

final

December 12, 2024

0.1 Abstract

0.2 Package import

```
[1]: import os

os.environ["KERAS_BACKEND"] = "tensorflow"

import ast
import numpy as np

from tensorflow import keras
```

```
[2]: #from tensorflow.keras import ops
from tensorflow.keras import layers
import pandas as pd

from sklearn.model_selection import train_test_split
import matplotlib.pyplot as plt
from rdkit import Chem, RDLogger
from rdkit.Chem import BondType
from rdkit.Chem.Draw import MolsToGridImage
from rdkit.Chem import Draw
from rdkit import Chem
from rdkit.Chem import rdmolops, AllChem
from tensorflow.keras.regularizers import l1_l2
RDLogger.DisableLog("rdApp.*")
```

```
[3]: import tensorflow as tf
print("TensorFlow version:", tf.__version__)
print("GPU available:", tf.config.list_physical_devices('GPU'))
print("GPU in use:", tf.test.gpu_device_name())
```

TensorFlow version: 2.16.2

GPU available: []

GPU in use:

0.3 Database pharsing

```
[4]: '''  
    read the entire dataset  
    '''  
  
df = pd.read_csv('dataset1.csv')  
df.drop([0,1,2,3,4], inplace=True)  
df=df.rename(columns = {'PUBCHEM_EXT_DATASOURCE_SMILES':  
    ↳ 'SMILES', 'PUBCHEM_ACTIVITY_OUTCOME': 'Activity', 'PUBCHEM_ACTIVITY_SCORE':  
    ↳ 'Score'})  
columns_to_drop = [col for col in df.columns if col not in ['SMILES',  
    ↳ 'Activity', 'Score', 'Potency', 'Efficacy']]  
df = df.drop(columns = columns_to_drop)  
#df=df.drop(['Unnamed: 3', 'Unnamed: 4', 'Unnamed: 5'], axis=1)  
df = df.dropna(subset=['SMILES'])  
  
df=df.fillna(0)  
print(df.head())  
print(df.info())
```

	SMILES	Activity	Score \
5	CNCC1=NC2=C(C=C(C=C2)C1)C(=N1)C3=CC=CN3	Inactive	0.0
6	CCSC(=NC1=CC=C(C=C1)C(F)(F)F)N.C1	Inactive	0.0
7	CCN(CC1=CC(=CC=C1)S(=O)(=O)[O-])C2=CC=C(C=C2)C...	Inactive	0.0
8	CC1=CC=C(C=C1)S(=O)(=O)N2CCN(CC2)C3=NC(=NC4=CC...	Inactive	0.0
9	CC1=CC=C(C=C1)S(=O)(=O)N2CCN(CC2)C3=NC(=NC4=CC...	Inactive	0.0

	Potency	Efficacy
5	0.0	0.0
6	0.0	0.0
7	0.0	0.0
8	0.0	0.0
9	0.0	0.0

```
<class 'pandas.core.frame.DataFrame'>  
Index: 342051 entries, 5 to 342072  
Data columns (total 5 columns):  
#   Column      Non-Null Count  Dtype  
---  -  
0   SMILES      342051 non-null object  
1   Activity    342051 non-null object  
2   Score       342051 non-null float64  
3   Potency     342051 non-null float64  
4   Efficacy    342051 non-null float64  
dtypes: float64(3), object(2)  
memory usage: 15.7+ MB  
None
```

```
[5]: valid_indices = []
# Loop through each SMILES string in the DataFrame
for i in range(len(df)):
    smiles = df.iloc[i]['SMILES'] # Use iloc for positional indexing

    # Convert SMILES to molecule
    mol = Chem.MolFromSmiles(smiles)

    # Check if the molecule is valid and has <= 50 atoms
    if mol is not None and mol.GetNumAtoms() <= 50:
        valid_indices.append(i)
# Filter the DataFrame to include only valid molecules
df_50 = df.iloc[valid_indices]
```

```
[6]: df_50
```

```
[6]:
```

	SMILES	Activity	Score	\
5	CNCC1=NC2=C(C=C(C=C2)C1)C(=N1)C3=CC=CN3	Inactive	0.0	
6	CCSC(=NC1=CC=C(C=C1)C(F)(F)F)N.Cl	Inactive	0.0	
8	CC1=CC=C(C=C1)S(=O)(=O)N2CCN(CC2)C3=NC(=NC4=CC...	Inactive	0.0	
9	CC1=CC=C(C=C1)S(=O)(=O)N2CCN(CC2)C3=NC(=NC4=CC...	Inactive	0.0	
10	C1CN(CCN1C2=NC(=NC3=CC=CC=C32)C4=CC=CS4)S(=O)(...	Inactive	0.0	
...	
342068	CC(=O)NC1=CC=C(C=C1)OCC2=C(C=CC(=C2)CN(CC3=CC=...	Inactive	0.0	
342069	CC(=O)NC1=CC=C(C=C1)OCC2=C(C=CC(=C2)CN(CC3=CC=...	Inactive	0.0	
342070	CC(=O)NC1=CC=C(C=C1)OCC2=C(C=CC(=C2)CN(CC3=CC=...	Inactive	0.0	
342071	CC(=O)NC1=CC=C(C=C1)C(=O)N(CC2=CC=CC=C2)CC3=CC...	Inactive	0.0	
342072	CC(=O)NC1=CC=C(C=C1)OCC2=C(C=CC(=C2)CN(CC3=CC=...	Inactive	0.0	

	Potency	Efficacy
5	0.0	0.0
6	0.0	0.0
8	0.0	0.0
9	0.0	0.0
10	0.0	0.0
...
342068	0.0	0.0
342069	0.0	0.0
342070	0.0	0.0
342071	0.0	0.0
342072	0.0	0.0

[341260 rows x 5 columns]

```
[7]: def is_charged(smiles):
    mol = Chem.MolFromSmiles(smiles)
    if not mol:
```

```

    return False # Invalid SMILES
    return any(atom.GetFormalCharge() != 0 for atom in mol.GetAtoms())

# Test the function
print(is_charged("CC1=C(SC(=C1C#N)NC(=O)C2=CC(C=C2)OC) [N+] (=O)"))

```

True

```
[8]: df_50['Charged'] = df_50['SMILES'].apply(is_charged)
```

```

uncharged = df_50[df_50['Charged'] == False]
uncharged

```

/var/folders/jn/kkchdc94t50xrmycsvkq2x80000gn/T/ipykernel_85584/162626946.py:1:

SettingWithCopyWarning:

A value is trying to be set on a copy of a slice from a DataFrame.

Try using .loc[row_indexer,col_indexer] = value instead

See the caveats in the documentation: https://pandas.pydata.org/pandas-docs/stable/user_guide/indexing.html#returning-a-view-versus-a-copy

```
df_50['Charged'] = df_50['SMILES'].apply(is_charged)
```

```
[8]:
```

	SMILES	Activity	Score	\
5	CNCC1=NC2=C(C=C(C=C2)C1)C(=N1)C3=CC=CN3	Inactive	0.0	
6	CCSC(=NC1=CC=C(C=C1)C(F)(F)F)N.C1	Inactive	0.0	
8	CC1=CC=C(C=C1)S(=O)(=O)N2CCN(CC2)C3=NC(=NC4=CC...	Inactive	0.0	
9	CC1=CC=C(C=C1)S(=O)(=O)N2CCN(CC2)C3=NC(=NC4=CC...	Inactive	0.0	
10	C1CN(CCN1C2=NC(=NC3=CC=CC=C32)C4=CC=CS4)S(=O)(...	Inactive	0.0	
...	
342068	CC(=O)NC1=CC=C(C=C1)OCC2=C(C=CC(=C2)CN(CC3=CC=...	Inactive	0.0	
342069	CC(=O)NC1=CC=C(C=C1)OCC2=C(C=CC(=C2)CN(CC3=CC=...	Inactive	0.0	
342070	CC(=O)NC1=CC=C(C=C1)OCC2=C(C=CC(=C2)CN(CC3=CC=...	Inactive	0.0	
342071	CC(=O)NC1=CC=C(C=C1)C(=O)N(CC2=CC=CC=C2)CC3=CC...	Inactive	0.0	
342072	CC(=O)NC1=CC=C(C=C1)OCC2=C(C=CC(=C2)CN(CC3=CC=...	Inactive	0.0	

	Potency	Efficacy	Charged
5	0.0	0.0	False
6	0.0	0.0	False
8	0.0	0.0	False
9	0.0	0.0	False
10	0.0	0.0	False
...
342068	0.0	0.0	False
342069	0.0	0.0	False
342070	0.0	0.0	False
342071	0.0	0.0	False
342072	0.0	0.0	False

[322199 rows x 6 columns]

```
[9]: # Picking all "Active" molecules from the dataset
active_df = uncharged[uncharged['Activity'] == 'Active']
active_df.info()

# Picking all "Inactive" molecules from the dataset
inactive_df = uncharged[uncharged['Activity'] == 'Inactive']
inactive_df.info()

# Randomly sample from inactive_df to match the size of active_df
inactive_sampled = inactive_df.sample(n=len(active_df), random_state=42)

# Combine the active and sampled inactive molecules
balanced_df = pd.concat([active_df, inactive_sampled])

# Shuffle the combined dataset
balanced_df = balanced_df.sample(frac=1, random_state=42).reset_index(drop=True)

balanced_df.info()
```

```
<class 'pandas.core.frame.DataFrame'>
```

```
Index: 6273 entries, 13 to 341825
```

```
Data columns (total 6 columns):
```

#	Column	Non-Null Count	Dtype
0	SMILES	6273 non-null	object
1	Activity	6273 non-null	object
2	Score	6273 non-null	float64
3	Potency	6273 non-null	float64
4	Efficacy	6273 non-null	float64
5	Charged	6273 non-null	bool

```
dtypes: bool(1), float64(3), object(2)
```

```
memory usage: 300.2+ KB
```

```
<class 'pandas.core.frame.DataFrame'>
```

```
Index: 304069 entries, 5 to 342072
```

```
Data columns (total 6 columns):
```

#	Column	Non-Null Count	Dtype
0	SMILES	304069 non-null	object
1	Activity	304069 non-null	object
2	Score	304069 non-null	float64
3	Potency	304069 non-null	float64
4	Efficacy	304069 non-null	float64
5	Charged	304069 non-null	bool

```
dtypes: bool(1), float64(3), object(2)
```

```
memory usage: 14.2+ MB
```

```
<class 'pandas.core.frame.DataFrame'>
```

```

RangeIndex: 12546 entries, 0 to 12545
Data columns (total 6 columns):
#   Column      Non-Null Count  Dtype
---  -
0   SMILES      12546 non-null  object
1   Activity    12546 non-null  object
2   Score       12546 non-null  float64
3   Potency     12546 non-null  float64
4   Efficacy    12546 non-null  float64
5   Charged     12546 non-null  bool
dtypes: bool(1), float64(3), object(2)
memory usage: 502.5+ KB

```

```
[10]: filtered_df = balanced_df
      filtered_df
```

```

[10]:
      SMILES  Activity  Score \
0      CC1=C(C=CC=C1Br)NC(=O)C2=C(C=CS2)N3C=CC=C3  Active    82.0
1      CCCCC(C(C)CC(=O)NC1CCCCC1)C(=O)O  Active    43.0
2      CC1=CC=C(C=C1)S(=O)(=O)NC2=NN3C(C=C(NC3=N2)C)C...  Active    41.0
3      CC1=CC(=O)OC2=C1C=C(C=C2)OCC(=O)NC3=CC=CC(=C3)...  Inactive    0.0
4      CC1=CC(=C(N1C)C)C(=O)COC(=O)C23CC4CC(C2)CC(C4)...  Inactive    0.0
...
12541  C1CN(CCN1C(=O)C2=CC=CC=C2CC3=CC=CC=C3)S(=O)(=O...  Inactive    0.0
12542      C1=CC=C(C=C1)OC2=NC=NC(=C2)N3C=NC=N3  Active    64.0
12543  CC1=C(C(=CC=C1)N2CCN(CC2)C3=NC4=CC=CC=C4C(=O)N...  Active    42.0
12544  CCC(C)NC(=O)CSC1=NC2=CC=CC=C2C3=NC(C(=O)N31)C4...  Active    42.0
12545      CC(C)C1=CC=C(C=C1)S(=O)(=O)NC2CCCC2  Inactive    0.0

      Potency  Efficacy  Charged
0      8.9125  140.7280   False
1     12.5893  136.6590   False
2     22.3872  166.6580   False
3      0.0000   0.0000   False
4      0.0000   0.0000   False
...
12541   0.0000   0.0000   False
12542   2.8184   74.9734   False
12543  17.7828  126.5240   False
12544  15.8489  139.3040   False
12545   0.0000   0.0000   False

[12546 rows x 6 columns]

```

0.4 Parameter setting

```
[11]: '''
      scan through all the molecules to obtain unique atom types
      '''
      smiles = filtered_df['SMILES'].tolist()
      search_elements=[]
      for smile in smiles:
          mol = Chem.MolFromSmiles(smile)
          atoms = list(set([atom.GetSymbol() for atom in mol.GetAtoms()]))
          search_elements += atoms
          search_elements = list(set(search_elements))
      search_elements.append("H")
      print(search_elements)

['C', 'F', 'N', 'I', 'O', 'P', 'Br', 'B', 'S', 'Cl', 'As', 'H']

[12]: '''
      Setting up the atom mapping and bond mapping.
      Code adopted from https://keras.io/examples/generative/molecule\_generation/
      '''
      SMILE_CHARSET = str(search_elements)
      bond_mapping = {"SINGLE": 0, "DOUBLE": 1, "TRIPLE": 2, "AROMATIC": 3}
      bond_mapping.update(
          {0: BondType.SINGLE, 1: BondType.DOUBLE, 2: BondType.TRIPLE, 3: BondType.
           ↪AROMATIC}
      )
      SMILE_CHARSET = ast.literal_eval(SMILE_CHARSET)

      MAX_MOLSIZE = max(filtered_df['SMILES'].str.len())
      SMILE_to_index = dict((c, i) for i, c in enumerate(SMILE_CHARSET))
      index_to_SMILE = dict((i, c) for i, c in enumerate(SMILE_CHARSET))
      atom_mapping = dict(SMILE_to_index)
      atom_mapping.update(index_to_SMILE)
      print(atom_mapping)
      print("Max molecule size: {}".format(MAX_MOLSIZE))
      print("Character set Length: {}".format(len(SMILE_CHARSET)))

{'C': 0, 'F': 1, 'N': 2, 'I': 3, 'O': 4, 'P': 5, 'Br': 6, 'B': 7, 'S': 8, 'Cl':
9, 'As': 10, 'H': 11, 0: 'C', 1: 'F', 2: 'N', 3: 'I', 4: 'O', 5: 'P', 6: 'Br',
7: 'B', 8: 'S', 9: 'Cl', 10: 'As', 11: 'H'}
Max molecule size: 117
Character set Length: 12
```

0.5 Hyperparameters

```
[13]: '''  
      Defining the Hyperparameters of the model  
      '''  
  
      NUM_ATOMS = 50 #Max number of atoms  
      ATOM_DIM = len(SMILE_CHARSET) # Number of atom types  
      BOND_DIM = 5 # Number of bond types
```

0.6 Molecule featurization

```
[14]: '''  
      Defining functions to convert smiles string into node graph and recover  
      ↪ molecule structure from it.  
      Code referenced from: https://keras.io/examples/generative/molecule\_generation/  
      '''  
  
      def smiles_to_graph(smiles):  
          '''  
          Reference: https://keras.io/examples/generative/wgan-graphs/  
          '''  
          # Converts SMILES to molecule object  
          molecule = Chem.MolFromSmiles(smiles)  
          #molecule = Chem.AddHs(molecule)  
          # Initialize adjacency and feature tensor  
          adjacency = np.zeros((BOND_DIM, NUM_ATOMS, NUM_ATOMS), "float32")  
          features = np.zeros((NUM_ATOMS, ATOM_DIM), "float32")  
  
          # loop over each atom in molecule  
          for atom in molecule.GetAtoms():  
              i = atom.GetIdx()  
              atom_type = atom_mapping[atom.GetSymbol()]  
              features[i] = np.eye(ATOM_DIM)[atom_type]  
              # loop over one-hop neighbors  
              for neighbor in atom.GetNeighbors():  
                  j = neighbor.GetIdx()  
                  bond = molecule.GetBondBetweenAtoms(i, j)  
                  bond_type_idx = bond_mapping[bond.GetBondType().name]  
                  adjacency[bond_type_idx, [i, j], [j, i]] = 1  
  
              # Where no bond, add 1 to last channel (indicating "non-bond")  
              # Notice: channels-first  
              adjacency[-1, np.sum(adjacency, axis=0) == 0] = 1  
  
          # Where no atom, add 1 to last column (indicating "non-atom")
```



```

features[np.where(np.sum(features, axis=1) == 0)[0], -1] = 1

return adjacency, features

def graph_to_molecule(adjacency, features):
    # RWMol is a molecule object intended to be edited
    molecule = Chem.RWMol()
    # Remove "no atoms" & atoms with no bonds
    keep_idx = np.where(
        (np.argmax(features, axis=1) != ATOM_DIM - 1)
        & (np.sum(adjacency[:-1], axis=(0, 1)) > 0))[0]

    features = features[keep_idx]
    adjacency = adjacency[:, keep_idx][:, :, keep_idx]

    # Add atoms to molecule
    for atom_type_idx in np.argmax(features, axis=1):
        atom = Chem.Atom(atom_mapping[atom_type_idx])
        _ = molecule.AddAtom(atom)

    added_bonds = set()
    (bonds_ij, atoms_i, atoms_j) = np.where(np.triu(adjacency) == 1)
    for (bond_ij, atom_i, atom_j) in zip(bonds_ij, atoms_i, atoms_j):
        if atom_i == atom_j or bond_ij == BOND_DIM - 1:
            continue
        bond_type = bond_mapping.get(bond_ij, None)
        if (atom_i, atom_j) in added_bonds or (atom_j, atom_i) in added_bonds:
            continue
        molecule.AddBond(int(atom_i), int(atom_j), bond_type)
        added_bonds.add((atom_i, atom_j))

    # Sanitize without Kekulization
    try:
        Chem.SanitizeMol(molecule, sanitizeOps=Chem.SanitizeFlags.SANITIZE_ALL_
↪ Chem.SanitizeFlags.SANITIZE_KEKULIZE)
    except Exception as e:
        print(f"Sanitization failed: {e}")
        return None

    # Add explicit hydrogens
    molecule_with_h = Chem.AddHs(molecule)

    # Fix aromaticity in aromatic rings
    for atom in molecule_with_h.GetAtoms():
        if atom.GetIsAromatic():

```

```

        atom.SetIsAromatic(False) # Clear aromaticity if needed

# Force Kekulization to alternate bond orders in aromatic rings
try:
    Chem.Kekulize(molecule_with_h, clearAromaticFlags=True)
except Chem.KekulizeException as e:
    print(f"Kekulization failed: {e}")
    return molecule_with_h # Return molecule without Kekulé bonds

return molecule_with_h

```

0.7 Building model

```

[15]: '''
    Defining GCN
    Reference: https://keras.io/examples/generative/wgan-graphs/
    The Encoder takes as input a molecule's graph adjacency matrix and feature_
    ↪matrix.
    '''
class RelationalGraphConvLayer(keras.layers.Layer):
    def __init__(
        self,
        units=128,
        activation="relu",
        use_bias=False,
        kernel_initializer="glorot_uniform",
        bias_initializer="zeros",
        kernel_regularizer=None,
        bias_regularizer=None,
        **kwargs
    ):
        super().__init__(**kwargs)

        self.units = units
        self.activation = keras.activations.get(activation)
        self.use_bias = use_bias
        self.kernel_initializer = keras.initializers.get(kernel_initializer)
        self.bias_initializer = keras.initializers.get(bias_initializer)
        self.kernel_regularizer = keras.regularizers.get(kernel_regularizer)
        self.bias_regularizer = keras.regularizers.get(bias_regularizer)

    def build(self, input_shape):
        bond_dim = input_shape[0][1]
        atom_dim = input_shape[1][2]

        self.kernel = self.add_weight(
            shape=(bond_dim, atom_dim, self.units),

```

```

        initializer=self.kernel_initializer,
        regularizer=self.kernel_regularizer,
        trainable=True,
        name="W",
        dtype=tf.float32,
    )

    if self.use_bias:
        self.bias = self.add_weight(
            shape=(bond_dim, 1, self.units),
            initializer=self.bias_initializer,
            regularizer=self.bias_regularizer,
            trainable=True,
            name="b",
            dtype=tf.float32,
        )

    self.built = True

    def call(self, inputs, training=False):
        adjacency, features = inputs
        # Aggregate information from neighbors
        x = tf.matmul(adjacency, features[:, None, :, :])
        # Apply linear transformation
        x = tf.matmul(x, self.kernel)
        if self.use_bias:
            x += self.bias
        # Reduce bond types dim
        x_reduced = tf.reduce_sum(x, axis=1)
        # Apply non-linear transformation
        return self.activation(x_reduced)

```

0.8 Build the Encoder and Decoder

```

[16]: '''
    defining function to build encoder and decoder.
    Code adopted and modified from https://keras.io/examples/generative/
    ↪ molecule_generation/
    '''

    def get_encoder(gconv_units, latent_dim, adjacency_shape, feature_shape,
    ↪ dense_units, dropout_rate, regularizer=None):
        adjacency = keras.layers.Input(shape=adjacency_shape,
    ↪ name="adjacency_input")
        features = keras.layers.Input(shape=feature_shape, name="feature_input")
        scores = keras.layers.Input(shape=(1,), name="score_input") # Conditional
    ↪ input (scalar)

```

```

# Graph convolution layers
features_transformed = features
for units in gconv_units:
    features_transformed = RelationalGraphConvLayer(units)(
        adjacency, features_transformed
    )

# Reduce 2D representation to 1D
x = keras.layers.GlobalAveragePooling1D()(features_transformed)

# Concatenate the score (condition) to the reduced graph representation
x = keras.layers.Concatenate()([x, scores])

# Fully connected layers
for units in dense_units:
    x = layers.Dense(units, activation="relu",
        ↪kernel_regularizer=regularizer)(x)
    x = layers.Dropout(dropout_rate)(x)

# Latent space
z_mean = layers.Dense(latent_dim, name="z_mean")(x)
z_log_var = layers.Dense(latent_dim, name="z_log_var")(x)

# Create encoder model
encoder = keras.Model(inputs=[adjacency, features, scores],
    ↪outputs=[z_mean, z_log_var], name="encoder")
encoder.summary()
return encoder

class SymmetrizeLayer(layers.Layer):
    def call(self, x):
        return (x + tf.transpose(x, (0, 1, 3, 2))) / 2

def get_decoder(dense_units, latent_dim, adjacency_shape, feature_shape,
    ↪dropout_rate, regularizer=None):
    latent_input = keras.Input(shape=(latent_dim,), name="latent_input")
    scores = keras.Input(shape=(1,), name="score_input") # Conditional input
    ↪(scalar)

    # Concatenate latent input with the conditional score
    x = keras.layers.Concatenate()([latent_input, scores])

    # Dense layers
    for units in dense_units:

```

```

        x = keras.layers.Dense(units, activation="tanh",
kernel_regularizer=regularizer)(x)
        x = keras.layers.Dropout(dropout_rate)(x)

        # Adjacency reconstruction
        adj_output = keras.layers.Dense(tf.math.reduce_prod(adjacency_shape).
numpy().astype(int))(x)
        adj_output = keras.layers.Reshape(adjacency_shape)(adj_output)
        adj_output = SymmetrizeLayer()(adj_output)
        adj_output = keras.layers.Softmax(axis=1)(adj_output)

        # Feature reconstruction
        feat_output = keras.layers.Dense(tf.math.reduce_prod(feature_shape).numpy().
astype(int))(x)
        feat_output = keras.layers.Reshape(feature_shape)(feat_output)
        feat_output = keras.layers.Softmax(axis=2)(feat_output)

        # Create decoder model
        decoder = keras.Model(inputs=[latent_input, scores], outputs=[adj_output,
feat_output], name="decoder")
        decoder.summary()
        return decoder

```

0.9 Build the VAE

```

[17]: '''
defining the VAE
Code adopted and modified from https://keras.io/examples/generative/
molecule_generation/
'''

class VAE(keras.Model):
    def __init__(self, encoder, decoder, beta=1.0, **kwargs):
        super(VAE, self).__init__(**kwargs)
        self.encoder = encoder
        self.decoder = decoder
        self.beta = beta

    def call(self, inputs):
        adjacency, features, scores = inputs
        z_mean, z_log_var = self.encoder([adjacency, features, scores])
        z = self.reparameterize(z_mean, z_log_var)
        return self.decoder([z, scores])

    def sampling(self, args):
        """
        Reparameterization trick: Sample from a Gaussian distribution using

```

```

        z = z_mean + epsilon * exp(z_log_var / 2), where epsilon is sampled
        ↪ from  $N(0, 1)$ .
        """
        z_mean, z_log_var = args
        batch = tf.shape(z_mean)[0]
        dim = tf.shape(z_mean)[1]
        epsilon = tf.keras.backend.random_normal(shape=(batch, dim)) #
        ↪ Standard normal noise
        return z_mean + tf.exp(0.5 * z_log_var) * epsilon

```

0.10 Model training

```

[18]: '''
        splitting the dataset into training and testing
        '''

train, test = train_test_split(filtered_df, test_size=0.2, random_state=42)
train_df, val_df = train_test_split(train, test_size=0.2, random_state=42)
train_df.reset_index(drop=True, inplace=True)
val_df.reset_index(drop=True, inplace=True)
test.reset_index(drop=True, inplace=True)

adj_train, fea_train, score_train = [], [], []
adj_val, fea_val, score_val = [], [], []

for idx in range(len(train_df)):
    adjacency, features = smiles_to_graph(train_df.loc[idx]["SMILES"])
    score = train_df.loc[idx]["Score"]
    adj_train.append(adjacency)
    fea_train.append(features)
    score_train.append(score)

for idx in range(len(val_df)):
    adjacency, features = smiles_to_graph(val_df.loc[idx]["SMILES"])
    score = val_df.loc[idx]["Score"]
    adj_val.append(adjacency)
    fea_val.append(features)
    score_val.append(score)

adj_train = np.array(adj_train)
fea_train = np.array(fea_train)
score_train_ = np.array(score_train).reshape(-1,1)

adj_val = np.array(adj_val)
fea_val = np.array(fea_val)
score_val_ = np.array(score_val).reshape(-1,1)

```

```
[19]: from sklearn.preprocessing import MinMaxScaler

scaler = MinMaxScaler()

score_train_n = scaler.fit_transform(score_train_)
score_val_n = scaler.transform(score_val_)
```

```
[20]: print(adj_train.shape)
print(fea_train.shape)
print(score_train_.shape)
print(adj_val.shape)
print(fea_val.shape)
print(score_val_.shape)
```

```
(8028, 5, 50, 50)
(8028, 50, 12)
(8028, 1)
(2008, 5, 50, 50)
(2008, 50, 12)
(2008, 1)
```

```
[21]: print(np.max(score_train_n))
```

```
0.9999999999999999
```

```
[22]: #Hyperparameters
BATCH_SIZE = 64
EPOCHS = 25
VAE_LR = 3e-4 # changed to 1e-3
LATENT_DIM = 128 # Size of the latent space
```

```
[23]: '''
compiling the VAE
'''

encoder = get_encoder(
    gconv_units=[16],
    adjacency_shape=(BOND_DIM, NUM_ATOMS, NUM_ATOMS),
    feature_shape=(NUM_ATOMS, ATOM_DIM),
    latent_dim=LATENT_DIM,
    dense_units=[256, 512],
    dropout_rate=0,
    regularizer=l1_l2(l1=1e-6, l2=1e-3)
)
decoder = get_decoder(
    dense_units=[128, 256, 512],
    dropout_rate=0.3,
    latent_dim=LATENT_DIM,
```

```

adjacency_shape=(BOND_DIM, NUM_ATOMS, NUM_ATOMS),
feature_shape=(NUM_ATOMS, ATOM_DIM),
regularizer=l1_l2(l1=1e-4, l2=1e-2)
)
vae = VAE(encoder, decoder)

vae.compile(optimizer=keras.optimizers.Adam(learning_rate=VAE_LR))

```

Model: "encoder"

Layer (type)	Output Shape	Param #	Connected to
adjacency_input (InputLayer)	(None, 5, 50, 50)	0	-
feature_input (InputLayer)	(None, 50, 12)	0	-
relational_graph_c... (RelationalGraphCo...)	(None, 50, 16)	960	adjacency_input[... feature_input[0]...
global_average_poo... (GlobalAveragePool...)	(None, 16)	0	relational_graph...
score_input (InputLayer)	(None, 1)	0	-
concatenate (Concatenate)	(None, 17)	0	global_average_p... score_input[0][0]
dense (Dense)	(None, 256)	4,608	concatenate[0][0]
dropout (Dropout)	(None, 256)	0	dense[0][0]
dense_1 (Dense)	(None, 512)	131,584	dropout[0][0]
dropout_1 (Dropout)	(None, 512)	0	dense_1[0][0]
z_mean (Dense)	(None, 128)	65,664	dropout_1[0][0]
z_log_var (Dense)	(None, 128)	65,664	dropout_1[0][0]

Total params: 268,480 (1.02 MB)

Trainable params: 268,480 (1.02 MB)

Non-trainable params: 0 (0.00 B)

Model: "decoder"

Layer (type)	Output Shape	Param #	Connected to
latent_input (InputLayer)	(None, 128)	0	-
score_input (InputLayer)	(None, 1)	0	-
concatenate_1 (Concatenate)	(None, 129)	0	latent_input[0] [...] score_input[0][0]
dense_2 (Dense)	(None, 128)	16,640	concatenate_1[0]...
dropout_2 (Dropout)	(None, 128)	0	dense_2[0][0]
dense_3 (Dense)	(None, 256)	33,024	dropout_2[0][0]
dropout_3 (Dropout)	(None, 256)	0	dense_3[0][0]
dense_4 (Dense)	(None, 512)	131,584	dropout_3[0][0]
dropout_4 (Dropout)	(None, 512)	0	dense_4[0][0]
dense_5 (Dense)	(None, 12500)	6,412,500	dropout_4[0][0]
reshape (Reshape)	(None, 5, 50, 50)	0	dense_5[0][0]
dense_6 (Dense)	(None, 600)	307,800	dropout_4[0][0]
symmetrize_layer (SymmetrizeLayer)	(None, 5, 50, 50)	0	reshape[0][0]
reshape_1 (Reshape)	(None, 50, 12)	0	dense_6[0][0]
softmax (Softmax)	(None, 5, 50, 50)	0	symmetrize_layer...
softmax_1 (Softmax)	(None, 50, 12)	0	reshape_1[0][0]

Total params: 6,901,548 (26.33 MB)

Trainable params: 6,901,548 (26.33 MB)

Non-trainable params: 0 (0.00 B)

```
[24]: val_loss_list = []
      train_loss_list = []
      kl_theshold = 1.0

[25]: train_dataset = tf.data.Dataset.from_tensor_slices((adj_train, fea_train, ↵
      ↵score_train_)).batch(BATCH_SIZE)
      val_dataset = tf.data.Dataset.from_tensor_slices((adj_val, fea_val, ↵
      ↵score_val_)).batch(BATCH_SIZE)

[26]: for epoch in range(EPOCHS):
      print(f"Epoch {epoch + 1}/{EPOCHS}")
      if epoch < 10:
          beta = 0.05
      else:
          beta = epoch*0.01
      # Training Loop
      train_loss = 0
      for (adjacency, features, scores) in train_dataset:
          with tf.GradientTape() as tape:
              # Forward pass
              z_mean, z_log_var = vae.encoder([adjacency, features, scores])
              z = vae.sampling([z_mean, z_log_var])
              adj_reconstruction, feature_reconstruction = vae.decoder([z, ↵
              ↵scores])

              # Compute losses
              adj_loss = tf.reduce_mean(
                  tf.reduce_sum(keras.losses.binary_crossentropy(adjacency, ↵
              ↵adj_reconstruction), axis=(1, 2))
              )
              feat_loss = tf.reduce_mean(
                  tf.reduce_sum(keras.losses.categorical_crossentropy(features, ↵
              ↵feature_reconstruction), axis=1)
              )
              reconstruction_loss = adj_loss + feat_loss
              kl_loss = -0.5 * tf.reduce_mean(
                  tf.reduce_sum(1 + z_log_var - tf.square(z_mean) - tf.
              ↵exp(z_log_var), axis=1)
              )
```

```

        total_loss = reconstruction_loss + beta * kl_loss

        # Backpropagation
        grads = tape.gradient(total_loss, vae.trainable_weights)
        vae.optimizer.apply_gradients(zip(grads, vae.trainable_weights))

        train_loss += total_loss

train_loss /= len(train_dataset)
train_loss_list.append(train_loss)

print(f"Train Loss: {train_loss.numpy()}, KL Loss: {kl_loss.numpy()},  

↪Reconstruction Loss: {reconstruction_loss.numpy()}")

# Validation Loop
val_loss = 0
for (val_adjacency, val_features, val_scores) in val_dataset:
    # Forward pass
    z_mean, z_log_var = vae.encoder([val_adjacency, val_features,  

↪val_scores])
    z = vae.sampling([z_mean, z_log_var])
    val_adj_reconstruction, val_feat_reconstruction = vae.decoder([z,  

↪val_scores])

    # Compute losses
    val_adj_loss = tf.reduce_mean(
        tf.reduce_sum(keras.losses.binary_crossentropy(val_adjacency,  

↪val_adj_reconstruction), axis=(1, 2))
    )
    val_feat_loss = tf.reduce_mean(
        tf.reduce_sum(keras.losses.categorical_crossentropy(val_features,  

↪val_feat_reconstruction), axis=1)
    )
    val_reconstruction_loss = val_adj_loss + val_feat_loss
    val_kl_loss = -0.5 * tf.reduce_mean(
        tf.reduce_sum(1 + z_log_var - tf.square(z_mean) - tf.  

↪exp(z_log_var), axis=1)
    )
    val_total_loss = val_reconstruction_loss + beta * val_kl_loss

    val_loss += val_total_loss

val_loss /= len(val_dataset)
val_loss_list.append(val_loss)

```

```

# Adjust beta if KL loss is very low
if kl_loss < kl_theshold:
    beta = 0.05
print(f"Validation Loss: {val_loss.numpy()}, KL Loss: {val_kl_loss.
numpy()}, Reconstruction Loss: {val_reconstruction_loss.numpy()}")
print('BETA is: ', beta)

```

Epoch 1/25

2024-12-12 17:13:48.636583: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 66.95944213867188, KL Loss: 5.409961223602295, Reconstruction Loss: 36.86543655395508

2024-12-12 17:13:49.281854: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 38.307167053222656, KL Loss: 7.239017486572266, Reconstruction Loss: 35.230045318603516

BETA is: 0.05

Epoch 2/25

2024-12-12 17:13:57.845969: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 38.08542251586914, KL Loss: 7.262415409088135, Reconstruction Loss: 35.81265640258789

2024-12-12 17:13:58.369698: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 37.226993560791016, KL Loss: 8.832179069519043, Reconstruction Loss: 34.20383071899414

BETA is: 0.05

Epoch 3/25

2024-12-12 17:14:07.423950: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 34.377647399902344, KL Loss: 13.722108840942383, Reconstruction Loss: 31.014507293701172

2024-12-12 17:14:07.964127: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 32.02784729003906, KL Loss: 14.46497631072998, Reconstruction Loss: 29.39027976989746

BETA is: 0.05

Epoch 4/25

2024-12-12 17:14:16.276468: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 31.324237823486328, KL Loss: 15.213322639465332, Reconstruction Loss: 29.61969566345215

2024-12-12 17:14:16.802566: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 30.837430953979492, KL Loss: 15.234996795654297, Reconstruction Loss: 27.43654441833496

BETA is: 0.05

Epoch 5/25

2024-12-12 17:14:24.980889: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 30.58317756652832, KL Loss: 14.706596374511719, Reconstruction Loss: 29.678884506225586

2024-12-12 17:14:25.507104: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 30.382816314697266, KL Loss: 16.113767623901367, Reconstruction Loss: 27.395017623901367

BETA is: 0.05

Epoch 6/25

2024-12-12 17:14:33.694724: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 30.195568084716797, KL Loss: 15.516021728515625, Reconstruction Loss: 28.656808853149414

2024-12-12 17:14:34.217531: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 30.03752899169922, KL Loss: 16.778310775756836, Reconstruction Loss: 27.231407165527344

BETA is: 0.05

Epoch 7/25

2024-12-12 17:14:42.781334: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 30.25861930847168, KL Loss: 16.683490753173828, Reconstruction Loss: 28.643312454223633

2024-12-12 17:14:43.308068: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.948104858398438, KL Loss: 17.50771141052246, Reconstruction Loss: 26.600330352783203

BETA is: 0.05

Epoch 8/25

2024-12-12 17:14:51.590812: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.655874252319336, KL Loss: 15.88936710357666, Reconstruction Loss: 28.086673736572266

2024-12-12 17:14:52.134219: W tensorflow/core/framework/local_rendevvous.cc:404]
Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.685688018798828, KL Loss: 16.187108993530273, Reconstruction Loss: 26.570858001708984

BETA is: 0.05

Epoch 9/25

2024-12-12 17:15:00.586432: W tensorflow/core/framework/local_rendevvous.cc:404]
Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.42094612121582, KL Loss: 15.739115715026855, Reconstruction Loss: 28.297420501708984

2024-12-12 17:15:01.104570: W tensorflow/core/framework/local_rendevvous.cc:404]
Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.49125862121582, KL Loss: 16.52348518371582, Reconstruction Loss: 26.6302433013916

BETA is: 0.05

Epoch 10/25

2024-12-12 17:15:09.567166: W tensorflow/core/framework/local_rendevvous.cc:404]
Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.231225967407227, KL Loss: 15.30255126953125, Reconstruction Loss: 27.82144546508789

2024-12-12 17:15:10.087751: W tensorflow/core/framework/local_rendevvous.cc:404]
Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.279537200927734, KL Loss: 16.1846981048584, Reconstruction Loss: 26.073829650878906

BETA is: 0.05

Epoch 11/25

2024-12-12 17:15:18.996780: W tensorflow/core/framework/local_rendevvous.cc:404]
Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.884469985961914, KL Loss: 12.192437171936035, Reconstruction Loss: 28.23736000061035

2024-12-12 17:15:19.550752: W tensorflow/core/framework/local_rendevvous.cc:404]
Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.96446990966797, KL Loss: 13.384326934814453, Reconstruction Loss: 26.216266632080078

BETA is: 0.1

Epoch 12/25

2024-12-12 17:15:27.985540: W tensorflow/core/framework/local_rendevvous.cc:404]
Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.867341995239258, KL Loss: 10.197785377502441, Reconstruction Loss: 28.79987907409668

2024-12-12 17:15:28.528380: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.928882598876953, KL Loss: 10.827570915222168, Reconstruction Loss: 28.151025772094727

BETA is: 0.11

Epoch 13/25

2024-12-12 17:15:36.862064: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.78401756286621, KL Loss: 9.285300254821777, Reconstruction Loss: 28.00973892211914

2024-12-12 17:15:37.384640: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.772369384765625, KL Loss: 9.544482231140137, Reconstruction Loss: 26.41633415222168

BETA is: 0.12

Epoch 14/25

2024-12-12 17:15:45.740881: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 30.474695205688477, KL Loss: 9.579269409179688, Reconstruction Loss: 28.297019958496094

2024-12-12 17:15:46.264128: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.842233657836914, KL Loss: 9.950072288513184, Reconstruction Loss: 26.075977325439453

BETA is: 0.13

Epoch 15/25

2024-12-12 17:15:54.413123: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.662763595581055, KL Loss: 7.968265533447266, Reconstruction Loss: 27.789501190185547

2024-12-12 17:15:54.943067: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 30.198326110839844, KL Loss: 8.257922172546387, Reconstruction Loss: 26.271183013916016

BETA is: 0.14

Epoch 16/25

2024-12-12 17:16:03.272911: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 30.350461959838867, KL Loss: 8.190351486206055, Reconstruction Loss: 28.928070068359375

2024-12-12 17:16:03.803879: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.776681900024414, KL Loss: 8.088484764099121, Reconstruction Loss: 26.190954208374023

BETA is: 0.15

Epoch 17/25

2024-12-12 17:16:12.172904: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 30.108983993530273, KL Loss: 7.2249579429626465, Reconstruction Loss: 27.830188751220703

2024-12-12 17:16:12.691355: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.608184814453125, KL Loss: 7.540590763092041, Reconstruction Loss: 26.069196701049805

BETA is: 0.16

Epoch 18/25

2024-12-12 17:16:20.934304: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.48594856262207, KL Loss: 6.611556053161621, Reconstruction Loss: 27.594058990478516

2024-12-12 17:16:21.456986: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.519359588623047, KL Loss: 6.855264186859131, Reconstruction Loss: 27.162221908569336

BETA is: 0.17

Epoch 19/25

2024-12-12 17:16:29.780382: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.709335327148438, KL Loss: 6.697078704833984, Reconstruction Loss: 27.527141571044922

2024-12-12 17:16:30.301756: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.652938842773438, KL Loss: 6.887571811676025, Reconstruction Loss: 26.06425666809082

BETA is: 0.18

Epoch 20/25

2024-12-12 17:16:38.987568: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 30.029027938842773, KL Loss: 6.425731182098389, Reconstruction Loss: 27.666391372680664

2024-12-12 17:16:39.494257: W tensorflow/core/framework/local_rendevvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.664268493652344, KL Loss: 6.584661960601807, Reconstruction Loss: 26.06829261779785

BETA is: 0.19

Epoch 21/25

2024-12-12 17:16:47.847076: W tensorflow/core/framework/local_rendevvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.68038558959961, KL Loss: 6.201575756072998, Reconstruction Loss: 27.925024032592773

2024-12-12 17:16:48.368611: W tensorflow/core/framework/local_rendevvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.514310836791992, KL Loss: 6.663267612457275, Reconstruction Loss: 25.973363876342773

BETA is: 0.2

Epoch 22/25

2024-12-12 17:16:56.768037: W tensorflow/core/framework/local_rendevvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.492443084716797, KL Loss: 6.026613712310791, Reconstruction Loss: 27.56798553466797

2024-12-12 17:16:57.305910: W tensorflow/core/framework/local_rendevvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.594175338745117, KL Loss: 6.642364501953125, Reconstruction Loss: 25.876943588256836

BETA is: 0.21

Epoch 23/25

2024-12-12 17:17:05.655918: W tensorflow/core/framework/local_rendevvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.715835571289062, KL Loss: 9.752555847167969, Reconstruction Loss: 27.58233642578125

2024-12-12 17:17:06.196618: W tensorflow/core/framework/local_rendevvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 30.531145095825195, KL Loss: 12.062098503112793, Reconstruction Loss: 25.871355056762695

BETA is: 0.22

Epoch 24/25

2024-12-12 17:17:14.603124: W tensorflow/core/framework/local_rendevvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.64225196838379, KL Loss: 5.31318998336792, Reconstruction Loss: 27.660545349121094

2024-12-12 17:17:15.133052: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Validation Loss: 29.5415096282959, KL Loss: 5.519102096557617, Reconstruction Loss: 26.132801055908203

BETA is: 0.23

Epoch 25/25

2024-12-12 17:17:23.421472: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

Train Loss: 29.305187225341797, KL Loss: 5.078959941864014, Reconstruction Loss: 27.42971420288086

Validation Loss: 29.36880874633789, KL Loss: 5.2280707359313965, Reconstruction Loss: 25.697792053222656

BETA is: 0.24

2024-12-12 17:17:23.960953: W tensorflow/core/framework/local_rendezvous.cc:404] Local rendezvous is aborting with status: OUT_OF_RANGE: End of sequence

```
[27]: '''  
Checking the model's ability to reconstruct a molecule from the training dataset  
'''
```

```
i=10  
adjacency_check, features_check = smiles_to_graph(train_df.loc[i]["SMILES"])  
score_check = [train_df.loc[i]["Score"]]  
molobj = Chem.MolFromSmiles(train_df.loc[i]["SMILES"])  
adj0 = np.expand_dims(adjacency_check,axis=0)  
feature0 = np.expand_dims(features_check,axis=0)  
score0 = np.expand_dims(score_check,axis=0)  
print(adj0.shape)  
print(feature0.shape)  
print(score0.shape)
```

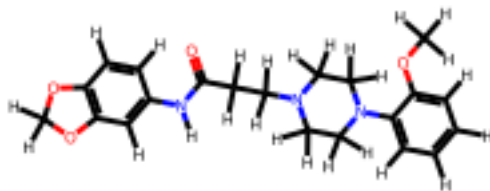
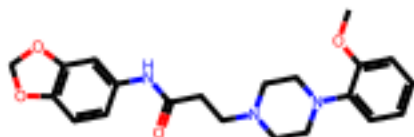
(1, 5, 50, 50)

(1, 50, 12)

(1, 1)

```
[28]: mole_pred = graph_to_molecule(adj0[0], feature0[0])  
Draw.MolsToGridImage([molobj,mole_pred], molsPerRow=2,)
```

[28]:



```
[29]: plt.figure(figsize=(10, 6))
plt.plot(range(EPOCHS), train_loss_list, label='Training Loss', marker='o')
plt.plot(range(EPOCHS), val_loss_list, label='Validation Loss', marker='s')

# Add title and labels
plt.title('Training and Validation Loss Over Epochs', fontsize=16)
plt.xlabel('Epochs', fontsize=14)
plt.ylabel('Loss', fontsize=14)
plt.legend()
plt.show()
```



0.11 Visualize latent space

```
[30]: adj_test, fea_test, score_test = [], [], []

for idx in range(len(test)):
    adjacency, features = smiles_to_graph(test.loc[idx]["SMILES"])
    score = test.loc[idx]["Score"]
    adj_test.append(adjacency)
    fea_test.append(features)
    score_test.append(score)

adj_test = np.array(adj_test)
fea_test = np.array(fea_test)
score_test_ = np.array(score_test).reshape(-1,1)

score_test_n = scaler.transform(score_test_)

[31]: ls_train = vae.encoder.predict([adj_train, fea_train, score_train_])
ls_test = vae.encoder.predict([adj_test, fea_test, score_test_])
```

```
251/251          0s 1ms/step
79/79           0s 949us/step
```

```
[32]: ls_train_ = np.array(ls_train)
ls_test_ = np.array(ls_test)
```

```
[33]: z_mean, _ = vae.encoder.predict([adj_test, fea_test, score_test_])
```

79/79 0s 940us/step

```
[34]: latent_noise = np.random.normal(scale=0.1, size=z_mean.shape) # Adjust scale
      ↪ as needed
adj_pred, feature_pred = vae.decoder.predict([z_mean, score_test_])
print("Shape of adj_pred:", adj_pred.shape)
print("Shape of feature_pred:", feature_pred.shape)

# Reconstruct molecules
gen_molecules = [
    graph_to_molecule(adj_pred[i], feature_pred[i])
    for i in range(adj_pred.shape[0])
]
```

79/79 0s 4ms/step
Shape of adj_pred: (2510, 5, 50, 50)
Shape of feature_pred: (2510, 50, 12)

```
[35]: from scipy.stats import pearsonr

# Correlate latent dimensions with molecular scores
correlations = [pearsonr(z_mean[:, i], score_test_.flatten())[0] for i in
      ↪ range(z_mean.shape[1])]
print("Correlations between latent dimensions and scores:", correlations)
```

Correlations between latent dimensions and scores: [-0.7850554335112523,
-0.6306591561123174, 0.7122755575526665, -0.8384445636545532,
-0.03329708578326372, -0.5128290639442652, -0.3503247027277769,
0.7418109621038158, 0.2802030373831089, 0.7067002149616265, -0.5584619795273105,
0.5889388012397376, -0.8340726274211244, -0.4677558525565457,
0.6963222732860327, -0.7603323000244127, 0.32131010068755494,
-0.025614220248301664, -0.35092335340120984, 0.21160725495285485,
-0.5861931198978443, -0.07654051722532812, 0.9091821075312202,
0.28357791502671437, 0.2875836981517368, -0.14522348188021028,
0.3581287029268357, 0.5954473220122992, 0.21280458240518516,
-0.24607783952217777, -0.3427265948240112, -0.866325531669493,
-0.0015810939225799243, -0.5049793164094454, -0.2222107454554603,
-0.03660636535373199, 0.8123520450428507, 0.7805129301284202,
-0.38020402282392707, -0.3575034298355539, -0.6015176190984987,
-0.2671712290984119, 0.7173269356104413, 0.7925537909553199,
-0.1711200109054484, -0.47689979486434786, -0.6228145335431732,
-0.10161235621550897, 0.26871469638137496, 0.4015489560252962,
0.43007015780231955, 0.8316529522907639, 0.28551271612804724,
0.5885021337372118, -0.6034185116847695, 0.7621254941182303, -0.479002061446721,

```

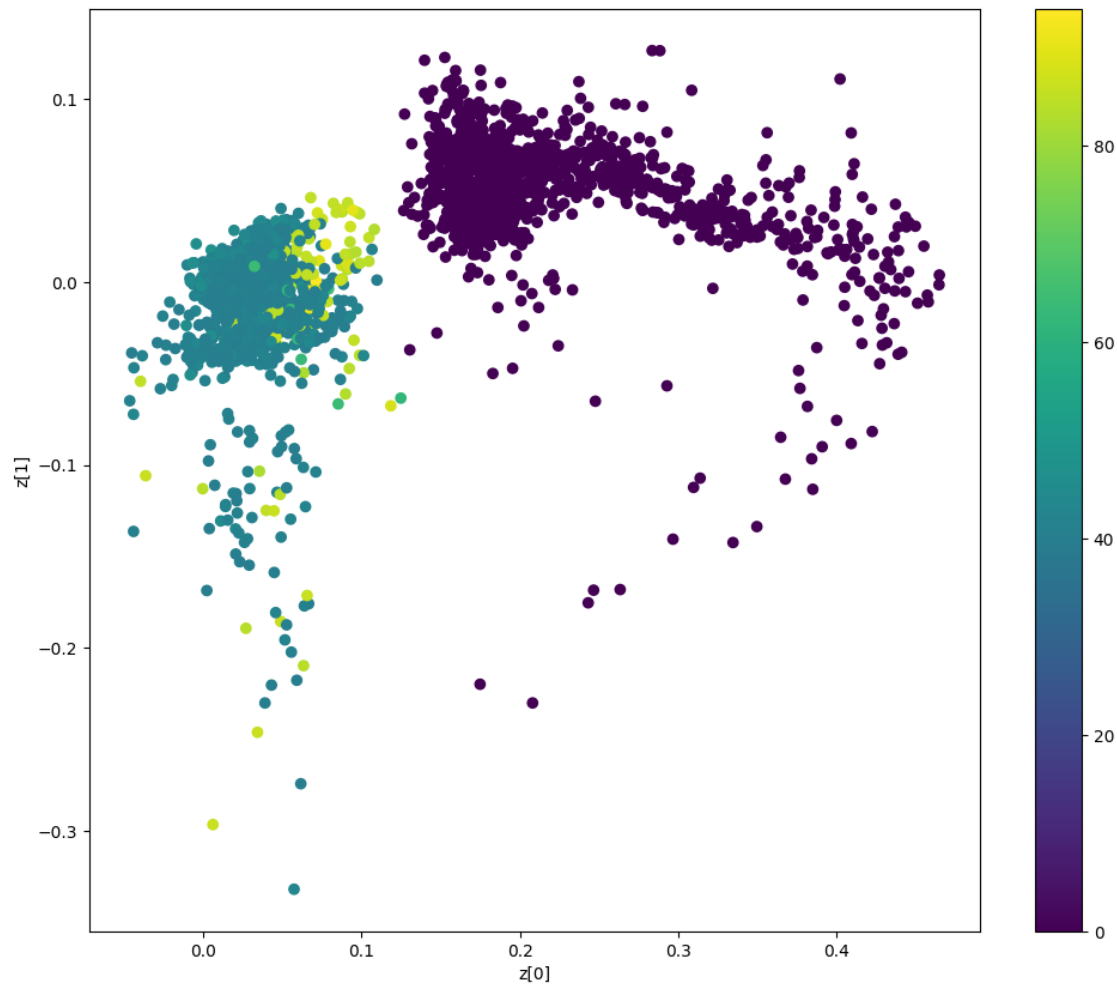
0.7086164492993798, 0.10362830692577474, -0.7154291969187991,
0.3289321510301183, 0.23354268279607698, -0.48967506805129474,
-0.11147307340246222, -0.4152879280556193, 0.2056593068335026,
0.1587648121037009, -0.8834500734164302, 0.29542531971612696,
0.4870743434230065, -0.12363232695242646, -0.013637741457221957,
-0.7754597517650521, -0.7327582691785419, -0.6256190122659977,
-0.714954677197715, 0.7890813489887689, -0.35633482994510884,
0.48801617925020113, 0.7360983673056387, -0.8001255948127852,
0.7061483860019719, 0.3824897644281933, 0.6199756543338328,
-0.22146019425794355, -0.03098114082251182, -0.5862217171707669,
-0.6391098003826348, 0.6026825853282776, -0.11006482487092431,
-0.7789412180675428, -0.46014937111018805, -0.30433131924323065,
-0.4758355433790832, 0.6251182532113478, -0.8814038613010716,
-0.4733470438418641, -0.5292510461568108, -0.74863584678701,
0.23352610233819618, 0.10242448041906899, -0.5309709488749564,
-0.7786801118734887, -0.6781878683546873, -0.15878406641304382,
-0.4466108024473777, 0.17192119083836072, 0.19463701849277398,
0.7019410208600592, -0.6527598279255493, 0.6438038557634551,
-0.7942390856552993, 0.7392068994549414, 0.4816113309242206, 0.5804107934560877,
-0.7167182671675619, 0.6972193487622782, 0.7071969203679852,
0.09713808146575374, 0.2176402763163122, -0.8495411644126821,
-0.4543197179974845, 0.30193883016333073, -0.8500648473261383,
-0.32764039233664466, 0.6982855275392338, 0.319589124112355,
-0.36319398549868454]

```

```

[36]: plt.figure(figsize=(12, 10))
plt.scatter(z_mean[:, 0], z_mean[:, 1], c=score_test_)
plt.colorbar()
plt.xlabel("z[0]")
plt.ylabel("z[1]")
plt.show()

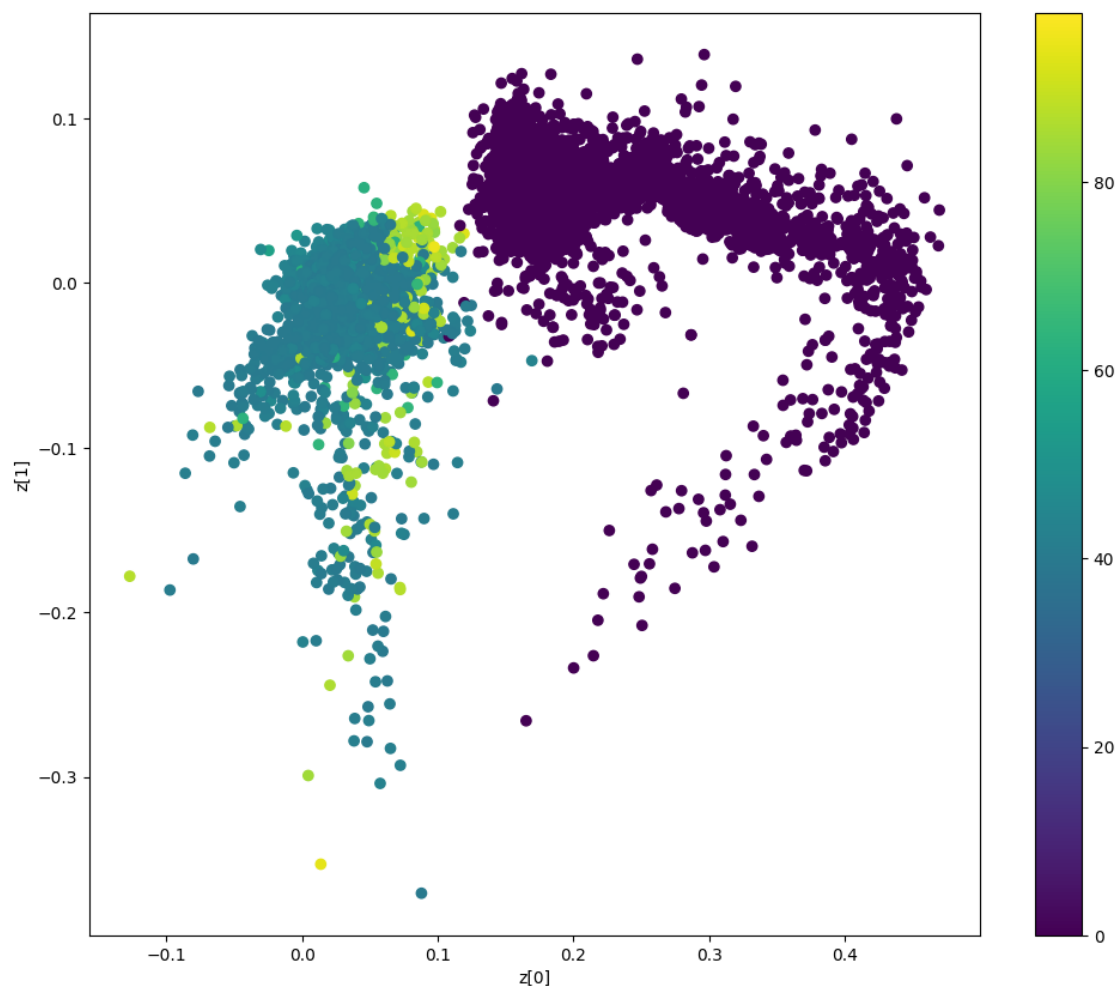
```



```
[37]: z_mean2, _2 = vae.encoder.predict([adj_train, fea_train, score_train_])
```

251/251 0s 912us/step

```
[38]: plt.figure(figsize=(12, 10))
plt.scatter(z_mean2[:, 0], z_mean2[:, 1], c=score_train_)
plt.colorbar()
plt.xlabel("z[0]")
plt.ylabel("z[1]")
plt.show()
```



[]:

0.12 Model Inferencing

We would be inferring our model to predict over random latent space and try to generate 100 new valid molecules.

0.12.1 Generate unique Molecules with the model

```
[39]: def inference(model=vae, batch_size=1000, dim = LATENT_DIM, activity=10):
    z = np.random.normal(size=(batch_size, dim))
    activityarray = (np.zeros(batch_size) + activity).reshape(-1,1)

    reconstruction_adjacency, reconstruction_features = model.decoder.
    ↪predict([z,activityarray])
    # obtain one-hot encoded adjacency tensor
```



```

adjacency = tf.argmax(reconstruction_adjacency, axis=1)
adjacency = tf.one_hot(adjacency, depth=BOND_DIM, axis=1)
# Remove potential self-loops from adjacency
adjacency = tf.linalg.set_diag(adjacency, tf.zeros(tf.shape(adjacency)[:
↪-1]))
# obtain one-hot encoded feature tensor
features = tf.argmax(reconstruction_features, axis=2)
features = tf.one_hot(features, depth=ATOM_DIM, axis=2)

return [
    graph_to_molecule(adjacency[i].numpy(), features[i].numpy())
    for i in range(batch_size)
]

```

```

[40]: gen_mols = inference(batch_size=1000,activity=10)
MolsToGridImage([m for m in gen_mols if m is not None][:1000], molsPerRow=5,
↪subImgSize=(260, 160))

```

32/32 0s 4ms/step

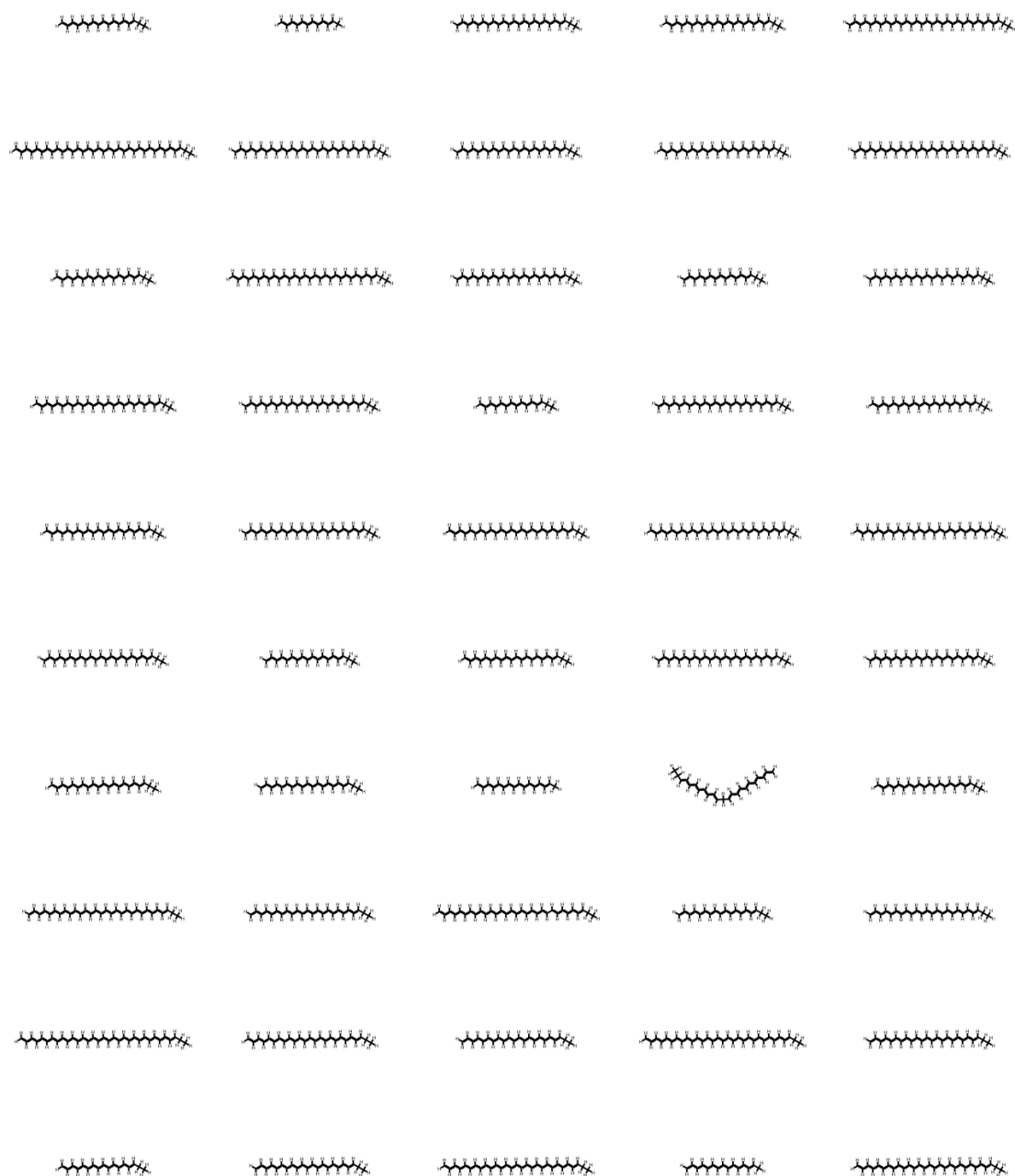
Sanitization failed: Explicit valence for atom # 23 C, 8, is greater than permitted

Sanitization failed: Explicit valence for atom # 2 N, 9, is greater than permitted

Sanitization failed: Explicit valence for atom # 0 S, 101, is greater than permitted

/Users/thinh/Library/Python/3.12/lib/python/site-packages/rdkit/Chem/Draw/IPythonConsole.py:261: UserWarning: Truncating the list of molecules to be displayed to 50. Change the maxMols value to display more.
warnings.warn(

[40]:



[]: