

Setup

```
gpu_info = !nvidia-smi
gpu_info = '\n'.join(gpu_info)
if gpu_info.find('failed') >= 0:
    print('Not connected to a GPU')
else:
    print(gpu_info)
```

Sun Mar 10 19:50:49 2024

| NVIDIA-SMI 535.104.05 | | | | Driver Version: 535.104.05 | | CUDA Version: 12.2 | |
|-----------------------|----------|------|---------------|----------------------------|--------------|--------------------|-------------------|
| GPU | Name | Perf | Persistence-M | Bus-Id | Disp.A | Volatile | Uncorr. ECC |
| Fan | Temp | | Pwr:Usage/Cap | | Memory-Usage | GPU-Util | Compute M. MIG M. |
| 0 | Tesla T4 | | Off | 00000000:00:04.0 | Off | 0 | 0 |
| N/A | 51C | P8 | 10W / 70W | 0MiB / 15360MiB | | 0% | Default N/A |

| Processes: | | | | | | | |
|----------------------------|----|----|-----|------|--------------|------------|--|
| GPU | GI | CI | PID | Type | Process name | GPU Memory | |
| | ID | ID | | | | Usage | |
| No running processes found | | | | | | | |

```
from psutil import virtual_memory
ram_gb = virtual_memory().total / 1e9
print('Your runtime has {:.1f} gigabytes of available RAM\n'.format(ram_gb))
```

```
if ram_gb < 20:
    print('Not using a high-RAM runtime')
else:
    print('You are using a high-RAM runtime!')
```

Your runtime has 13.6 gigabytes of available RAM

Not using a high-RAM runtime



```
!pip install chemprop
!pip install rdkit-pypi # should be included in above after Chemprop v1.6 release
```

```
import chemprop
import pandas as pd
import matplotlib.pyplot as plt
from matplotlib.offsetbox import AnchoredText
from sklearn.metrics import mean_absolute_error, mean_squared_error
from sklearn.decomposition import PCA
```

```
Requirement already satisfied: imagesize in /usr/local/lib/python3.10/dist-packages (from sphinx>=3.1.2->chemprop)
Requirement already satisfied: requests>=2.5.0 in /usr/local/lib/python3.10/dist-packages (from sphinx>=3.1.2->ch
Requirement already satisfied: protobuf>=3.20 in /usr/local/lib/python3.10/dist-packages (from tensorboardX>=2.0-
Requirement already satisfied: filelock in /usr/local/lib/python3.10/dist-packages (from torch>=1.4.0->chemprop)
Requirement already satisfied: typing-extensions in /usr/local/lib/python3.10/dist-packages (from torch>=1.4.0->c
Requirement already satisfied: sympy in /usr/local/lib/python3.10/dist-packages (from torch>=1.4.0->chemprop) (1.
Requirement already satisfied: fsspec in /usr/local/lib/python3.10/dist-packages (from torch>=1.4.0->chemprop) (2
Requirement already satisfied: triton==2.1.0 in /usr/local/lib/python3.10/dist-packages (from torch>=1.4.0->chemp
Collecting typing-inspect>=0.7.1 (from typed-argument-parser>=1.6.1->chemprop)
  Downloading typing_inspect-0.9.0-py3-none-any.whl (8.8 kB)
Collecting docstring-parser>=0.15 (from typed-argument-parser>=1.6.1->chemprop)
  Downloading docstring_parser-0.15-py3-none-any.whl (36 kB)
Requirement already satisfied: MarkupSafe>=2.0 in /usr/local/lib/python3.10/dist-packages (from Jinja2>=3.0->flas
Requirement already satisfied: charset-normalizer<4,>=2 in /usr/local/lib/python3.10/dist-packages (from requests
Requirement already satisfied: idna<4,>=2.5 in /usr/local/lib/python3.10/dist-packages (from requests>=2.5.0->sph
Requirement already satisfied: urllib3<3,>=1.21.1 in /usr/local/lib/python3.10/dist-packages (from requests>=2.5.
Requirement already satisfied: certifi>=2017.4.17 in /usr/local/lib/python3.10/dist-packages (from requests>=2.5.
Collecting mpy-extensions>=0.3.0 (from typing-inspect>=0.7.1->typed-argument-parser>=1.6.1->chemprop)
  Downloading mpy_extensions-1.0.0-py3-none-any.whl (4.7 kB)
Requirement already satisfied: mpmath>=0.19 in /usr/local/lib/python3.10/dist-packages (from sympy->torch>=1.4.0-
Building wheels for collected packages: typed-argument-parser
  Building wheel for typed-argument-parser (setup.py) ... done
  Created wheel for typed-argument-parser: filename=typed_argument_parser-1.9.0-py3-none-any.whl size=25615 sha25
  Stored in directory: /root/.cache/pip/wheels/f0/94/0f/9539f578bed7e1bd423c702e403712f5ee8989f831a71db000
Successfully built typed-argument-parser
Installing collected packages: tensorboardX, rdkit, mpy-extensions, docstring-parser, typing-inspect, typed-argu
Successfully installed chemprop-1.6.1 docstring-parser-0.15 mpy-extensions-1.0.0 pandas-flavor-0.6.0 rdkit-2023.
Collecting rdkit-pypi
  Downloading rdkit_pypi-2022.9.5-cp310-cp310-manylinux_2_17_x86_64.manylinux2014_x86_64.whl (29.4 MB)
29.4/29.4 MB 22.2 MB/s eta 0:00:00
Requirement already satisfied: numpy in /usr/local/lib/python3.10/dist-packages (from rdkit-pypi) (1.25.2)
Requirement already satisfied: Pillow in /usr/local/lib/python3.10/dist-packages (from rdkit-pypi) (9.4.0)
Installing collected packages: rdkit-pypi
Successfully installed rdkit-pypi-2022.9.5
```



```
from IPython.core.interactiveshell import InteractiveShell
InteractiveShell.ast_node_interactivity = "all"
```

```
hiv_df = pd.read_csv("HIV.csv")
hiv_df.head()
```

| | smiles | activity | HIV_active |  |
|---|---|----------|------------|---|
| 0 | CCC1=[O+][Cu-3]2([O+]=C(CC)C1)[O+]=C(CC)CC(CC)... | Cl | 0 |  |
| 1 | C(=Cc1cccc1)C1=[O+][Cu-3]2([O+]=C(C=Cc3cccc3... | Cl | 0 | |
| 2 | CC(=O)N1c2cccc2Sc2c1ccc1cccc21 | Cl | 0 | |
| 3 | Nc1ccc(C=Cc2ccc(N)cc2S(=O)(=O)O)c(S(=O)(=O)O)c1 | Cl | 0 | |
| 4 | O=S(=O)(O)CCS(=O)(=O)O | Cl | 0 | |

Next steps:  [View recommended plots](#)

```
hiv_df.describe()
```

| | HIV_active |  |
|-------|--------------|---|
| count | 41127.000000 |  |
| mean | 0.035086 | |
| std | 0.184001 | |
| min | 0.000000 | |
| 25% | 0.000000 | |
| 50% | 0.000000 | |
| 75% | 0.000000 | |
| max | 1.000000 | |

```
unique_values = hiv_df['HIV_active'].unique()
print(f"Unique values in 'HIV_active': {unique_values}")
```

Unique values in 'HIV_active': [0 1]



```
unique_values = hiv_df['smiles'].unique()
print(f"Unique values in 'smiles': {unique_values}")
print(f"length of unique value: {len(unique_values)}")
```

```
Unique values in 'smiles': ['CC1=[O+] [Cu-3]2([O+]=C(CC)C1)[O+]=C(CC)CC(CC)=[O+]2'
'C(=Cc1cccc1)C1=[O+] [Cu-3]2([O+]=C(C=Cc3cccc3)CC(c3cccc3)=[O+]2)[O+]=C(c2cccc2)C1'
'CC(=O)N1c2cccc2Sc2c1ccc1cccc21' ...
'Cc1ccc(N2C(=O)C3c4[nH]c5cccc5c4C4CCC(C(C)(C)C)CC4C3C2=O)cc1'
'Cc1cccc(N2C(=O)C3c4[nH]c5cccc5c4C4CCC(C(C)(C)C)CC4C3C2=O)c1'
'CCCCC=C(c1cc(Cl)c(OC)c(-c2nc(C)no2)c1)c1cc(Cl)c(OC)c(-c2nc(C)no2)c1']
length of unique value: 41127
```

```
# Filter rows where 'your_column' is not equal to 1 or 0
filtered_df = hiv_df[(hiv_df['HIV_active'] != 1) & (hiv_df['HIV_active'] != 0)]
filtered_df
```

smiles activity HIV_active 

```
# Filter rows where 'target_column' is equal to 1h
hiv_df_filtered_active = hiv_df[hiv_df['HIV_active'] == 1]
hiv_df_filtered_active
```

| | smiles | activity | HIV_active |  |
|-------|--|----------|------------|---|
| 11 | O=C(O)Cc1ccc(SSc2ccc(CC(=O)O)cc2)cc1 | CM | 1 |  |
| 16 | NNP(=S)(NN)c1cccc1 | CM | 1 | |
| 80 | O=Nc1ccc(O)c(N=O)c1O | CM | 1 | |
| 203 | Oc1ccc(Cl)cc1C(c1cc(Cl)ccc1O)C(Cl)(Cl)Cl | CM | 1 | |
| 234 | NNC(=O)c1cccc1SSc1cccc1C(=O)NN | CM | 1 | |
| ... | ... | ... | ... | |
| 41090 | Cc1cn(COCCCOCC(=O)c2cccc2)c(=O)[nH]c1=O | CM | 1 | |
| 41092 | Cc1cn(C2CC3C(COC(CCC[Se]c4cccc4)N3O)O2)c(=O)[... | CM | 1 | |
| 41093 | Cc1cn(C2CC3C(COC(CCCC[Se]c4cccc4)N3O)O2)c(=O)... | CM | 1 | |
| 41098 | Cc1cn(C2CC3C(COC(CC[Se]C#N)N3O)O2)c(=O)[nH]c1=O | CM | 1 | |
| 41099 | C[Se]CCC1OCC2OC(n3cc(C)c(=O)[nH]c3=O)CC2N1O | CA | 1 | |

1443 rows x 3 columns

Next steps:  [View recommended plots](#)

```
# Filter rows where 'target_column' is equal to 1h
hiv_df_filtered_inactive = hiv_df[hiv_df['HIV_active'] == 0]
hiv_df_filtered_inactive = hiv_df_filtered_inactive.sample(n=1500, axis=0, random_state=42)
hiv_df_filtered_inactive
```

| | smiles | activity | HIV_active |  |
|-------|---|----------|------------|---|
| 2428 | O=C1c2cccc2-c2nc3cccc3nc21 | CI | 0 |  |
| 6197 | O=C(CSc1cc(-c2ccc(Cl)cc2)s[s+])1c1cccc1 | CI | 0 | |
| 17138 | O=C(C=Nc1cccc1C(=O)O)c1ccco1 | CI | 0 | |
| 12261 | CCCCCCCCCCCCCCCC[N+](C)(C)Cc1ccc(C[N+](C)(C)... | CI | 0 | |
| 3588 | N#CSC1CCCCCCC1SC#N | CI | 0 | |
| ... | ... | ... | ... | |
| 18477 | CC(=O)OC1(C#N)CC2OC1C1C2N1C(=O)OC(C)(C)C | CI | 0 | |
| 1189 | CCOC(=O)C1Cc2cc(C)c(C)cc2N(C)C1=O | CI | 0 | |
| 36657 | CCOC(=O)N1CCN(c2ccc3c(C)cc(C)nc3n2)CC1 | CI | 0 | |
| 27919 | CN(C)C=Nc1ccc2c3c(cccc13)-c1cccc1-2 | CI | 0 | |
| 13479 | CCC1CC2CC3c4[nH]c5ccc(OC)cc5c4CCN(C2)C13.Cl | CI | 0 | |

1500 rows x 3 columns

Next steps:  [View recommended plots](#)

```
hiv_df_sampled = pd.concat([hiv_df_filtered_active, hiv_df_filtered_inactive], axis=0, ignore_index=True)
hiv_df_sampled
```

| | smiles | activity | HIV_active | |
|------|--|----------|------------|--|
| 0 | <chem>O=C(O)Cc1ccc(SSc2ccc(CC(=O)O)cc2)cc1</chem> | CM | 1 | |
| 1 | <chem>NNP(=S)(NN)c1cccc1</chem> | CM | 1 | |
| 2 | <chem>O=Nc1ccc(O)c(N=O)c1O</chem> | CM | 1 | |
| 3 | <chem>Oc1ccc(Cl)cc1C(c1cc(Cl)ccc1O)C(Cl)(Cl)Cl</chem> | CM | 1 | |
| 4 | <chem>NNC(=O)c1cccc1SSc1cccc1C(=O)NN</chem> | CM | 1 | |
| ... | ... | ... | ... | |
| 2938 | <chem>CC(=O)OC1(C#N)CC2OC1C1C2N1C(=O)OC(C)(C)C</chem> | CI | 0 | |
| 2939 | <chem>CCOC(=O)C1Cc2cc(C)c(C)cc2N(C)C1=O</chem> | CI | 0 | |
| 2940 | <chem>CCOC(=O)N1CCN(c2ccc3c(C)cc(C)nc3n2)CC1</chem> | CI | 0 | |
| 2941 | <chem>CN(C)C=Nc1ccc2c3c(cccc13)-c1cccc1-2</chem> | CI | 0 | |
| 2942 | <chem>CCC1CC2CC3c4[nH]c5ccc(OC)cc5c4CCN(C2)C13.Cl</chem> | CI | 0 | |

2943 rows x 3 columns

Next steps: [View recommended plots](#)

```
# Randomly shuffle rows
hiv_df_sampled = hiv_df_sampled.sample(frac=1, random_state=42)
```

```
hiv_df_sampled.head()
```

| | smiles | activity | HIV_active | |
|------|---|----------|------------|--|
| 240 | <chem>Cc1cc2c(c(=O)o1)C1=S(SC(c3cccc3)=C1)S2</chem> | CM | 1 | |
| 2325 | <chem>N#CN1CCC=C(c2cc3cccc3[nH]2)C1</chem> | CI | 0 | |
| 1676 | <chem>CCC1SC(C)C(=O)NC1=O</chem> | CI | 0 | |
| 1952 | <chem>O=C1CC2(CCN(Cc3cccc3)CC2)CC(=O)N1</chem> | CI | 0 | |
| 677 | <chem>CC(=O)OC1SC(c2c(F)cccc2F)n2c1nc1cccc12</chem> | CM | 1 | |

Next steps: [View recommended plots](#)

```
hiv_df_sampled.to_csv('HIV_2.csv', index=False)
# .drop(['activity'], axis=1).
hiv_df_sampled_2 = pd.read_csv("HIV_2.csv")
hiv_df_sampled_2.head()
hiv_df_sampled_2.tail()
```

| | smiles | activity | HIV_active | |
|---|---|----------|------------|--|
| 0 | <chem>Cc1cc2c(c(=O)o1)C1=S(SC(c3cccc3)=C1)S2</chem> | CM | 1 | |
| 1 | <chem>N#CN1CCC=C(c2cc3cccc3[nH]2)C1</chem> | CI | 0 | |
| 2 | <chem>CCC1SC(C)C(=O)NC1=O</chem> | CI | 0 | |
| 3 | <chem>O=C1CC2(CCN(Cc3cccc3)CC2)CC(=O)N1</chem> | CI | 0 | |
| 4 | <chem>CC(=O)OC1SC(c2c(F)cccc2F)n2c1nc1cccc12</chem> | CM | 1 | |

| | smiles | activity | HIV_active | |
|------|--|----------|------------|--|
| 2938 | <chem>O=C(CS)Nc1cccc(O)c1</chem> | CI | 0 | |
| 2939 | <chem>O=C(Nc1ccc(N=Nc2ccc(S(=O)(=O)O)cc2)cc1)c1ccc(N...</chem> | CM | 1 | |
| 2940 | <chem>NC(=O)CCN(CCC(N)=O)CCC(N)=O</chem> | CI | 0 | |
| 2941 | <chem>Cn1nc2cccc2cc1=O</chem> | CI | 0 | |
| 2942 | <chem>C[n+]1c(C=NNC(=O)c2ccc(C(=O)NN=Cc3cn4cccc4[n+...</chem> | CI | 0 | |

```
arguments = [  
    '--data_path', 'HIV_2.csv',  
    '--dataset_type', 'classification',  
    '--save_dir', 'test_checkpoints_multimolecule',  
    '--epochs', '30',  
    '--save_smiles_splits',  
    '--quiet',  
    '--batch_size', '64',  
    '--ignore_columns', 'activity',  
    '--depth', '5',  
    '--hidden_size', '300'  
]  
  
args = chemprop.args.TrainArgs().parse_args(arguments)
```

```
mean_score, std_score = chemprop.train.cross_validate(args=args, train_func=chemprop.train.run_training)
```

```
78% ██████████ | 29/37 [00:02<00:00, 10.27it/s]  
84% ██████████ | 31/37 [00:03<00:00, 10.35it/s]  
89% ██████████ | 33/37 [00:03<00:00, 10.40it/s]  
95% ██████████ | 35/37 [00:03<00:00, 10.11it/s]  
100% ██████████ | 37/37 [00:03<00:00, 9.64it/s]
```

```
0% | | 0/5 [00:00<?, ?it/s]  
20% █ | 1/5 [00:00<00:00, 5.84it/s]  
40% ███ | 2/5 [00:00<00:00, 5.72it/s]  
60% █████ | 3/5 [00:00<00:00, 5.86it/s]  
80% ███████ | 4/5 [00:00<00:00, 6.10it/s]  
100% ██████████ | 5/5 [00:00<00:00, 6.85it/s]  
67% ██████████ | 20/30 [01:39<00:45, 4.60s/it]  
0% | | 0/37 [00:00<?, ?it/s]  
3% █ | 1/37 [00:00<00:06, 5.50it/s]  
5% █ | 2/37 [00:00<00:06, 5.45it/s]  
8% █ | 3/37 [00:00<00:06, 5.31it/s]  
11% █ | 4/37 [00:00<00:06, 5.11it/s]  
14% █ | 5/37 [00:00<00:06, 5.18it/s]  
16% █ | 6/37 [00:01<00:05, 5.37it/s]  
19% █ | 7/37 [00:01<00:05, 5.43it/s]  
22% █ | 8/37 [00:01<00:04, 6.31it/s]  
24% █ | 9/37 [00:01<00:03, 7.06it/s]  
27% █ | 10/37 [00:01<00:03, 7.65it/s]  
32% ███ | 12/37 [00:01<00:02, 8.84it/s]  
35% ███ | 13/37 [00:01<00:02, 9.05it/s]  
38% ███ | 14/37 [00:02<00:02, 8.79it/s]  
41% ███ | 15/37 [00:02<00:02, 8.78it/s]  
43% ███ | 16/37 [00:02<00:02, 8.61it/s]  
46% ███ | 17/37 [00:02<00:02, 8.91it/s]  
51% ████ | 19/37 [00:02<00:01, 9.61it/s]  
54% ████ | 20/37 [00:02<00:01, 9.58it/s]  
59% ████ | 22/37 [00:02<00:01, 10.09it/s]  
62% ████ | 23/37 [00:02<00:01, 9.95it/s]  
68% ████ | 25/37 [00:03<00:01, 10.05it/s]  
70% ████ | 26/37 [00:03<00:01, 9.66it/s]  
73% ████ | 27/37 [00:03<00:01, 9.39it/s]  
76% ████ | 28/37 [00:03<00:00, 9.44it/s]  
78% ████ | 29/37 [00:03<00:00, 9.47it/s]  
84% █████ | 31/37 [00:04<00:01, 5.16it/s]  
86% █████ | 32/37 [00:04<00:00, 5.70it/s]  
92% █████ | 34/37 [00:04<00:00, 6.98it/s]  
95% █████ | 35/37 [00:04<00:00, 7.32it/s]  
97% █████ | 36/37 [00:04<00:00, 7.54it/s]
```

```
0% | | 0/5 [00:00<?, ?it/s]  
20% █ | 1/5 [00:00<00:00, 9.60it/s]  
60% █████ | 3/5 [00:00<00:00, 10.74it/s]  
100% ██████████ | 5/5 [00:00<00:00, 11.80it/s]  
70% ██████████ | 21/30 [01:45<00:43, 4.81s/it]  
0% | | 0/37 [00:00<?, ?it/s]  
3% █ | 1/37 [00:00<00:03, 9.65it/s]  
5% █ | 2/37 [00:00<00:03, 9.30it/s]  
8% █ | 3/37 [00:00<00:03, 9.40it/s]  
11% █ | 4/37 [00:00<00:03, 9.40it/s]  
14% █ | 5/37 [00:00<00:03, 9.30it/s]  
19% █ | 7/37 [00:00<00:03, 9.91it/s]  
22% █ | 8/37 [00:00<00:02, 9.79it/s]
```

```
mean_score, std_score  
(0.8328736900165471, 0.0)
```

```
bp_df = pd.read_csv("BBBP.csv")
bp_df.head()
```

| | num | name | p_np | smiles |
|---|-----|----------------------|------|---|
| 0 | 1 | Propanolol | 1 | [Cl].CC(C)NCC(O)COc1cccc2ccccc12 |
| 1 | 2 | Terbutylchlorambucil | 1 | C(=O)(OC(C)(C)C)CCCc1ccc(cc1)N(CCGI)CCCI |
| 2 | 3 | 40730 | 1 | c12c3c(N4CCN(C)CC4)c(F)cc1c(c(C(O)=O)cn2C(C)CO... |
| 3 | 4 | 24 | 1 | C1CCN(CC1)Cc1cccc(c1)OCCCN(C(=O)C |
| 4 | 5 | cloxacillin | 1 | Cc1onc(c2ccccc2Cl)c1C(=O)N[C@H]3[C@H]4SC(C)(C)... |

Next steps: [View recommended plots](#)

```
bp_df.tail()
```

| | num | name | p_np | smiles |
|------|------|----------------------------------|------|---|
| 2045 | 2049 | licostinel | 1 | C1=C(Cl)C(=C(C2=C1NC(=O)C(N2)=O)[N+](=O)[O-])Cl |
| 2046 | 2050 | ademetonine(adenosyl-methionine) | 1 | [C@H]3([N]2C1=C(C(=NC=N1)N)N=C2)[C@@H]([C@@H](... |
| 2047 | 2051 | mesocarb | 1 | [O+]1=N[N](C=C1[N-]C(NC2=CC=CC=C2)=O)C(CC3=CC=... |
| 2048 | 2052 | ... | ... | C1=C(OC)C(=CC2=C1C(=N+)(C(=C2CC)C) |

```
bp_df.drop(['num', 'name', 'p_np'], axis=1).to_csv('BBBP_2.csv', index=False)
```

```
bp_df_2 = pd.read_csv("BBBP_2.csv")
bp_df_2.head()
bp_df_2.tail()
```

| | smiles |
|------|---|
| 0 | [Cl].CC(C)NCC(O)COc1cccc2ccccc12 |
| 1 | C(=O)(OC(C)(C)C)CCCc1ccc(cc1)N(CCGI)CCCI |
| 2 | c12c3c(N4CCN(C)CC4)c(F)cc1c(c(C(O)=O)cn2C(C)CO... |
| 3 | C1CCN(CC1)Cc1cccc(c1)OCCCN(C(=O)C |
| 4 | Cc1onc(c2ccccc2Cl)c1C(=O)N[C@H]3[C@H]4SC(C)(C)... |
| 2045 | C1=C(Cl)C(=C(C2=C1NC(=O)C(N2)=O)[N+](=O)[O-])Cl |
| 2046 | [C@H]3([N]2C1=C(C(=NC=N1)N)N=C2)[C@@H]([C@@H](... |
| 2047 | [O+]1=N[N](C=C1[N-]C(NC2=CC=CC=C2)=O)C(CC3=CC=... |
| 2048 | C1=C(OC)C(=CC2=C1C(=N+)(C(=C2CC)C)[NH-])C3=CC... |
| 2049 | [N+](=NCC(=O)N[C@@H]([C@H](O)C1=CC=C([N+])([O-])... |

```
arguments = [
    '--test_path', 'BBBP_2.csv',
    '--preds_path', 'BBBP_preds.csv',
    '--checkpoint_dir', 'test_checkpoints_multimolecule'
]
```

```
args = chemprop.args.PredictArgs().parse_args(arguments)
preds = chemprop.train.make_predictions(args=args)
```

```
[19:56:48] Explicit valence for atom # 1 N, 4, is greater than permitted
[19:56:48] WARNING: not removing hydrogen atom without neighbors
[19:56:48] Explicit valence for atom # 6 N, 4, is greater than permitted
[19:56:48] WARNING: not removing hydrogen atom without neighbors
[19:56:48] WARNING: not removing hydrogen atom without neighbors
```

```
bp_preds_df = pd.read_csv("BBBP_preds.csv")
bp_preds_df.head()
```



| | smiles | HIV_active |
|---|--|----------------------|
| 0 | <chem>[Cl].CC(C)NCC(O)COc1cccc2cccc12</chem> | 0.07791923731565475 |
| 1 | <chem>C(=O)(OC(C)(C)C)CCCc1ccc(cc1)N(CCCl)CCCl</chem> | 0.052813779562711716 |
| 2 | <chem>c12c3c(N4CCN(C)CC4)c(F)cc1c(c(C(O)=O)cn2C(C)CO...</chem> | 0.5678612589836121 |
| 3 | <chem>C1CCN(CC1)Cc1cccc(c1)OCCNC(=O)C</chem> | 0.0569111704826355 |
| 4 | <chem>Cc1onc(2cccc2Cl)c1C(=O)N[C@H]3[C@H]4SC(C)(C)...</chem> | 0.4187469184398651 |

Next steps: ☒ View recommended plots



```
bp_preds_df.tail()
```

| | smiles | HIV_active |
|------|--|---------------------|
| 2045 | <chem>C1=C(Cl)C(=C(C2=C1NC(=O)C(N2)=O)[N+](=O)[O-])Cl</chem> | 0.1297084391117096 |
| 2046 | <chem>[C@H]3([N]2C1=C(C(=NC=N1)N)N=C2)[C@H]([C@H](...</chem> | 0.23322035372257233 |
| 2047 | <chem>[O+]1=N[N](C=C1[N-]C(NC2=CC=CC=C2)=O)C(CC3=CC=...</chem> | 0.3322729766368866 |
| 2048 | <chem>C1=C(OC)C(=CC2=C1C(=[N+](C(=C2CC)C)[NH-])C3=CC...</chem> | 0.3129490911960602 |
| 2049 | <chem>[N+](=NCC(=O)N[C@H]([C@H](O)C1=CC=C([N+](O)...</chem> | 0.2837357521057129 |

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| | smiles | HIV_active |  |
|--------|---|------------|--|
| count | 2050 | 2050 |  |
| unique | 2050 | 2004 | |
| top | [N+](=NCC(=O)N[C@@H]([C@H](O)C1=CC=C([N+](O-)... Invalid SMILES | | |
| freq | 1 | 11 | |

```
bp_preds_df = bp_preds_df[bp_preds_df['HIV_active'] != "Invalid SMILES"]
bp_preds_df.describe()
```

| | smiles | HIV_active |  |
|--------|---|------------|---|
| count | 2039 | 2039 |  |
| unique | 2039 | 2003 | |
| top | [N+](=NCC(=O)N[C@@H]([C@H](O)C1=CC=C([N+](O-)... 0.0426582507789135 | | |
| freq | 1 | 3 | |



```
bp_preds_df['HIV_active'] = bp_preds_df['HIV_active'].astype(float)
```

```
bp_preds_df['HIV_active_2'] = bp_preds_df['HIV_active'].apply(lambda x: 1 if x > 0.8 else 0)
bp_preds_df.head()
```

| | smiles | HIV_active | HIV_active_2 |  |
|---|---|------------|--------------|---|
| 0 | [Cl].CC(C)NCC(O)COc1cccc2ccccc12 | 0.077919 | 0 |  |
| 1 | C(=O)(OC(C)(C)C)CCc1ccc(cc1)N(CCC)CCCl | 0.052814 | 0 | |
| 2 | c12c3c(N4CCN(C)CC4)c(F)cc1c(c(C(O)=O)cn2C(C)CO... | 0.567861 | 0 | |
| 3 | C1CCN(CC1)Cc1cccc(c1)OCCCN(C=O)C | 0.056911 | 0 | |
| 4 | Cc1onc(c2ccccc2Cl)c1C(=O)N[C@H]3[C@H]4SC(C)(C)... | 0.418747 | 0 | |

Next steps:  [View recommended plots](#)

```
bp_preds_df.describe()
```

| | HIV_active | HIV_active_2 |  |
|-------|-------------|--------------|---|
| count | 2039.000000 | 2039.000000 |  |
| mean | 0.263883 | 0.009809 | |
| std | 0.201033 | 0.098576 | |
| min | 0.000762 | 0.000000 | |
| 25% | 0.097246 | 0.000000 | |
| 50% | 0.208566 | 0.000000 | |
| 75% | 0.408756 | 0.000000 | |
| max | 0.988302 | 1.000000 | |

```
# Filter rows where 'target_column' is equal to 1
bp_preds_df_filtered = bp_preds_df[bp_preds_df['HIV_active_2'] == 1]
bp_preds_df_filtered
```


1 to 20 of 20 entries

Filter

?

| index | smiles | HIV_active | H |
|-------|--|--------------------|---|
| 11 | CC1=CN([C@H]2C[C@H](F)[C@@H](CO)O2)C(=O)NC1=O | 0.9441607594490051 | |
| 235 | OC[C@H]1O[C@H](C[C@@H]1O)N2C=C(F)C(=O)NC2=O | 0.8643155097961426 | |
| 270 | Cc1nnc2CN=C(c3ccccc3)c4cc(Cl)ccc4n12 | 0.8021654486656189 | |
| 279 | OC[C@H]1O[C@@H](OC2=C(Oc3cc(O)cc(O)c3C2=O)c4ccc(O)c(O)c4)[C@H](O)[C@@H](O)[C@@H]1O | 0.8800746202468872 | |
| 289 | OC[C@@H]1CC[C@H](O1)n2cnc3C(=O)N=CNc23 | 0.8122326731681824 | |
| 319 | CC1=CN([C@@H]2O[C@H](CO)C=C2)C(=O)NC1=O | 0.9302963614463806 | |
| 445 | CC1=CN([C@H]2C[C@H](N=[N+]=[N-])[C@@H](CO)O2)C(=O)NC1=O | 0.9854286313056946 | |
| 540 | Cc1nnc2CN=C(c3ccccc3Cl)c4cc(Cl)ccc4n12 | 0.8569784164428711 | |
| 581 | CN1CCC23C4CCC(O)C2Oc5c(O)ccc(CC14)c35 | 0.816520631313324 | |
| 711 | Cc1ncc2CN=C(c3ccccc3F)c4cc(Cl)ccc4n12.Oc(=O)\C=C/C(O)=O | 0.8424991369247437 | |
| 771 | Cn1nnnc1SCC2=C(N3[C@H](SC2)[C@H](NC(=O)CS\C=C/C#N)C3=O)C(O)=O | 0.808535635471344 | |
| 785 | CC1=C(N2[C@H](SC1)[C@H](NC(=O)Cc3ccc(cc3)C4=NCCCN4)C2=O)C(O)=O | 0.8538467288017273 | |
| 912 | CO[C@H]1/C=C/O[C@@]2(C)Oc3c(C)c(O)c4c(O)c(NC(=O)C(=C\C=C\C[C@H](C)[C@H](O)[C@@H](C)[C@@H](O)[C@@H](C)[C@H](OC(C)=O)[C@@H]1C)/C)cc(O)c4c3C2=O | 0.8292530179023743 | |
| 945 | Oc1ccc(cc1)/C=C([N+]#[C-])/C(=C/c2ccc(O)cc2)[N+]#[C-] | 0.8675630688667297 | |
| 973 | COC1(C(F)(F)C1(F)Cl)F | 0.9545143246650696 | |
| 1052 | C1=C(Br)SC2=C1C(=NCC3=NN=C([N]23)C)C4=C(C=CC=C4)Cl | 0.8791234493255615 | |
| 1086 | C1=C(SC2=C1C(=NCC3=NN=C([N]23)C4CCCC4)C5=CC=CC=C5Cl)Br | 0.8506717085838318 | |
| 1365 | [C@@H]25CC1=CC=C(C4=C1[C@@]3([C@H]2CC[C@@H]([C@@H]3O4)OC)CCN5C)O | 0.8105615973472595 | |
| 1822 | C1=C(SC2=C1C(=NCC3=NN=C([N]23)C)C4=CC=CC=C4Cl)CC | 0.8752243518829346 | |
| 2044 | [N+](=[N-])=O | 0.9883024096488953 | |

Next steps:

View recommended plots

```
smiles_to_check = bp_preds_df_filtered['smiles'].to_list()

hiv_df_sampled_2[hiv_df_sampled_2['smiles'].isin(smiles_to_check)]

smiles activity HIV_active
hiv_df[hiv_df['smiles'].isin(smiles_to_check)]

smiles activity HIV_active
bp_df[bp_df['smiles'].isin(smiles_to_check)]
```

| num | | name | p_np | |
|------|------|-----------------------|------|--|
| 11 | 12 | alovudine | 1 | <chem>CC1=CN([C@H]2C[C@H](F)[C@@H](CO)O2)C(=O)N</chem> |
| 235 | 237 | floxuridine | 0 | <chem>OC[C@H]1O[C@H](C[C@@H]1O)N2C=C(F)C(=O)N</chem> |
| 270 | 272 | Alprazolam | 1 | <chem>Cc1nnc2CN=C(c3ccccc3)c4cc(Cl)cc12</chem> |
| 279 | 281 | Isoquercitrin | 0 | <chem>OC[C@H]1O[C@@H](OC2=C(Oc3cc(O)cc(O)c3C2=O)C[C@H]1O</chem> |
| 289 | 291 | Didanosine | 0 | <chem>OC[C@@H]1CC[C@@H](O1)n2cnc3C(=O)N=CN=C3</chem> |
| 319 | 321 | Stavudine | 1 | <chem>CC1=CN([C@@H]2O[C@H](CO)C=C2)C(=O)N</chem> |
| 445 | 447 | zidovudine | 0 | <chem>CC1=CN([C@H]2C[C@H](N=[N+]=[N-])[C@@H](CO)N2)C(=O)N</chem> |
| 540 | 542 | Triazolam | 1 | <chem>Cc1nnc2CN=C(c3ccccc3Cl)c4cc(Cl)cc12</chem> |
| 581 | 583 | dihydromorphine | 1 | <chem>CN1CCC23C4CCC(O)C2Oc5c(O)ccc(CC4)C13</chem> |
| 711 | 713 | midazolam maleate | 1 | <chem>Cc1ncc2CN=C(c3ccccc3F)c4cc(Cl)ccc4n12.O=C(=O)OC</chem> |
| 771 | 773 | cefivitril | 0 | <chem>Cn1nnnc1SCC2=C(N3[C@H](SC2)[C@H](NC(=O)CS)C3)C(=O)N</chem> |
| 785 | 787 | cefrotil | 0 | <chem>CC1=C(N2[C@H](SC1)[C@H](NC(=O)Cc3ccc(cc3)C(=O)N)C2)C(=O)N</chem> |
| 912 | 914 | rifamycin | 0 | <chem>CO[C@H]1/C=C/O[C@@]2(C)Oc3c(C)c(O)c4c(O)c(CO)cc321</chem> |
| 945 | 947 | xantocillin | 0 | <chem>Oc1ccc(cc1)/C=C([N+]#[C-])/C(=C/c2ccc(O)cc2)C</chem> |
| 973 | 975 | aliflurane | 1 | <chem>COC1(C(F)(F)F)C(F)(F)F1</chem> |
| 1052 | 1054 | brotizolam | 1 | <chem>C1=C(Br)SC2=C1C(=NCC3=NN=C([N]23)C)C4=C(C=CC4)C</chem> |
| 1086 | 1089 | ciclotizolam | 1 | <chem>C1=C(SC2=C1C(=NCC3=NN=C([N]23)C4CCCCC4)C5=C(C(=CC5)C)C2)C</chem> |
| 1365 | 1369 | methyldihydromorphine | 1 | <chem>[C@@H]25CC1=CC=C(C4=C1[C@@]3([C@H]2CC[C@@H]3C)C=C4)C1</chem> |
| 1822 | 1826 | etizolam | 1 | <chem>C1=C(SC2=C1C(=NCC3=NN=C([N]23)C)C4=CC=CC=C4C2)C</chem> |
| 2044 | 2048 | nitrous-oxide | 1 | <chem>[N+]</chem> |



```
bp_df_final = pd.merge(bp_df[bp_df['smiles'].isin(smiles_to_check)], bp_preds_df_filtered, on='smiles')
bp_df_final
```

| num | | name | p_np | s |
|-----|------|-----------------------|------|--|
| 0 | 12 | alovudine | 1 | <chem>CC1=CN([C@H]2C[C@H](F)[C@@H](CO)O2)C(=O)N</chem> |
| 1 | 237 | floxuridine | 0 | <chem>OC[C@H]1O[C@H](C[C@@H]1O)N2C=C(F)C(=O)N</chem> |
| 2 | 272 | Alprazolam | 1 | <chem>Cc1nnc2CN=C(c3ccccc3)c4cc(Cl)cc12</chem> |
| 3 | 281 | Isoquercitrin | 0 | <chem>OC[C@H]1O[C@@H](OC2=C(Oc3cc(O)cc(O)c3C2=O)C[C@H]1O</chem> |
| 4 | 291 | Didanosine | 0 | <chem>OC[C@@H]1CC[C@@H](O1)n2cnc3C(=O)N=CN=C3</chem> |
| 5 | 321 | Stavudine | 1 | <chem>CC1=CN([C@@H]2O[C@H](CO)C=C2)C(=O)N</chem> |
| 6 | 447 | zidovudine | 0 | <chem>CC1=CN([C@H]2C[C@H](N=[N+]=[N-])[C@@H](CO)N2)C(=O)N</chem> |
| 7 | 542 | Triazolam | 1 | <chem>Cc1nnc2CN=C(c3ccccc3Cl)c4cc(Cl)cc12</chem> |
| 8 | 583 | dihydromorphine | 1 | <chem>CN1CCC23C4CCC(O)C2Oc5c(O)ccc(CC4)C13</chem> |
| 9 | 713 | midazolam maleate | 1 | <chem>Cc1ncc2CN=C(c3ccccc3F)c4cc(Cl)ccc4n12.O=C(=O)OC</chem> |
| 10 | 773 | cefivitril | 0 | <chem>Cn1nnnc1SCC2=C(N3[C@H](SC2)[C@H](NC(=O)CS)C3)C(=O)N</chem> |
| 11 | 787 | cefrotil | 0 | <chem>CC1=C(N2[C@H](SC1)[C@H](NC(=O)Cc3ccc(cc3)C(=O)N)C2)C(=O)N</chem> |
| 12 | 914 | rifamycin | 0 | <chem>CO[C@H]1/C=C/O[C@@]2(C)Oc3c(C)c(O)c4c(O)c(CO)cc321</chem> |
| 13 | 947 | xantocillin | 0 | <chem>Oc1ccc(cc1)/C=C([N+]#[C-])/C(=C/c2ccc(O)cc2)C</chem> |
| 14 | 975 | aliflurane | 1 | <chem>COC1(C(F)(F)F)C(F)(F)F1</chem> |
| 15 | 1054 | brotizolam | 1 | <chem>C1=C(Br)SC2=C1C(=NCC3=NN=C([N]23)C)C4=C(C=CC4)C</chem> |
| 16 | 1089 | ciclotizolam | 1 | <chem>C1=C(SC2=C1C(=NCC3=NN=C([N]23)C4CCCCC4)C5=C(C(=CC5)C)C2)C</chem> |
| 17 | 1369 | methyldihydromorphine | 1 | <chem>[C@@H]25CC1=CC=C(C4=C1[C@@]3([C@H]2CC[C@@H]3C)C=C4)C1</chem> |
| 18 | 1826 | etizolam | 1 | <chem>C1=C(SC2=C1C(=NCC3=NN=C([N]23)C)C4=CC=CC=C4C2)C</chem> |
| 19 | 2048 | nitrous-oxide | 1 | <chem>[N+](=O)</chem> |

Next steps: [View recommended plots](#)



```
bp_df_final.to_csv('HIV_result.csv', index=False)
```

```
sub_df = pd.read_csv("substances.csv")
sub_df.head()
```

| | zinc_id | smiles |  |
|---|-------------------|---|---|
| 0 | ZINC0000000000027 | <chem>N[C@@H](CCc1ccc(N(CCCl)CCCl)cc1)C(=O)O</chem> |  |
| 1 | ZINC000016090786 | <chem>N[C@H](CCc1ccc(N(CCCl)CCCl)cc1)C(=O)O</chem> | |
| 2 | ZINC000001763088 | <chem>N[C@H](CCc1ccc(N(CCCl)CCCl)cc1)C(=O)O</chem> | |
| 3 | ZINC000002033385 | <chem>N[C@@H](CCc1ccc(N(CCCl)CCCl)cc1)C(=O)O</chem> | |
| 4 | ZINC000000001673 | <chem>N[C@@H](Cc1ccc(N(CCCl)CCCl)cc1)C(=O)O</chem> | |

Next steps:  [View recommended plots](#)

```
sub_df.tail()
```

| | zinc_id | smiles |  |
|----|------------------|---|---|
| 46 | ZINC000196349655 | <chem>O=C(O)CCSc1ccc(N(CCCl)CCCl)cc1</chem> |  |
| 47 | ZINC000064454242 | <chem>N=NCCCc1ccc(N(CCCl)CCCl)cc1</chem> | |
| 48 | ZINC000005161807 | <chem>O=C(O)C/C=C/c1ccc(N(CCCl)CCCl)cc1</chem> | |
| 49 | ZINC000001682294 | <chem>O=C(O)CCOc1ccc(N(CCCl)CCCl)cc1</chem> | |
| 50 | ZINC000079564304 | <chem>O=C(O)CNC(=O)c1ccc(N(CCCl)CCCl)cc1</chem> | |

```
sub_df.info()
```

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 51 entries, 0 to 50
Data columns (total 2 columns):
#   Column  Non-Null Count  Dtype
---  ---
0    zinc_id  51 non-null         object
1    smiles   51 non-null         object
dtypes: object(2)
memory usage: 944.0+ bytes
```

```
arguments = [
    '--test_path', 'substances.csv',
    '--preds_path', 'substances_preds.csv',
    '--checkpoint_dir', 'test_checkpoints_multimolecule',
    '--smiles_columns', 'smiles'
]
```

```
args = chemprop.args.PredictArgs().parse_args(arguments)
preds = chemprop.train.make_predictions(args=args)
```

```
Loading training args
Setting molecule featurization parameters to default.
Loading data
51it [00:00, 60426.41it/s]
100%|██████████| 51/51 [00:00<00:00, 76780.15it/s]
/usr/local/lib/python3.10/dist-packages/torch/utils/data/dataloader.py:557: UserWarning: This DataLoader will create
  warnings.warn(_create_warning_msg(
Validating SMILES
Test size = 51
0%|          | 0/1 [00:00<?, ?it/s]Loading pretrained parameter "encoder.encoder.0.cached_zero_vector".
Loading pretrained parameter "encoder.encoder.0.W_i.weight".
Loading pretrained parameter "encoder.encoder.0.W_h.weight".
Loading pretrained parameter "encoder.encoder.0.W_o.weight".
Loading pretrained parameter "encoder.encoder.0.W_o.bias".
Loading pretrained parameter "readout.1.weight".
Loading pretrained parameter "readout.1.bias".
Loading pretrained parameter "readout.4.weight".
Loading pretrained parameter "readout.4.bias".
Moving model to cuda

0%|          | 0/2 [00:00<?, ?it/s]
50%|██████    | 1/2 [00:00<00:00, 2.54it/s]
100%|██████████| 1/1 [00:01<00:00, 1.20s/it]Saving predictions to substances_preds.csv
```

Elapsed time = 0:00:01

```
fda_df = pd.read_csv("fda_approved.csv")
fda_df.head()
```

| | zinc_id | smiles | |
|---|------------------|--|--|
| 0 | ZINC000001530427 | <chem>C[C@@H]1O[C@@H]1P(=O)(O)O</chem> | |
| 1 | ZINC000003807804 | <chem>Clc1cccc1C(c1cccc1)(c1cccc1)n1ccnc1</chem> | |
| 2 | ZINC000000120286 | <chem>Nc1nc(N)c2nc(-c3ccccc3)c(N)nc2n1</chem> | |
| 3 | ZINC000242548690 | <chem>C[C@H]1O[C@@H](O[C@H]2[C@@H](O)C[C@H](O[C@H]3[...])</chem> | |
| 4 | ZINC000000008492 | <chem>Oc1cccc2cccn12</chem> | |

Next steps: [View recommended plots](#)

```
arguments = [
    '--test_path', 'fda_approved.csv',
    '--preds_path', 'fda_approved_preds.csv',
    '--checkpoint_dir', 'test_checkpoints_multimolecule',
    '--smiles_columns', 'smiles'
]
```

```
args = chemprop.args.PredictArgs().parse_args(arguments)
preds = chemprop.train.make_predictions(args=args)
```



```
Loading training args
Setting molecule featurization parameters to default.
Loading data
892it [00:00, 90156.61it/s]
100%|██████████| 892/892 [00:00<00:00, 51982.98it/s]
/usr/local/lib/python3.10/dist-packages/torch/utils/data/dataloader.py:557: UserWarning: This DataLoader will create
warnings.warn(_create_warning_msg)
Validating SMILES
Test size = 892
0%|          | 0/1 [00:00<?, ?it/s]Loading pretrained parameter "encoder.encoder.0.cached_zero_vector".
Loading pretrained parameter "encoder.encoder.0.W_i.weight".
Loading pretrained parameter "encoder.encoder.0.W_h.weight".
Loading pretrained parameter "encoder.encoder.0.W_o.weight".
Loading pretrained parameter "encoder.encoder.0.W_o.bias".
Loading pretrained parameter "readout.1.weight".
Loading pretrained parameter "readout.1.bias".
Loading pretrained parameter "readout.4.weight".
Loading pretrained parameter "readout.4.bias".
Moving model to cuda

0%|          | 0/18 [00:00<?, ?it/s]
6%|          | 1/18 [00:02<00:48, 2.86s/it]
11%|         | 2/18 [00:03<00:20, 1.28s/it]
50%|██████    | 9/18 [00:03<00:01, 4.71it/s]
78%|████████  | 14/18 [00:03<00:00, 8.14it/s]
100%|██████████| 1/1 [00:03<00:00, 3.84s/it]Saving predictions to fda_approved_preds.csv
Elapsed time = 0:00:04
```

```
fda_preds_df = pd.read_csv("fda_approved_preds.csv")
fda_preds_df.head()
```

| | zinc_id | smiles | HIV_active | |
|---|------------------|--|------------|--|
| 0 | ZINC000001530427 | <chem>C[C@@H]1O[C@@H]1P(=O)(O)O</chem> | 0.085422 | |
| 1 | ZINC000003807804 | <chem>Clc1cccc1C(c1cccc1)(c1cccc1)n1ccnc1</chem> | 0.303771 | |
| 2 | ZINC000000120286 | <chem>Nc1nc(N)c2nc(-c3ccccc3)c(N)nc2n1</chem> | 0.093442 | |
| 3 | ZINC000242548690 | <chem>C[C@H]1O[C@@H](O[C@H]2[C@@H](O)C[C@H](O[C@H]3[...])</chem> | 0.565260 | |
| 4 | ZINC000000008492 | <chem>Oc1cccc2cccn12</chem> | 0.089892 | |

Next steps: [View recommended plots](#)

| HIV_active | |  | | | |
|------------|------------------|---|------------|--------------|---|
| count | 892.000000 |  | | | |
| mean | 0.257912 | | | | |
| std | 0.204399 | | | | |
| min | 0.003855 | | | | |
| 25% | 0.097084 | | | | |
| 50% | 0.190803 | | | | |
| 75% | 0.389287 | | | | |
| max | 0.992705 | | | | |
| zinc_id | | smiles | HIV_active | HIV_active_2 |  |
| 0 | ZINC000001530427 | <chem>C[C@@H]1O[C@@H]1P(=O)(O)O</chem> | 0.085422 | 0 | |
| 1 | ZINC000003807804 | <chem>Clc1ccccc1C(c1ccccc1)(c1ccccc1)n1ccnc1</chem> | 0.303771 | 0 | |
| 2 | ZINC000000120286 | <chem>Nc1nc(N)c2nc(-c3ccccc3)c(N)nc2n1</chem> | 0.093442 | 0 | |

```
# Filter rows where 'target_column' is equal to 1
fda_preds_df_filtered = fda_preds_df[fda_preds_df['HIV_active_2'] == 1]
fda_preds_df_filtered
```

| index | zinc_id | smiles | HIV_acti |
|-------|------------------|---|--------------|
| 31 | ZINC000003816287 | CNC(=O)c1ccccc1Sc1ccc2c(/C=C/c3ccccc3)n[nH]c2c1 | 0.8191019296 |
| 47 | ZINC000003813010 | O=c1[nH]c(=O)n([C@H]2C[C@H](O)[C@@H](CO)O2)cc1F | 0.8848749995 |
| 81 | ZINC000003818726 | O=C/C=C/c1cccc(S(=O)(=O)Nc2ccccc2)c1)NO | 0.8046247959 |
| 94 | ZINC000068153186 | CC(C)(C)c1nc(-c2cccc(NS(=O)=O)c3c(F)cccc3F)c2F)c(-c2ccnc(N)n2)s1 | 0.9387590289 |
| 197 | ZINC000000005423 | Cc1nc(-c2ccc(OCC(C)C)c(C#N)c2)sc1C(=O)O | 0.8545335531 |
| 276 | ZINC000000002212 | Cc1nnc2n1-c1ccc(Cl)cc1C(c1ccccc1Cl)=NC2 | 0.8569784164 |
| 321 | ZINC000013597823 | O=c1[nH]cnc2c1ncn2[C@H]1CC[C@H](CO)O1 | 0.851600170 |
| 340 | ZINC000001530621 | CCN[C@H]1C[C@H](C)S(=O)(=O)c2sc(S(N)(=O)=O)cc21 | 0.824345052 |
| 499 | ZINC000000896717 | COc1cc(/C(C)=N/S(=O)(=O)c2ccccc2)ccc1Cc1cn(C)c2ccc(NC(=O)OC3CCCC3)cc12 | 0.8092177510 |
| 540 | ZINC000004474564 | CC/C=C/C/C=C/C=C/C=C/C=C/C=C/C=C/C=C/C=CCCC(=O)O | 0.824715793 |
| 542 | ZINC000005733652 | COc1ccc(-c2cc(=O)c3c(O)cc(O)cc3o2)cc1O | 0.8031748533 |
| 626 | ZINC000003830993 | N[C@@H](Cc1cc(I)c(Oc2cc(I)c(O)c(I)c2)c(I)c1)C(=O)O | 0.8241854906 |
| 715 | ZINC000169289767 | Cc1cc(-c2ccc(/N=N/c3c(S(=O)(=O)O)cc4cc(S(=O)(=O)O)cc(N)c4c3O)c(C)c2)ccc1/N=N/c1c(S(=O)(=O)O)cc2cc(S(=O)(=O)O)cc(N)c2c1O | 0.9927045702 |
| 727 | ZINC000003807172 | C[C@H]1CNc2c(cccc2S(=O)(=O)N[C@@H](CCNC(=N)N)C(=O)N2CC[C@@H](C)C[C@@H]2C(=O)O)C1 | 0.8672307729 |
| 806 | ZINC000000000903 | Cc1nnc2n1-c1ccc(Cl)cc1C(c1ccccc1)=NC2 | 0.8021654486 |
| 819 | ZINC000000137884 | Cc1cn([C@H]2C=C[C@@H](CO)O2)c(=O)[nH]c1=O | 0.9346132278 |
| 821 | ZINC000000897244 | CC1(C)[C@H](C(=O)O)N2C(=O)C[C@H]2S1(=O)=O | 0.8134081363 |
| 879 | ZINC000003779042 | Cc1cn([C@H]2C[C@H](N=[N+]=[N-])[C@@H](CO)O2)c(=O)[nH]c1=O | 0.9854286313 |

```
smiles to check: ['CNC(=O)c1ccccc1Sc1ccc2c(/C=C/c3ccccc3)n[nH]c2c1', 'O=c1[nH]c(=O)n([C@H]2C[C@H](O)[C@@H](CO)O2)
```

smiles activity HIV active smiles activity HIV_active

| num | name | p_np | smiles |
|-----|------|------|--------|
|-----|------|------|--------|

| | zinc_id | smiles |
|-----|------------------|---|
| 31 | ZINC000003816287 | CNC(=O)c1cccc1Sc1ccc2c(/C=C/c3cccn3)n[nH]c2c1 |
| 47 | ZINC000003813010 | O=c1[nH]c(=O)n([C@H]2[C@@H](O)[C@@H](CO)O2)cc1F |
| 81 | ZINC000003818726 | O=C/C=C/c1cccc(S(=O)(=O)Nc2ccccc2c1)NO |
| 94 | ZINC000068153186 | CC(C)(C)c1nc(-c2cccc(NS(=O)(=O)c3c(F)cccc3F)c2... |
| 197 | ZINC000000005423 | Cc1nc(-c2ccc(OCC(C)C)c(C#N)c2)sc1C(=O)O |
| 276 | ZINC000000002212 | Cc1nnc2n1-c1ccc(Cl)cc1C(c1ccccc1Cl)=NC2 |
| 321 | ZINC000013597823 | O=c1[nH]cnc2c1ncn2[C@H]1CC[C@@H](CO)O1 |
| 340 | ZINC000001530621 | CCN[C@H]1C[C@H](C)S(=O)(=O)c2sc(S(N)(=O)=O)cc21 |
| 499 | ZINC000000896717 | COc1cc(/C(O)=N/S(=O)(=O)c2ccccc2C)ccc1Cc1cn(C)... |
| 540 | ZINC000004474564 | CC/C=C\C/C=C\C/C=C\C/C=C\C/C=C\C/C=C\CCC(=O)O |
| 542 | ZINC000005733652 | COc1ccc(-c2cc(=O)c3c(O)cc(O)cc3o2)cc1O |
| 626 | ZINC000003830993 | N[C@@H](Cc1cc(l)c(Oc2cc(l)c(O)c(l)c2)c(l)c1)C(... |
| 715 | ZINC000169289767 | Cc1cc(-c2ccc(/N=N/c3c(S(=O)(=O)O)cc4cc(S(=O)(=... |
| 727 | ZINC000003807172 | C[C@H]1CNc2c(cccc2S(=O)(=O)N[C@@H](CCGN(C=N)N)... |
| 806 | ZINC000000000903 | Cc1nnc2n1-c1ccc(Cl)cc1C(c1ccccc1)=NC2 |
| 819 | ZINC000000137884 | Cc1cn([C@H]2C=C[C@@H](CO)O2)c(=O)[nH]c1=O |
| 821 | ZINC000000897244 | CC1(C)[C@H](C(=O)O)N2C(=O)C[C@H]2S1(=O)=O |
| 879 | ZINC000003779042 | Cc1cn([C@H]2[C@H](N=[N+]=[N-])C[C@@H](CO)O2)c(... |

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1 to 18 of 18 entries

| index | zinc_id_x | smiles | zinc_id_y |
|-------|------------------|--|------------------|
| 0 | ZINC000003816287 | <chem>CNC(=O)c1ccccc1Sc1ccc2c(/C=C/c3ccccc3)n[nH]c2c1</chem> | ZINC000003816287 |
| 1 | ZINC000003813010 | <chem>O=c1[nH]c(=O)n([C@H]2C[C@H](O)[C@@H](CO)O2)cc1F</chem> | ZINC000003813010 |
| 2 | ZINC000003818726 | <chem>O=C(/C=C/c1cccc(S(=O)(=O)Nc2ccccc2)c1)NO</chem> | ZINC000003818726 |
| 3 | ZINC000068153186 | <chem>CC(C)(C)c1nc(-c2cccc(NS(=O)(=O)c3c(F)cccc3F)c2F)c(-c2ccnc(N)n2)s1</chem> | ZINC000068153186 |
| 4 | ZINC000000005423 | <chem>Cc1nc(-c2ccc(OCC(C)C)c(C#N)c2)sc1C(=O)O</chem> | ZINC000000005423 |
| 5 | ZINC000000002212 | <chem>Cc1nnc2n1-c1ccc(Cl)cc1C(c1ccccc1Cl)=NC2</chem> | ZINC000000002212 |
| 6 | ZINC000013597823 | <chem>O=c1[nH]cnc2c1ncn2[C@H]1CC[C@H](CO)O1</chem> | ZINC000013597823 |
| 7 | ZINC000001530621 | <chem>CCN[C@H]1C[C@H](C)S(=O)(=O)c2sc(S(N)(=O)=O)cc21</chem> | ZINC000001530621 |
| 8 | ZINC000000896717 | <chem>COc1cc(/C(O)=N/S(=O)(=O)c2ccccc2C)ccc1Cc1cn(C)c2ccc(NC(=O)OC3CCCC3)cc12</chem> | ZINC000000896717 |
| 9 | ZINC000004474564 | <chem>CC/C=C/C/C=C/C=C/C=C/C=C/C=C/C=C/C=C/C=CCCC(=O)O</chem> | ZINC000004474564 |
| 10 | ZINC000005733652 | <chem>COc1ccc(-c2cc(=O)c3c(O)cc(O)cc3o2)cc1O</chem> | ZINC000005733652 |
| 11 | ZINC000003830993 | <chem>N[C@@H](Cc1cc(I)c(Oc2cc(I)c(O)c(O)c2)c(I)c1)C(=O)O</chem> | ZINC000003830993 |
| 12 | ZINC000169289767 | <chem>Cc1cc(-c2ccc(/N=N/c3c(S(=O)(=O)O)cc4cc(S(=O)(=O)O)cc(N)c4c3O)c(C)c2)ccc1/N=N/c1c(S(=O)(=O)O)cc2cc(S(=O)(=O)O)cc(N)c2c1O</chem> | ZINC000169289767 |
| 13 | ZINC000003807172 | <chem>C[C@H]1CNc2c(cccc2S(=O)(=O)N[C@@H](CCNC(=N)N)C(=O)N2CC[C@@H](C)C[C@@H]2C(=O)O)C1</chem> | ZINC000003807172 |
| 14 | ZINC000000000903 | <chem>Cc1nnc2n1-c1ccc(Cl)cc1C(c1ccccc1)=NC2</chem> | ZINC000000000903 |
| 15 | ZINC000000137884 | <chem>Cc1cn([C@H]2C=C[C@@H](CO)O2)c(O)[nH]c1=O</chem> | ZINC000000137884 |
| 16 | ZINC000000897244 | <chem>CC1(C)[C@H](C(=O)O)N2C(=O)C[C@H]2S1(=O)=O</chem> | ZINC000000897244 |
| 17 | ZINC000003779042 | <chem>Cc1cn([C@H]2C[C@H](N=[N+]=[N-])[C@@H](CO)O2)c(=O)[nH]c1=O</chem> | ZINC000003779042 |

Show 25 per page

Next steps: ☒ View recommended plots

```
fda df final.to_csv('fda approved result.csv', index=False)
```

```
# !wget https://zinc15.docking.org/substances/subsets/named.csv
```

```
--2024-03-10 05:58:28-- https://zinc15.docking.org/substances/subsets/named.csv
Resolving zinc15.docking.org (zinc15.docking.org)... 169.230.75.4
Connecting to zinc15.docking.org (zinc15.docking.org)|169.230.75.4|:443... connected.
HTTP request sent, awaiting response... 200 OK
Length: unspecified [text/csv]
Saving to: 'named.csv.1'

named.csv.1          [ <=>          ]  9.28K  --.-KB/s    in 0.04s

2024-03-10 05:58:29 (242 KB/s) - 'named.csv.1' saved [9499]
```

```
zinc_df = pd.read_csv("named.csv")
zinc_df.head()
zinc_df.tail()
```

| | zinc_id | smiles |
|-------|-------------------|--|
| 0 | ZINC000030727788 | <chem>C=C[C@]1(C)C[C@@H](OC(=O)CSC(C)(C)CNC(=O)[C@H]...</chem> |
| 1 | ZINC000150377216 | <chem>CCCCC/C=C\C/C=C\CCCCCCCC(=O)OC[C@H](CCCCCCCC...</chem> |
| 2 | ZINC000100780125 | <chem>CC(=O)O[C@H]1C[C@](C)(O)[C@@H]2CC=C(C)[C@@H]2[...</chem> |
| 3 | ZINC000006580536 | <chem>O=C(O)[C@H](Cc1ccccc1)N(CCCl)CCCl</chem> |
| 4 | ZINC000150351802 | <chem>O=C1C[C@H](c2ccc(O)c(O)c2)Oc2c1c(O)cc(O[C@H]1O...</chem> |
| | zinc_id | smiles |
| 34595 | ZINC000005999135 | <chem>COc1cc([C@@H]2Oc3c(OC)cc(/C=C/CO)cc3[C@H]2CO)c...</chem> |
| 34596 | ZINC000084710404 | <chem>COC(=O)c1cc(OC)c2cc(OC)c(OC)c(O)c2c1O</chem> |
| 34597 | ZINC000150369761 | <chem>CC/C=C\C/C=C\C/C=C\C/C=C\C/C=C\C/C=C\C\CCC(=O)O[...</chem> |
| 34598 | ZINC000095098911 | <chem>CCCCCCCCCCCC[C@H](O)[C@H]1CC[C@H]([C@H](O)CCCC...</chem> |
| 34599 | ZINC0000000001009 | <chem>C[N+](C)([O-])CC/C=C1c2ccccc2C=Cc2c(Cl)cccc21</chem> |

```
arguments = [
    '--test_path', 'named.csv',
    '--preds_path', 'named_preds.csv',
    '--checkpoint_dir', 'test_checkpoints_multimolecule',
    '--smiles_columns', 'smiles'
]
```

```
args = chemprop.args.PredictArgs().parse_args(arguments)
preds = chemprop.train.make_predictions(args=args)
```

```
84% ██████████ | 581/692 [02:16<00:19, 5.65it/s]
84% ██████████ | 584/692 [02:16<00:17, 6.35it/s]
85% ██████████ | 586/692 [02:17<00:18, 5.70it/s]
85% ██████████ | 587/692 [02:17<00:25, 4.09it/s]
85% ██████████ | 589/692 [02:17<00:21, 4.71it/s]
85% ██████████ | 591/692 [02:18<00:16, 6.06it/s]
86% ██████████ | 592/692 [02:18<00:24, 4.00it/s]
86% ██████████ | 594/692 [02:19<00:22, 4.31it/s]
86% ██████████ | 595/692 [02:20<00:36, 2.64it/s]
86% ██████████ | 597/692 [02:20<00:25, 3.71it/s]
87% ██████████ | 600/692 [02:20<00:18, 4.89it/s]
87% ██████████ | 602/692 [02:20<00:16, 5.30it/s]
87% ██████████ | 603/692 [02:21<00:23, 3.74it/s]
87% ██████████ | 605/692 [02:21<00:17, 4.87it/s]
88% ██████████ | 608/692 [02:22<00:14, 5.78it/s]
88% ██████████ | 610/692 [02:22<00:12, 6.42it/s]
88% ██████████ | 611/692 [02:23<00:26, 3.00it/s]
89% ██████████ | 613/692 [02:23<00:19, 3.96it/s]
89% ██████████ | 616/692 [02:24<00:14, 5.10it/s]
89% ██████████ | 618/692 [02:24<00:12, 5.87it/s]
89% ██████████ | 619/692 [02:24<00:18, 3.92it/s]
90% ██████████ | 621/692 [02:25<00:14, 4.98it/s]
90% ██████████ | 624/692 [02:25<00:12, 5.66it/s]
90% ██████████ | 626/692 [02:25<00:11, 5.85it/s]
91% ██████████ | 627/692 [02:26<00:15, 4.28it/s]
91% ██████████ | 629/692 [02:26<00:11, 5.68it/s]
91% ██████████ | 632/692 [02:26<00:09, 6.55it/s]
92% ██████████ | 634/692 [02:27<00:08, 6.71it/s]
92% ██████████ | 635/692 [02:27<00:12, 4.41it/s]
92% ██████████ | 637/692 [02:27<00:09, 5.58it/s]
92% ██████████ | 640/692 [02:28<00:08, 6.09it/s]
93% ██████████ | 642/692 [02:28<00:07, 6.39it/s]
93% ██████████ | 643/692 [02:29<00:11, 4.33it/s]
93% ██████████ | 646/692 [02:29<00:06, 6.69it/s]
94% ██████████ | 648/692 [02:29<00:06, 6.46it/s]
94% ██████████ | 650/692 [02:30<00:06, 6.05it/s]
94% ██████████ | 651/692 [02:31<00:18, 2.25it/s]
94% ██████████ | 653/692 [02:32<00:12, 3.03it/s]
95% ██████████ | 655/692 [02:32<00:08, 4.13it/s]
95% ██████████ | 657/692 [02:32<00:08, 4.10it/s]
95% ██████████ | 658/692 [02:32<00:08, 4.21it/s]
95% ██████████ | 659/692 [02:33<00:10, 3.02it/s]
96% ██████████ | 661/692 [02:33<00:07, 4.17it/s]
96% ██████████ | 664/692 [02:34<00:05, 5.20it/s]
96% ██████████ | 666/692 [02:34<00:04, 5.65it/s]
96% ██████████ | 667/692 [02:34<00:06, 3.93it/s]
97% ██████████ | 669/692 [02:35<00:04, 5.17it/s]
97% ██████████ | 672/692 [02:35<00:03, 5.79it/s]
97% ██████████ | 674/692 [02:35<00:02, 6.22it/s]
98% ██████████ | 675/692 [02:36<00:04, 4.22it/s]
98% ██████████ | 677/692 [02:36<00:02, 5.55it/s]
98% ██████████ | 680/692 [02:36<00:01, 6.79it/s]
99% ██████████ | 683/692 [02:37<00:01, 8.29it/s]
99% ██████████ | 688/692 [02:37<00:00, 13.13it/s]
100% ██████████ | 691/692 [02:37<00:00, 14.99it/s]
100% ██████████ | 1/1 [02:38<00:00, 158.26s/it]
```

Saving predictions to named_preds.csv

Elapsed time = 0:02:52

```
zinc_preds_df = pd.read_csv("named_preds.csv")
zinc_preds_df.head()
zinc_preds_df = zinc_preds_df[zinc_preds_df['HIV_active'] != "Invalid SMILES"]
zinc_preds_df.describe()
zinc_preds_df['HIV_active'] = zinc_preds_df['HIV_active'].astype(float)
zinc_preds_df['HIV_active_2'] = zinc_preds_df['HIV_active'].apply(lambda x: 1 if x > 0.8 else 0)
zinc_preds_df.head()
```


| | zinc_id | smiles | HIV_active | |
|---|------------------|--|------------|--|
| 0 | ZINC000030727788 | <chem>C=C[C@]1(C)C[C@@H](OC(=O)CSC(C)(C)CNC(=O)[C@H]...</chem> | 0.345203 | |
| 1 | ZINC000150377216 | <chem>CCCCC/C=C/C/C=C\CCCCCCCC(=O)OC[C@H](COCCCCCCC...</chem> | 0.619356 | |
| 2 | ZINC000100780125 | <chem>CC(=O)O[C@H]1C[C@](C)(O)[C@@H]2CC=C(C)[C@@H]2[...</chem> | 0.260698 | |
| 3 | ZINC000006580536 | <chem>O=C(O)[C@H](Cc1ccccc1)N(CCCl)CCCl</chem> | 0.043098 | |
| 4 | ZINC000150351802 | <chem>O=C1C[C@H](c2ccc(O)c(O)c2)Oc2c1c(O)cc(O[C@H]1O...</chem> | 0.716069 | |

| | HIV_active | |
|-------|--------------|--|
| count | 34600.000000 | |
| mean | 0.466442 | |
| std | 0.243481 | |
| min | 0.000041 | |
| 25% | 0.251037 | |
| 50% | 0.499464 | |
| 75% | 0.673119 | |
| max | 0.998601 | |

Next steps:

zinc_id

View recommended plots

smiles

View recommended plots

HIV_active

View recommended plots

HIV_ac

View recommended plots

```
0 ZINC000030727788 C=C[C@]1(C)C[C@@H](OC(=O)CSC(C)(C)CNC(=O)[C@H]... 0.345203
# Filter rows where 'target_column' is equal to 1
zinc_preds_df_filtered = zinc_preds_df[zinc_preds_df['HIV_active_2'] == 1]
zinc_preds_df_filtered
```

1 to 25 of 2092 entries

| index | zinc_id | smiles | HIV_active | HIV_active_2 |
|-------|------------------|--|--------------------|--------------|
| 21 | ZINC000040753343 | <chem>COc1cc(/C=C/c2cc(O)c(CC=C(C)C)c(O)c2)cc2c1O[C@]1(C)CC[C@@H](O)C(C)(C)[C@H]1C2</chem> | 0.9342743158340454 | 1 |
| 37 | ZINC000014615844 | <chem>CC1(C)OC(=O)C=C[C@@]2(C)[C@@H]1CC(=O)[C@@]1(C)[C@@H]2[C@@H](O)[C@@]2(C)[C@@H](c3ccoc3)OC(=O)[C@@H]3O[C@@]321</chem> | 0.8453148603439331 | 1 |
| 91 | ZINC000150366864 | <chem>COc1cc(-c2[+][c3cc(O)cc(O)c3cc2O[C@H]2O[C@H](CO[C@H]3O[C@@H](CO)[C@@H](O)[C@@H](O)[C@@H]3O)[C@@H](O)[C@@H](O)[C@@H]2O)cc(O)c1O</chem> | 0.8436190485954285 | 1 |
| 110 | ZINC000049888739 | <chem>CCCCC(=O)c1c(O)c2c(c3c1O[C@@]1(O)[C@H]([C@@H]3C(C)C)[C@@H](O)C(C)(C)C(=O)C1(C)C)OC1=C(C(=O)C(C)C(=O)C1(C)C)[C@@H]2C(C)C</chem> | 0.8854982256889343 | 1 |
| 124 | ZINC000150343906 | <chem>CCCCC/C=C/C/C=C\CCCCCCCC(=O)OC[C@H](CO[P@@](=O)(O)O)[C@H]1[C@@H](O)[C@@H](O)[C@@H](O)[C@@H](O)[C@@H](OP(=O)(O)O)[C@@H]1O)OC(=O)CCCCCCCCCCCCCCCCC</chem> | 0.8228963017463684 | 1 |
| 128 | ZINC000008220075 | <chem>Nc1nc(O)c(N)c(N[C@@H]2O[C@H](CO[P@@](=O)(O)O)[P@@](=O)(O)OP(=O)(O)O)[C@@H](O)[C@H]2O)n1</chem> | 0.9743053317070008 | 1 |
| 138 | ZINC000096888407 | <chem>CC(C)[C@H]1C=C[C@]2(C)Oc3c(C(=O)/C=C/c4ccccc4)c(O)cc(O)c3[C@@H]2C1</chem> | 0.8139744997024536 | 1 |
| 149 | ZINC000014680924 | <chem>Cc1c(Cl)c(O)cc(O)c1C(=O)O[C@H]1C[C@]2(C)[C@@H]3[C@@H](O)C(C)(C)C[C@@]3(O)C=C(C=O)[C@@]12O</chem> | 0.9280908703804016 | 1 |
| 165 | ZINC000150369401 | <chem>Oc1ccc([C@@H]2Oc3ccc(/C=C/C4cc(O)[C@H]2O[C@H](c4cc(O)cc(O)c4)[C@H](c4ccc(O)cc4)O5)cc3[C@@H]2c2cc(O)cc3c2[C@H](c2cc(O)cc(O)c2)[C@H](c2ccc(O)cc2)O3)cc1</chem> | 0.9324991106987 | 1 |
| 185 | ZINC000230100259 | <chem>O=C(O)[C@H]1Cc2c(O)cc(O)cc2O[C@H]1c1cc(O)c(O)c(O)c1-c1c(OC(=O)[C@H]2C[C@@](O)(C(=O)O)C[C@@H](O)[C@H]2O)cc(O)c(O)c1O)c1cc(O)c(O)c(O)c1</chem> | 0.9309805631637572 | 1 |
| 194 | ZINC000082170194 | <chem>O=C(OC[C@H]1OC[C@H](OC(=O)c2cc(O)c(O)c(O)c2)[C@@H](O)[C@@H]1OC(=O)c1cc(O)c(O)c(O)c1)c1cc(O)c(O)c(O)c1</chem> | 0.9685450196266174 | 1 |
| 206 | ZINC000013380011 | <chem>C[C@H]1Oc2cc(O)c3c(=O)c(O)c(-c4cc(O)c5c(c4)C=CC(C)(C)O5)oc3c2C1(C)C</chem> | 0.9307474493980408 | 1 |
| 208 | ZINC000085742375 | <chem>O=C(/C=C/C1ccc(O)cc1)O[C@@H]1[C@@H](CO)O[C@H](Oc2c(-c3cc(O)cc3)oc3cc(O)cc(O)c3c2=O)[C@H](OC(=O)/C=C/c2ccc(O)cc2)[C@@H]1O</chem> | 0.8347569704055786 | 1 |
| 218 | ZINC000033979236 | <chem>O=c1ccn([C@H]2C[C@@H](O)[C@H](CO[P@@](=O)(O)OP(=O)(O)O)O2)c(=O)[nH]1</chem> | 0.98438960313797 | 1 |
| 232 | ZINC000049089629 | <chem>C[C@@H]1CC(=O)[C@H]2C[C@]1(O)[C@]2(C)COC(=O)c1cc(O)c(O)c(O)c1</chem> | 0.9498233199119568 | 1 |
| 297 | ZINC000230100275 | <chem>O=C(O)[C@H]1Cc2c(O)cc(O)cc2O[C@H]1c1cc(O)c(O)c(O)c1-c1c(OC(=O)[C@H]2C[C@@](O)(C(=O)O)C[C@@H](O)[C@H]2O)cc(O)c(O)c1O)c1cc(O)c(O)c(O)c1</chem> | 0.928594410419464 | 1 |
| 300 | ZINC000014727558 | <chem>CC(C)=CCc1c(-c2ccc(O)cc2O)cc2c(CC=C(C)C)c(O)cc(O)c2c1=O</chem> | 0.8761184215545654 | 1 |
| 309 | ZINC000040861309 | <chem>C=C(C)[C@H]1CC=C(CO[P@@](=O)(O)OC[C@@H](COC(=O)OC(=O)CC1</chem> | 0.8323433995246887 | 1 |
| 315 | ZINC000195445287 | <chem>CCCCCCCCCCCCC(=O)SCCNC(=O)CCNC(=O)[C@H](O)C(C)(C)CO[P@@](=O)(O)O[P@@](=O)(O)OC[C@H]1O[C@@H](n2cnc3c(N)ncnc32)[C@H](O)[C@@H]1OP(=O)(O)O</chem> | 0.8638664484024048 | 1 |
| 328 | ZINC000169339170 | <chem>O=C1OC[C@@H]2OC(=O)c3cc(O)c(O)c(O)c3-c3c(O)c(O)c(O)c4c3C(=O)O[C@H]([C@@H]2OC(=O)c2cc(O)c(O)c(O)c2-c2c1cc(O)c(O)c2O)[C@@H]1OC(=O)c2c-4c(O)c(O)c(O)c2[C@@H]1O</chem> | 0.9289804697036744 | 1 |