets)
   # calling the pytorch data loading utility for the
   # training set
   trainloader = torch.utils.data.DataLoader(train_dataset,_
→batch_size=BATCH_SIZE, shuffle=True)
   # calling the pytorch data loading utility for the
   # validation set
   validloader = torch.utils.data.DataLoader(valid_dataset,__
→batch_size=BATCH_SIZE, shuffle=False)
   # Declaring the model and can be tuned here
   # using the hyper parameters
   model = Model(
       num_features=num_features,
       num_targets=num_targets,
       hidden_size=hidden_size,
   )
   # moving the model to GPU if available,
   # else will run it on CPU itself
   model.to(device_code)
   # A standard optimizer. Adam optimizer is widely used
   # because it combines the advantages of the Adaptive gradient
   # algorithm and the root mean square propogation. Basically, it does
   # not stick to one learning rate and adapts it to the problem.
   # It is widely known to offer good results really fast.
   optimizer = torch.optim.Adam(model.parameters(), lr=lr,_
→weight_decay=weight_decay)
   # We use a learning rate scheduler to converge to the lowest
   # loss faster. This is also seen to provide higher accuracy.
   # This can be tuned.
   # Some of the optimizers I tried here are
   # optim.lr_scheduler.OneCycleLR
   # optim.lr_scheduler.StepLR
```

```
# scheduler = optim.lr_scheduler.StepLR(optimizer, step_size=1, qamma=0.1)
   scheduler = optim.lr_scheduler.OneCycleLR(optimizer=optimizer, pct_start=0.
\rightarrow05, div_factor=1.5e3,
                                             max_lr=1e-2, epochs=max_epochs,_
⇒steps per epoch=len(trainloader))
   # after research I saw that the Binary cross
   # entropy loss with sigmoid layer works well
   lossFunction = nn.BCEWithLogitsLoss()
   # stops when the error starts increaseing. Setting the counter
   # to track this
   steps_before_early_stop = 0
   # general out of fold array shape
   out_of_fold = np.zeros((len(train), target.iloc[:, 1:].shape[1]))
   # declaring a very high value as an
   # initial loss for each kth fold
   best_loss = np.inf
   # looping through the epochs
   for epoch in range(max_epochs):
       # training the model
       training_loss = trainingFunction(model, optimizer,scheduler,__
→lossFunction, trainloader, device_code)
       print('epoch : ',epoch,'>> training loss : ',training loss)
       train_loss_list.append(training_loss)
       validation_loss, validation_predictions = validationFunction(model,_
→lossFunction, validloader, device_code)
       print('epoch : ',epoch,'>> validation : ',validation_loss)
       valid_loss_list.append(validation_loss)
       # checking if the loss is decreasing
       if validation loss < best loss:</pre>
           best_loss = validation_loss
           best_loss_list.append(best_loss)
           # Updating the out of fold predictions
           out_of_fold[val_idx] = validation_predictions
           # saving the model and data for this kth fold
           torch.save(model.state_dict(), f"FOLD{fold}_.pth")
       # Handling the increasing loss by calling
       # early stopping
       elif(bool_early_stop == True):
           # breaks out of the loop when this happens
```

```
steps_before_early_stop += 1
           if (steps_before_early_stop >= steps_early_stopping):
               break
   # adding dummy variables to the test set
   test_ = addDummies(test)
   # extracting the x_test
   x_test = test_[feature_columns].values
   testdataset = TestDataset(x_test)
   testloader = torch.utils.data.DataLoader(testdataset,__
→batch_size=BATCH_SIZE, shuffle=False)
   model = Model(num_features=num_features,num_targets=num_targets,
       hidden_size=hidden_size,
   )
   # uploading the saved data for this kth fold
   model.load_state_dict(torch.load(f"FOLD{fold}_.pth"))
   # again uploading the model to GPU, if available
   model.to(device_code)
   predictions = np.zeros((len(test_), target.iloc[:, 1:].shape[1]))
   # evaluates the model
   predictions = inferenceFunction(model, testloader, device_code)
   return out_of_fold, predictions
```

```
[26]: def executeKFold(number_of_folds, seed):
    # standard size for the out of fold predictions
    out_of_fold = np.zeros((len(train), len(target_columns)))
    # same size for all of the predictions
    predictions = np.zeros((len(test), len(target_columns)))

for each_k_fold in range(number_of_folds):
    print('Fold Number : ', each_k_fold)
    out_of_fold_, pred_ = run_training(each_k_fold, seed)

# adding all the predictions
    predictions += pred_ / number_of_folds
    # adding all the out of fold predictions
    out_of_fold += out_of_fold_
        print("------")

k_th_prediction = predictions
    return out_of_fold, k_th_prediction
```

```
[27]: # setting a standard seed number
SEED = [55]
# general out of fold array shape
out_of_fold = np.zeros((len(train), len(target_columns)))
# general predictions array shape
predictions = np.zeros((len(test), len(target_columns)))

# for seed in SEED:
out_of_fold_, predictions_ = executeKFold(number_of_folds, SEED[0])
out_of_fold += out_of_fold_ / len(SEED)
predictions += predictions_ / len(SEED)

train[target_columns] = out_of_fold
test[target_columns] = predictions
Fold Number : 0
```

```
epoch: 0 >> training loss: 0.2788736069979875
epoch : 0 >> validation : 0.021623456240471066
epoch : 1 >> training_loss : 0.020091968256494272
epoch: 1 >> validation: 0.01922281719101914
epoch : 2 >> training_loss : 0.019972266159627747
epoch: 2 >> validation: 0.01886744482506966
epoch : 3 >> training_loss : 0.019043210343174313
epoch: 3 >> validation: 0.018996690303601068
epoch : 4 >> training_loss : 0.018774466394730235
epoch: 4 >> validation: 0.018413769980442935
epoch : 5 >> training_loss : 0.01869301253362842
epoch : 5 >> validation : 0.01853232747264977
epoch : 6 >> training_loss : 0.01865103912094365
epoch : 6 >> validation : 0.018168202376571196
epoch: 7 >> training loss: 0.018572327146387617
epoch: 7 >> validation: 0.01888334468520921
epoch : 8 >> training_loss : 0.018362585228422414
epoch : 8 >> validation : 0.018195184943234098
epoch : 9 >> training_loss : 0.018180621188619864
epoch: 9 >> validation: 0.01839715341941036
epoch : 10 >> training_loss : 0.017992399317090926
epoch: 10 >> validation: 0.017737267185644858
epoch : 11 >> training_loss : 0.01772677853865468
epoch: 11 >> validation: 0.017560527595723498
epoch : 12 >> training_loss : 0.017529506098641002
epoch: 12 >> validation: 0.01734824900932867
epoch : 13 >> training loss : 0.017258823679193208
epoch : 13 >> validation : 0.01717964628839801
epoch : 14 >> training loss : 0.01701616107121758
epoch: 14 >> validation: 0.01715383331837325
epoch: 15 >> training loss: 0.016877019105722074
```

\_\_\_\_\_ Fold Number: 1 epoch : 0 >> training\_loss : 0.2791394925635794 epoch : 0 >> validation : 0.021501707748092454 epoch: 1 >> training loss: 0.02014602518276028 epoch: 1 >> validation: 0.019920575047104525 epoch : 2 >> training\_loss : 0.019292888320658518 epoch: 2 >> validation: 0.020500779746036077 epoch : 3 >> training\_loss : 0.01898241023654523 epoch: 3 >> validation: 0.019408218819519568 epoch : 4 >> training\_loss : 0.01884553834957921 epoch: 4 >> validation: 0.018803634969838733 epoch : 5 >> training\_loss : 0.018704059833417767 epoch: 5 >> validation: 0.01929768891041649 epoch : 6 >> training\_loss : 0.018495980099491452 epoch: 6 >> validation: 0.018345469753418504 epoch : 7 >> training\_loss : 0.018502545138092144 epoch: 7 >> validation: 0.0183179613748758 epoch: 8 >> training loss: 0.01830926013543554 epoch: 8 >> validation: 0.018034464246111697 epoch: 9 >> training loss: 0.01814349767468546 epoch: 9 >> validation: 0.018158412088864838 epoch : 10 >> training\_loss : 0.01796842050617156 epoch: 10 >> validation: 0.017744970556091647 epoch : 11 >> training\_loss : 0.017735546859710113 epoch: 11 >> validation: 0.01759479866074077 epoch : 12 >> training\_loss : 0.01754570741854284 epoch: 12 >> validation: 0.017433413593419666 epoch : 13 >> training\_loss : 0.017299547661905702 epoch: 13 >> validation: 0.017249190232491697 epoch : 14 >> training\_loss : 0.017085145907881467 epoch: 14 >> validation: 0.017144549509574628 epoch : 15 >> training\_loss : 0.01694276431656402 epoch: 15 >> validation: 0.01715218401032275 \_\_\_\_\_ Fold Number: 2 epoch : 0 >> training\_loss : 0.28062024895587695 epoch: 0 >> validation: 0.021831682176682455 epoch : 1 >> training\_loss : 0.019925031827195832 epoch: 1 >> validation: 0.02110225957786215 epoch : 2 >> training\_loss : 0.019255759826172954 epoch: 2 >> validation: 0.01923344059494035 epoch : 3 >> training\_loss : 0.01897560269774302 epoch: 3 >> validation: 0.01963105608291667 epoch : 4 >> training\_loss : 0.01888560847905667 epoch: 4 >> validation: 0.018987211848384346 epoch : 5 >> training\_loss : 0.018881849539668663

epoch: 15 >> validation: 0.01710507392497926

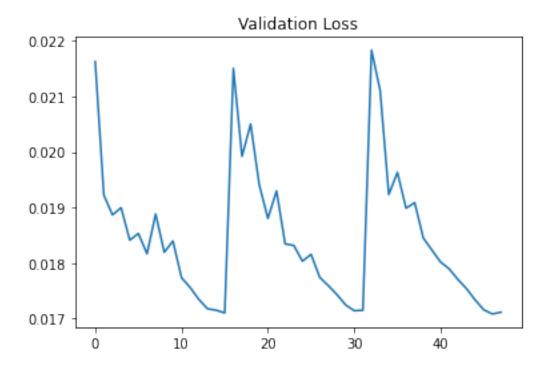
```
epoch : 5 >> validation : 0.019087364800788206
epoch : 6 >> training_loss : 0.01880092287193174
epoch: 6 >> validation: 0.018451588421032346
epoch : 7 >> training_loss : 0.01856426177141459
epoch: 7 >> validation: 0.018233612982620453
epoch : 8 >> training_loss : 0.018412961470692053
epoch: 8 >> validation: 0.018018708034450638
epoch : 9 >> training_loss : 0.018320308147889118
epoch: 9 >> validation: 0.01789587644603232
epoch : 10 >> training_loss : 0.018012238172409326
epoch: 10 >> validation: 0.017706910565752406
epoch : 11 >> training_loss : 0.01781968524598557
epoch : 11 >> validation : 0.01754140906870879
epoch : 12 >> training_loss : 0.017585085947876392
epoch: 12 >> validation: 0.017338143388644374
epoch : 13 >> training_loss : 0.01736602078637351
epoch: 13 >> validation: 0.017160253085452933
epoch : 14 >> training_loss : 0.01719682110554498
epoch: 14 >> validation: 0.017085314712262357
epoch: 15 >> training loss: 0.017056497793806635
epoch: 15 >> validation: 0.0171179612135065
```

## 9 Evaluating the logarithmic loss function applied to each drug-MoA annotation pair.

The Cross Validation loss is :>> 0.015734850054263057

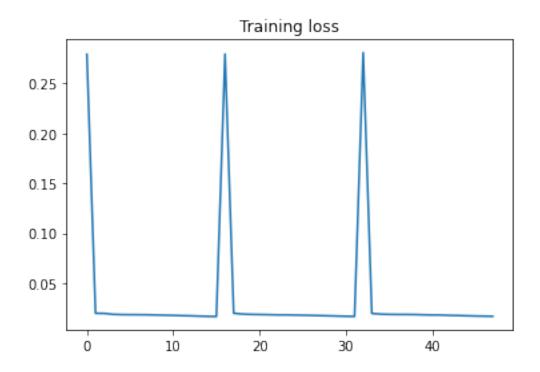
#### 10 Plotting the Validation loss for each fold

```
[29]: plt.plot(valid_loss_list)
    plt.title('Validation Loss')
    plt.savefig('valid_loss_list.png')
```



## 11 Plotting the Training loss for each fold

```
[30]: plt.plot(train_loss_list) plt.title('Training loss') plt.savefig('train_loss_list.png')
```



## 12 Plotting the Best recorded loss for each fold

```
[31]: plt.plot(best_loss_list)
    plt.title('Best Recorded loss')
    plt.savefig('best_loss_list.png')
```

