

## ▼ Prerequisite Packages

```
# prerequisites
!pip3 install cloudmesh-installer
!pip3 install cloudmesh-common
!pip3 install kaggle

Downloading https://files.pythonhosted.org/packages/87/89/479dc97e18549e21354893e4e
Requirement already satisfied: py>=1.5.0 in /usr/local/lib/python3.7/dist-packages (f
Requirement already satisfied: atomicwrites>=1.0 in /usr/local/lib/python3.7/dist-pac
Requirement already satisfied: attrs>=17.4.0 in /usr/local/lib/python3.7/dist-package
Requirement already satisfied: pluggy<0.8,>=0.5 in /usr/local/lib/python3.7/dist-pack
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Requirement already satisfied: zipp>=0.5 in /usr/local/lib/python3.7/dist-packages (f
Building wheels for collected packages: python-hostlist, ordered-set
  Building wheel for python-hostlist (setup.py) ... done
  Created wheel for python-hostlist: filename=python_hostlist-1.21-cp37-none-any.whl
  Stored in directory: /root/.cache/pip/wheels/0b/5b/55/ddcf52288f0b10f4564ca1b253159
  Building wheel for ordered-set (setup.py) ... done
  Created wheel for ordered-set: filename=ordered_set-4.0.2-py2.py3-none-any.whl size
  Stored in directory: /root/.cache/pip/wheels/e1/c6/9b/651d8a21d59b51a75ab9c070838f9
Successfully built python-hostlist ordered-set
ERROR: pytest-cov 2.11.1 has requirement coverage>=5.2.1, but you'll have coverage 3.
ERROR: pytest-cov 2.11.1 has requirement pytest>=4.6, but you'll have pytest 3.6.4 wh
Installing collected packages: bump2version, oyaml, simplejson, python-hostlist, colc
Successfully installed bump2version-1.0.0 cloudmesh-common-4.3.66 cloudmesh-installer
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Requirement already satisfied: idna<3,>=2.5 in /usr/local/lib/python3.7/dist-packages
```

## ▼ Sign Up for Kaggle Account and Generate API Token

1. Sign Up for Kaggle account at <https://www.kaggle.com>.
2. In Kaggle 'Profile'-'>'Account', generate api token by clicking 'Create New API Token'.

## Upload Token into Colab

3. Upload json file, with Kaggle api token and username, into Colab by running code below.  
Select downloaded 'kaggle.json' file when prompted and click 'Upload'.

NB third-party cookies should be enabled for upload to work.

```
from google.colab import files

# upload json file with api token and username
files.upload()

# create directory for token in Colab
!mkdir ~/.kaggle

# move file to directory
!mv kaggle.json ~/.kaggle/

# modify permissions on directory
!chmod 600 ~/.kaggle/kaggle.json
```

Choose Files kaggle.json

- **kaggle.json**(application/json) - 67 bytes, last modified: 5/3/2021 - 100% done  
Saving kaggle.json to kaggle.json

## ▼ Download Merck Molecular Activity Challenge Dataset

```
# download dataset
!kaggle competitions download -c MerckActivity

# make train and test directories for raw data
!mkdir -p /content/raw/train /content/raw/test

# unzip train and test datasets into respective folders; only overwrite exiting files if extr
!unzip /content/TrainingSet.zip -d /content/raw/train

!unzip /content/TestSet.zip -d /content/raw/test

# paths to raw data
path_to_raw_train_data = '/content/raw/train/TrainingSet/'
path_to_raw_test_data = '/content/raw/test/TestSet/'
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Warning: Looks like you're using an outdated API Version, please consider updating (ser  
TestSet.7z: Skipping, found more recently modified local copy (use --force to force dow  
Rsquared.R: Skipping, found more recently modified local copy (use --force to force dow  
ntree20\_benchmark.R: Skipping, found more recently modified local copy (use --force to  
ntree20\_basicBenchmark.csv.zip: Skipping, found more recently modified local copy (use  
TrainingSet.7z: Skipping, found more recently modified local copy (use --force to force  
TrainingSet.zip: Skipping, found more recently modified local copy (use --force to forc  
TestSet.zip: Skipping, found more recently modified local copy (use --force to force do  
Archive: /content/TrainingSet.zip

creating: /content/raw/train/TrainingSet/  
inflating: /content/raw/train/TrainingSet/ACT10\_competition\_training.csv  
inflating: /content/raw/train/TrainingSet/ACT11\_competition\_training.csv  
inflating: /content/raw/train/TrainingSet/ACT12\_competition\_training.csv  
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inflating: /content/raw/train/TrainingSet/ACT9\_competition\_training.csv

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inflating: /content/raw/test/TestSet/ACT2\_competition\_test.csv  
inflating: /content/raw/test/TestSet/ACT3\_competition\_test.csv  
inflating: /content/raw/test/TestSet/ACT4\_competition\_test.csv  
inflating: /content/raw/test/TestSet/ACT5\_competition\_test.csv  
inflating: /content/raw/test/TestSet/ACT6\_competition\_test.csv

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inflating: /content/raw/test/TestSet/ACT7_competition_test.csv  
inflating: /content/raw/test/TestSet/ACT8_competition_test.csv  
inflating: /content/raw/test/TestSet/ACT9_competition_test.csv
```

## ▼ Preprocess Dataset

```
import pandas as pd  
from cloudmesh.common.StopWatch import StopWatch  
  
# make train and test directories for preprocessed data  
!mkdir -p /content/preprocessed/train /content/preprocessed/test  
  
# paths to processed data  
path_to_preprocessed_train_data = '/content/preprocessed/train/'  
path_to_preprocessed_test_data = '/content/preprocessed/test/'  
  
# cycle through 15 data sets preprocessing them for learning  
StopWatch.start("preprocessing")  
dataset_file_no = 1  
  
while dataset_file_no <= 15:  
  
    dataset_train_file_name = 'ACT' + str(dataset_file_no) + '_competition_training.csv'  
    dataset_test_file_name = 'ACT' + str(dataset_file_no) + '_competition_test.csv'  
  
    train_filename = path_to_raw_train_data + dataset_train_file_name  
    test_filename = path_to_raw_test_data + dataset_test_file_name  
  
    train_filename_processed = path_to_preprocessed_train_data + dataset_train_file_name  
    test_filename_processed = path_to_preprocessed_test_data + dataset_test_file_name  
  
    print ('Preprocessing dataset ', 'ACT' + str(dataset_file_no))  
  
    train = pd.read_csv(train_filename)  
    test = pd.read_csv(test_filename)  
  
    print (len(train.columns.values))  
    print (len(test.columns.values))  
  
    train_inx_set = set(train.columns.values)  
    test_inx_set = set(test.columns.values)  
  
    # remove molecule label and columns that are not common to both training and test sets  
    train_inx = [inx for inx in train.columns.values if inx in set.intersection(train_inx_set,  
    test_inx = [inx for inx in test.columns.values if inx in set.intersection(train_inx_set,  
  
    train_inx.insert(0, 'Act')  
    train_inx.remove('MOLECULE')
```

```

test_inx.remove('MOLECULE')

#print (train_inx)
#print (test_inx)

train = train[train_inx]
test = test[test_inx]

#print (train.shape)
#print (test.shape)

# save data to csv
train.to_csv(train_filename_processed, index=False)
test.to_csv(test_filename_processed, index=False)

print(train.head(5))
print ('Preprocessing dataset ', 'ACT' + str(dataset_file_no), ' complete')

dataset_file_no += 1
StopWatch.stop("preprocessing")

StopWatch.benchmark()

```

```

    ....
      Act  D_6  D_37  D_38  D_39  ...  D_10963  D_10995  D_11012  D_11027  D_11030
0  8.6182    0    0    0    0  ...         0         0         0         0         0
1  9.8521    0    0    0    0  ...         0         0         0         0         0
2  8.3264    0    0    0    0  ...         0         0         0         0         0
3  8.2581    0    0    0    0  ...         0         0         0         0         0
4  7.3552    0    0    0    0  ...         0         0         0         0         0

```

```

[5 rows x 5553 columns]
Preprocessing dataset  ACT15  complete

```

Attribute	Value
BUG_REPORT_URL	<a href="https://bugs.launchpad.net/ubuntu/">https://bugs.launchpad.net/ubuntu/</a>
DISTRIB_CODENAME	bionic
DISTRIB_DESCRIPTION	"Ubuntu 18.04.5 LTS"
DISTRIB_ID	Ubuntu
DISTRIB_RELEASE	18.04
HOME_URL	<a href="https://www.ubuntu.com/">https://www.ubuntu.com/</a>
ID	ubuntu
ID_LIKE	debian
NAME	"Ubuntu"
PRETTY_NAME	"Ubuntu 18.04.5 LTS"
PRIVACY_POLICY_URL	<a href="https://www.ubuntu.com/legal/terms-and-policies/privacy-policy">https://www.ubuntu.com/legal/terms-and-policies/privacy-policy</a>
SUPPORT_URL	<a href="https://help.ubuntu.com/">https://help.ubuntu.com/</a>
UBUNTU_CODENAME	bionic
VERSION	"18.04.5 LTS (Bionic Beaver)"
VERSION_CODENAME	bionic
VERSION_ID	"18.04"
cpu_count	4
mem.active	1.0 GiB

```

| mem.available      | 24.4 GiB
| mem.free           | 938.3 MiB
| mem.inactive       | 22.8 GiB
| mem.percent        | 4.3 %
| mem.total          | 25.5 GiB
| mem.used           | 10.9 GiB
| platform.version   | #1 SMP Thu Jul 23 08:00:38 PDT 2020
| python             | 3.7.10 (default, May 3 2021, 02:48:31)
|                   | [GCC 7.5.0]
| python.pip         | 19.3.1
| python.version     | 3.7.10
| sys.platform       | linux
| uname.machine      | x86_64
| uname.node         | 7e0207301e13
| uname.processor    | x86_64
| uname.release      | 4.19.112+
| uname.system       | Linux
| uname.version      | #1 SMP Thu Jul 23 08:00:38 PDT 2020
| user               | collab

```

```

+-----+-----+-----+-----+-----+-----+-----+
| Name      | Status | Time  | Sum  | Start                | tag  | Node  |
+-----+-----+-----+-----+-----+-----+-----+
| preprocessing | ok      | 536.097 | 536.097 | 2021-05-07 17:08:11 |      | 7e0207
+-----+-----+-----+-----+-----+-----+-----+

```

## ▼ Predicting Molecular Activity

```

from __future__ import absolute_import
from __future__ import division
from __future__ import print_function
import numpy as np
import pandas as pd
import keras.backend as K
import tensorflow as tf
from keras.models import Sequential
from keras.layers import Dense, Activation, Dropout, InputLayer, Concatenate
from keras.utils import to_categorical, plot_model
from keras.utils.vis_utils import model_to_dot
from keras.optimizers import Adam
from cloudmesh.common.StopWatch import StopWatch

# make outputs directory
!mkdir -p /content/outputs/

path_to_train_data = '/content/preprocessed/train/'
path_to_test_data = '/content/preprocessed/test/'
path_to_outputs = '/content/outputs/'

```

```
# define parameters
```

```
# define parameters
hidden_units = 512
dropout = 0.45
BATCH_SIZE = 128
feature_dim = 128
opti = Adam(lr=0.0001, beta_1=0.5)

# define fully connected network/MLP
def fcn_model(input_shape=(feature_dim,)):
    model = Sequential()
    model.add(Activation('relu'))
    model.add(Dropout(dropout))
    model.add(Dense(hidden_units))
    model.add(Activation('relu'))
    model.add(Dropout(dropout))
    model.add(Dense(hidden_units))
    model.add(Activation('relu'))
    model.add(Dropout(dropout))
    model.add(Dense(hidden_units))
    model.add(Activation('relu'))
    model.add(Dropout(0.10))
    model.add(Dense(num_labels))
    model.add(Activation('softmax'))
    model.build(input_shape)
    model.summary()
    model.compile(loss='mean_squared_error', optimizer=opti, metrics=[Rsquared])
    print("\nTraining on dataset number", act_ds, " of 15:\n")
    model.fit(x_train, y_train, epochs=15, batch_size=BATCH_SIZE)
    loss, R2 = model.evaluate(x_test, y_test, batch_size=BATCH_SIZE)
    print("\nCorrelation coefficient:", R2)
    return model

# define correlation coefficient (R^2) formula
def Rsquared(x,y):

    # approach adopted from RuwanT
    # URL: https://github.com/RuwanT/merck/blob/master/main.py

    x = K.batch_flatten(x)
    y = K.batch_flatten(y)

    avx = K.mean(x)
    avy = K.mean(y)

    num = K.sum((x-avx) * (y-avy))
    num = num * num

    denom = K.sum((x-avx)*(x-avx)) * K.sum((y-avy)*(y-avy))

    return num/denom

# iterate through 15 distinct datasets of high throughput screening (HTS) assays
```

```
StopWatch.start("train-evaluate-predict")
act_ds = 1
while act_ds <= 15:
    print("\nReading from dataset", act_ds,"of 15:\n")
    data_train_main = pd.read_csv(path_to_train_data + 'ACT' + str(act_ds) + '_competition_tr
    data_ac = pd.read_csv(path_to_test_data + 'ACT' + str(act_ds) + '_competition_test.csv')

    # split each of the datasets into set for training (80%), set for testing/evaluation (10%)
    # set for use with validating prediction (10%)
    data_train = data_train_main.sample(frac = 0.8)
    data_test = data_train_main.drop(data_train.index)
    data_prediction = data_test.sample(frac = 0.5)
    data_test = data_test.drop(data_prediction.index)

    activity_inx = data_train.columns.get_loc('Act')
    feature_dim = data_train.shape[1] - (activity_inx+1)
    #print("no. of feature columns:", feature_dim)

    # identify molecular activity labels
    y_train = data_train['Act']
    #print("shape of y_train:", y_train.shape)
    num_labels = len(np.unique(y_train))
    #print("no. of unique labels:", num_labels)
    y_test = data_test['Act']

    # identify and filter for feature-set/molecular-substructure frequencies
    train_set_inx = set(data_train.columns.values)
    test_set_inx = set(data_ac.columns.values)

    train_inx = [inx for inx in data_train.columns.values if inx in set.intersection(train_set_inx, test_set_inx)]
    test_inx = [inx for inx in data_test.columns.values if inx in set.intersection(train_set_inx, test_set_inx)]
    predict_inx = [inx for inx in data_prediction.columns.values if inx in set.intersection(train_set_inx, test_set_inx)]

    data_train = data_train[train_inx]
    data_test = data_test[test_inx]
    data_prediction = data_prediction[predict_inx]
    #print(data_train.head(5))

    # format feature-set input data
    x_train = data_train[0:]
    x_train = np.asarray(x_train).astype('float32')
    x_test = data_test[0:]
    x_test = np.asarray(x_test).astype('float32')
    x_predict = data_prediction[0:]
    x_predict = np.asarray(x_predict).astype('float32')

    input_size = x_train.shape[0]

    # call fully connected network for learning and evaluation. Predict biological activity f
    y_predict = fcn_model(input_shape=(input_size, feature_dim)).predict(x_predict, batch_size=
```



```
# display inputs and outputs for the prediction step and write results to text file
```

```
prediction_file = open(path_to_outputs + "molecular_activity_prediction.txt", "a")
```

```
for i in range(len(x_predict)):
```

```
    print("Substructure/feature frequencies=%s, Predicted biological activity=%s" % (x_predict[i], y_predict[i]))
    prediction_file.write(str(x_predict[i]) + ", " + str(np.mean(y_predict[i])) + "\n")
```

```
    act_ds += 1
```

```
prediction_file.close()
```

```
StopWatch.stop("train-evaluate-predict")
```

```
StopWatch.benchmark()
```

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Substructure/feature frequencies=[0. 0. 0. ... 0. 0. 0.], Predicted biological activity=0.0
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Substructure/feature frequencies=[0. 0. 0. ... 0. 0. 0.], Predicted biological activity=0.0
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