In [ ]:
```
! pip install chembl_webresource_client
```

In [2]:
```
# Import necessary libraries
import pandas as pd
from chembl_webresource_client.new_client import new_client
```

In [3]:
```
#convert pubchemID to ChemBL ID
df_fla=pd.read_csv("fdb_molecules_22Nov.csv")
df_fla.head()
```

Out[3]:

	pubchem_id	molecular_weight	hbd_count	hba_count	xlogp	natural
0	4	75.111	2	2	-1.0	1
1	40	164.157	4	5	-1.6	0
2	47	130.143	1	3	1.1	0
3	49	116.116	1	3	0.7	1
4	51	146.098	2	5	-0.9	0

In [4]:
```
df_fla.info()
```

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 25595 entries, 0 to 25594
Data columns (total 6 columns):
 #   Column            Non-Null Count  Dtype  
---  ------            --------------  -----  
 0   pubchem_id        25595 non-null  int64  
 1   molecular_weight  25595 non-null  float64
 2   hbd_count         25595 non-null  int64  
 3   hba_count         25595 non-null  int64  
 4   xlogp             22932 non-null  float64
 5   natural           25595 non-null  int64  
dtypes: float64(2), int64(4)
memory usage: 1.2 MB
```

In [6]:
```
df_nat=df_fla.loc[df_fla['natural'] == 1]
df_nat.reset_index(drop=True, inplace=True)
df_nat.info()
df_nat.head()
```

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 2254 entries, 0 to 2253
Data columns (total 6 columns):
```

```
 #   Column            Non-Null Count  Dtype
---  ------            --------------  -----
 0   pubchem_id        2254 non-null   int64
 1   molecular_weight  2254 non-null   float64
 2   hbd_count         2254 non-null   int64
 3   hba_count         2254 non-null   int64
 4   xlogp             2203 non-null   float64
 5   natural           2254 non-null   int64
dtypes: float64(2), int64(4)
memory usage: 105.8 KB
```

Out[6]:

	pubchem_id	molecular_weight	hbd_count	hba_count	xlogp	natural
0	4	75.111	2	2	-1.0	1
1	49	116.116	1	3	0.7	1
2	58	102.089	1	3	0.1	1
3	70	130.143	1	3	0.9	1
4	72	154.121	3	4	1.1	1

Total fdb molecules were 25595 , natural molecules were only 2254 after conversion to chembl ids using Pubchem Identifier Exchange Service- Filename- "pubchem_tochembl.csv" . Only 1145 ids were left

In [10]:
```
df_nat.to_csv('fdb_natural_2254.csv', index=False)
```

In [ ]:
```
#####comverted pubchem is to chembl ids####
```

In [11]:
```
df_chem=pd.read_csv("pubchem_tochembl.csv")
df_chem.head()
```

Out[11]:

	pubchem_id	chembl_id
0	4	CHEMBL326602
1	49	CHEMBL146554
2	58	CHEMBL171246
3	70	CHEMBL445647
4	72	CHEMBL37537

In [12]:
```
df_chem.info()
```

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 2260 entries, 0 to 2259
Data columns (total 2 columns):
 #   Column      Non-Null Count  Dtype 
---  ------      --------------  ----- 
 0   pubchem_id  2260 non-null   int64 
 1   chembl_id   1145 non-null   object
dtypes: int64(1), object(1)
memory usage: 35.4+ KB
```

In [13]:
```
df_chem_nona=df_chem[df_chem.chembl_id.notna()]
df_chem_nona.reset_index(drop=True, inplace=True)
#df_chem_nona.chembl_id.str.strip()
df_chem_nona.info()
df_chem_nona.head()
```

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 1145 entries, 0 to 1144
Data columns (total 2 columns):
 #   Column      Non-Null Count  Dtype 
---  ------      --------------  ----- 
 0   pubchem_id  1145 non-null   int64 
 1   chembl_id   1145 non-null   object
dtypes: int64(1), object(1)
memory usage: 18.0+ KB
```

Out[13]:

	pubchem_id	chembl_id
0	4	CHEMBL326602
1	49	CHEMBL146554
2	58	CHEMBL171246
3	70	CHEMBL445647
4	72	CHEMBL37537

In [14]:
```
merge_dfchem_pub=df_nat.merge(df_chem_nona, on ="pubchem_id", how = "inner")
merge_dfchem_pub.info()
merge_dfchem_pub.head()
```

```
<class 'pandas.core.frame.DataFrame'>
Int64Index: 1145 entries, 0 to 1144
Data columns (total 7 columns):
 #   Column            Non-Null Count  Dtype 
---  ------            --------------  ----- 
```

```
 0   pubchem_id        1145 non-null   int64
 1   molecular_weight  1145 non-null   float64
 2   hbd_count         1145 non-null   int64
 3   hba_count         1145 non-null   int64
 4   xlogp             1126 non-null   float64
 5   natural           1145 non-null   int64
 6   chembl_id         1145 non-null   object
dtypes: float64(2), int64(4), object(1)
memory usage: 71.6+ KB
```

Out[14]:

	pubchem_id	molecular_weight	hbd_count	hba_count	xlogp	natural	
0	4	75.111	2	2	-1.0	1	CH
1	49	116.116	1	3	0.7	1	CH
2	58	102.089	1	3	0.1	1	CH
3	70	130.143	1	3	0.9	1	CH
4	72	154.121	3	4	1.1	1	C

In [10]:

```
molecules=list(merge_dfchem_pub['chembl_id'])
df2=merge_dfchem_pub
df2.to_csv('df2.csv', index= False)
```

In [11]:

```
import pandas as pd # uses pandas python module to view and analyse
data
import requests # this is used to access json files

#=====================================================================
==#

# using a list of natural fdb chembl ids bioactivity data scraping,
Limit=5 records for each CHembl Id were scraped with Standard_type-
IC50
# standard_type IC50 was selected , since it is most widely used fo
r bioactivity prediction.

def find_bioactivities_for_molecules(molecules):

    limit = 5 # limit the number of records pulled back for each ur
l call xxx
    standard_type='IC50'
    url_stem = "https://www.ebi.ac.uk" #This is the stem of the url
    url_full_string = url_stem + "/chembl/api/data/activity.json?mo
lecule_chembl_id__in={}&limit={}&standard_type={}".format(molecules
,limit,standard_type) #This is the full url with the specified inpu
```

```
t parameters
    url_full = requests.get( url_full_string ).json() #This calls t
he information back from the API using the 'requests' module, and c
onverts it to json format
    url_activities = url_full['activities'] #This is a list of the
 results for activities

    # This 'while' loop iterates over several pages of records (if
 required), and collates the list of results
    while url_full['page_meta']['next']:
        url_full = requests.get(url_stem + url_full['page_meta']['n
ext']).json()
        url_activities = url_activities + url_full['activities'] #A
dd result (as a list) to previous list of results

    # Convert the list of results into a Pandas dataframe:
    act_df = pd.DataFrame(url_activities)

    # Print out some useful information:
    print("This is the url string that calls the 'Activities' API w
ith the initial query specification:\n{}".format(url_full_string) )
    print("\nThese are the available columns for the Activities AP
I:\n{}".format(act_df.columns))

    #Specify which columns to keep so that the size of the datafram
e becomes more manageable:
    act_df = act_df[[  'target_chembl_id','target_organism', 'targe
t_pref_name'
                     ,'molecule_chembl_id','molecule_pref_name'
                     , 'pchembl_value', 'standard_type','standard_r
elation', 'standard_value', 'standard_units'
                     , 'assay_chembl_id']]

    return act_df
```

In [12]:

```
dic={}
df=pd.DataFrame()
id_list=[]
for i in range(len(df2)):
    k=df2.chembl_id[i]
    #dic["%s"%(k)]=find_bioactivities_for_molecules(k)
    try:
        data=find_bioactivities_for_molecules(k)
        df=df.append(data)
    #except KeyError as i:
    #print('Key Not Found in Employee Dictionary:',i)
    except KeyError:
```

```
        continue
    #id_list.append(dic["%s"%(i)])
#print(len(id_list))
```

```
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL326602&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL146554&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL171246&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
```

```
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL445647&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL37537&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL851&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
```

```
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL96&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
```

```
embl_id__in=CHEMBL202132&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14193&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
```

```
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL441343&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL539&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
```

```
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL170365&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL16081&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3561873&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14253&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
```

```
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1157&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
```

```
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL297569&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15972&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
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ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL541&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL720&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1182&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
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      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1478334&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2312529&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14245&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14227&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL116736&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1261&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1231821&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL6466&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183500&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL161577&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1232595&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL46931&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL424&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
```

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e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL298312&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL134886&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL324846&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1200739&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2105487&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL18407&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
```

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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL13766&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1233860&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL275626&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL12198&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1200559&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL365809&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL89306&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1229937&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL120433&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL504&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL545&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
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       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1162495&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL773&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL692&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL183419&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL71595&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
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d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL537&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL297800&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15844&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL82411&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
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mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL333714&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL274119&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL170721&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14688&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL16293&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL32571&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
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       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL26215&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL146755&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL82293&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL25080&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14060&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1233464&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1044&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL610&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1187&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
```

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       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL286398&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14687&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14021&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL266158&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1162144&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15580&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL23894&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL23194&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL576&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1235997&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL239243&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1547&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1236482&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL9113&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
```

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       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL439723&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL11257&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL13883&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL256087&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL328910&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
```

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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL271939&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1967951&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL346919&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
```

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       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL7303&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL93353&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1239&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL55285&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL113&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
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d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL706&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15134&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
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       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL485259&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1231862&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1397305&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL16645&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL107498&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL75124&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL272485&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL90039&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL42710&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
```

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       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1950582&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL48310&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
```

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       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL582&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL110309&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108766&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL416&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
```

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       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108545&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL481044&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL174778&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL8320&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
```

```
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL19224&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL45068&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL7983&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
```

```
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL24147&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL242273&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
```

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mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL46403&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1114&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1489254&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL11608&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1276010&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
```

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       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1682&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1222250&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
```

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      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL863&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL130&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL274323&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL930&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL8085&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
```

```
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL233248&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL253582&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
```

```
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL118722&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL448500&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL925&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL245067&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL752&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
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       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL291962&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL417016&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL301523&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL107&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL280331&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
```

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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1200941&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL25028&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
```

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e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1355&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL689&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1908365&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL17962&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14568&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL54976&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
```

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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1485&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14449&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL45967&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1743219&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL117080&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1897156&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL44658&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1439452&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL506184&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL464988&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185164&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1882894&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183871&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL25306&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL269630&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL1404017&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL45462&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15849&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14079&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108778&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
```

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e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2268550&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1651998&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1414114&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1513871&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL442565&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1329793&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL388558&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1370662&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1409791&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2260718&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1332922&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL184290&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1232797&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL401911&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
```

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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL30018&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL541939&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
```

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ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1719455&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2252746&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1898498&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL29411&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
```

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mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108925&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL383808&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL109652&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1668603&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14474&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
```

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       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL195895&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1412508&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14092&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108861&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
```

```
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108299&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182715&limit=5&standard_type=IC50
```

```
These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3730701&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1397202&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL193446&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL16435&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL379845&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL510714&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1200370&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2260711&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2252208&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL573781&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL510309&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
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```
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL45005&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1797280&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL192008&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL192591&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL2270060&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL45315&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL95681&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL1234459&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL29873&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL1351619&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:

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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1204680&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL295870&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186323&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188433&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL273782&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2287521&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL308187&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL189362&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3560314&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL328441&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
```

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mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL16217&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL508917&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL274467&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1385229&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15676&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL325372&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188459&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL449693&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2251642&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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```
embl_id__in=CHEMBL442915&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL29966&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL201083&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL401912&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
```

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       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL21932&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL371561&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL285235&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1224557&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL278024&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL192899&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL279305&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183743&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3189123&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL38934&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183476&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183441&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2252087&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3561355&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183179&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184059&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184025&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184929&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL152299&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187974&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL286062&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184606&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185876&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3559967&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL440161&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
```

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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL191935&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL195827&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL31637&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL190927&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
```

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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL227181&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL154155&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL294431&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL365316&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
```

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d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL44115&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1789215&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL177114&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL44800&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1338583&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL276218&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184928&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3181881&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1907995&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL447944&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187370&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184220&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187842&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185284&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL31561&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
```

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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL556489&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL537970&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL721&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
```

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d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1236818&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14409&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL234926&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
```

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       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186027&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL42003&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL45345&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186475&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL285323&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1608674&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1305819&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL195593&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
```

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       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL286727&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL22976&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1986332&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183097&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1797281&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187831&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182186&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL16200&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL32010&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL18850&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL119405&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15732&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
```

embl_id__in=CHEMBL479791&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL268736&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL44857&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL16102&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL13968&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3188256&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL1189&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:

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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL44215&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL16225&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL276521&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL46999&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL207602&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184829&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2137647&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL18893&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
```

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       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15939&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL18602&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2270393&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL15487&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183908&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL18549&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL320358&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14085&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL153339&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
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       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL273459&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL18104&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL1591973&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1894365&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL335900&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108436&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228455&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2141776&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1236582&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228456&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228458&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2270392&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL25363&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228377&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108030&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
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ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL335125&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL30959&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL444525&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1236576&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185269&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL24722&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL2228373&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL2228460&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL24832&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL24022&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL4245903&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL502773&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:

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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1160012&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL460124&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2251610&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL31422&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
```

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s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187012&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

```
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL12315&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL333179&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1088937&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
```

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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185936&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15841&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL271663&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
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       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL18620&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL508676&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL120568&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1762668&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL271871&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL387326&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1493986&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL225303&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL396295&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1868953&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182595&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL451923&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183988&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184678&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2104112&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL222021&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL1233714&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3560735&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL501246&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL470671&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL52084&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185956&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185221&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL395827&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
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      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14152&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185916&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184842&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL169176&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188902&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL14184&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL134658&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL139055&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228454&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL24563&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228457&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15245&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL80257&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL402812&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL544&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
```

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       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1668604&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1453648&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3907604&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL2252486&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL292303&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL311498&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
```

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d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15605&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15797&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1366925&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187172&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL117181&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3039186&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL46340&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1952257&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL116960&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2288022&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL470874&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL458690&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1672002&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL370688&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL166223&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4169727&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL281202&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL53566&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2288023&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL568737&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL332887&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183037&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL453797&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL443408&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3120653&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1560118&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL275638&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1487138&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL49732&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
```

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       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183607&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1445555&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL444711&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL23025&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL47483&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182933&limit=5&standard_type=IC50
```

```
These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187784&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183048&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL115725&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL45235&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2251727&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL449522&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL3561568&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL111077&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL134994&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2251601&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3183973&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3039425&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL444254&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL396000&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3184952&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL507795&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3183920&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')

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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL30106&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL192258&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1946183&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL454697&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184187&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL452729&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL487998&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3561586&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL450417&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL18360&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL478640&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL22585&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1235931&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL195861&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1814588&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2270394&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL170458&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
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       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL480097&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3360549&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL443470&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL192809&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2269086&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL276849&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
```

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       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL173373&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2230304&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL225153&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
```

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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1522321&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL46678&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15859&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188103&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1347061&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184082&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1416448&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL45425&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182412&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL47769&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3561042&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL454517&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL135694&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL135488&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1234557&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182381&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL26561&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185332&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL107874&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL532&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187976&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL225569&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL35297&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1230308&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4172502&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3189076&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182950&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL47127&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1566946&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL192458&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
```

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ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182050&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL95973&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182189&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL365740&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3561140&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2106111&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228473&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187896&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1200458&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2287316&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3222024&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL462861&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
```

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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2260721&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL460025&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL56395&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2229658&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1414126&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2424841&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182596&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186281&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL132474&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL11945&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL303697&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
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       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4218890&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1789306&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3560443&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3139347&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL455506&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2268554&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186822&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2104397&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL501351&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL441687&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
```

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       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3617994&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187231&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1487817&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1506228&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1907996&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1882821&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL134537&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2252095&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL123040&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2260713&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL469537&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4162810&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL366603&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187351&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL470670&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
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mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL501949&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL151649&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
```

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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL195215&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1355077&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187477&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182738&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL449810&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182794&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL290007&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186884&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL329881&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1236925&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL97794&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188615&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2269084&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL333298&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187568&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL452687&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183573&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL3187919&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL173521&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186302&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182720&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3187497&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL510068&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL1331840&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL1789229&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL1887227&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL2270634&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL2260710&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')

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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3234723&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL105912&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL96425&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2268553&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL15799&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3222023&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL45065&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228472&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187323&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2260717&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL333973&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3305989&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188292&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1558498&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1644029&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2106566&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1206211&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL3561573&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228568&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL94558&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL95123&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL506854&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186524&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL97525&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL95102&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL328856&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL276311&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL320146&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2271411&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187840&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL327303&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL94736&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL254585&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1904383&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3273568&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1159645&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1730107&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3560993&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188662&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183519&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL25894&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185609&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1867360&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL473366&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1606391&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL3187866&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108475&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL161598&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
```

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d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2270061&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL369243&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187336&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL94709&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL455491&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182048&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL372396&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2229585&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL22830&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL284391&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL42013&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL207555&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL207549&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228376&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228569&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185292&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL146&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
```

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       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL257332&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188899&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188160&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL97355&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL575060&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL328465&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL96852&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186540&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL503899&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188696&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182352&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1991946&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3275138&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184459&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL319497&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2260709&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL364588&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185401&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1303827&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184371&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184983&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL453815&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182849&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185940&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183495&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL321211&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
```

embl_id__in=CHEMBL2260714&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3186579&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3188181&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3186660&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3184950&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL1374805&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3560312&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:

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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL47155&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2252121&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2270391&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL372447&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL371371&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL225343&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185204&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3273567&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2268884&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186486&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL248594&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
```

```
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185758&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188412&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL358850&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1894194&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188960&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183686&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL119603&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188298&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186212&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182614&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
```

embl_id__in=CHEMBL3186269&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL1789171&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3185445&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3183829&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3183410&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3181874&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_chembl_id__in=CHEMBL3188389&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the initial query specification:

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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184429&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4297450&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186484&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL264141&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
```

```
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3728427&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187606&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182254&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL349881&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186900&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1812853&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186512&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1870000&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185688&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1097205&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1614854&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

```
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL297453&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL125743&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
```

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       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL266625&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL517016&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
```

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      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1974890&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1409937&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL372105&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL450288&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL118062&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4288485&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL73930&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL108862&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
```

```
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3593941&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1079147&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL21824&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1814589&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

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d', 'type',
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      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1232386&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL350966&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
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      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2272092&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL285123&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL66693&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL104875&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1076637&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
```

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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
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mbl_value',
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ion', 'src_id',
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ue',
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e',
       'standard_value', 'target_chembl_id', 'target_organism',
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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL379630&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1201068&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL583912&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL47386&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
```

```
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL366460&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
```

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embl_id__in=CHEMBL76447&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2368547&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
```

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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1080178&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2208211&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1081721&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1947232&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL97593&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3593942&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL298717&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3593576&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2285739&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2151143&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL277871&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL464709&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182701&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1236329&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL29757&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL118504&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
```

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       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1234268&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185240&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3593943&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186132&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2004717&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
```

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       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL88244&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL198877&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
```

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       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2268547&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1236183&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2287241&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4210821&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
```

```
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL510535&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL518542&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3120654&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL469654&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL171281&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1990994&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1988732&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3189103&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2071440&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3120645&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3560208&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1592541&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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```
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1968792&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183970&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3732724&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2289230&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1907993&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3109299&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1530328&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3728760&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186375&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL448805&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2229207&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
```

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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL330546&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL36327&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
```

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d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL486422&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183581&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL19344&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1408113&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL52416&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL226507&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL54922&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1232207&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
```

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d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL303483&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL573448&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
```

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       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1661&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL504760&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
```

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mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228955&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL487213&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
```

```
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL574688&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL237994&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
```

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ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2288883&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL517449&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3593944&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3593947&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2236867&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL221542&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
```

```
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL349605&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2059292&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL226915&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2004366&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3560504&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL185885&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL70518&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183593&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL205268&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2269087&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187980&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186295&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1444078&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL364713&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
```

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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1522900&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL54161&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2007486&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2001434&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3273403&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL9352&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL236688&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1907991&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL300520&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1233058&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL462997&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1232258&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
```

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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL66&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2229268&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1232369&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL449062&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL232202&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184774&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3109297&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL445759&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1614877&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2093058&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1519430&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL227934&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
```

```
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL218693&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL47244&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

```
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1096927&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2287242&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL451532&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL455041&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2924219&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2426622&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL235672&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2268548&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL447597&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL415690&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
```

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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL153658&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL27246&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL503160&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
```

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       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL25308&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
```

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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL165&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL453509&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL8659&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL501174&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1086445&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL63558&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL460647&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL492828&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
```

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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1354&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
```

```
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1356&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183483&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188345&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL117865&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL94514&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1589555&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183389&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL93500&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185064&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4159021&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL333306&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3356397&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL293492&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
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d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL460657&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL55060&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL66879&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
```

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       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL452630&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1576086&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL25719&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL318196&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL73639&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL109341&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
```

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       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL465829&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL487603&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4099976&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1080997&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL559945&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL43185&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL486193&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
```

```
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL115668&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL129795&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184169&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL253896&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187813&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL452173&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL83159&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL249592&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
```

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e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3126829&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL347285&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
```

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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL452683&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL55941&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL105424&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2286136&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL193598&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL304461&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186705&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
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       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL446299&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL294199&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2251602&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2268549&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1369384&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1468832&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL445206&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1164609&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1994615&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL398921&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
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embl_id__in=CHEMBL2387747&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1449245&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1095954&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL486207&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL508894&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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e',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL399036&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
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       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1873035&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2251453&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL170&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1200429&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228956&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL190613&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL52267&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
```

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e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182150&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2165230&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1900511&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL361197&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL170190&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3883497&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL50&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1644111&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL28&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
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       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL151&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL267476&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL463088&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL118958&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3120646&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL214321&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL250450&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL150&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
```

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       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL44&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2229606&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL299599&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2251452&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228570&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL128000&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL486625&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL12014&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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e',
```

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       'standard_value', 'target_chembl_id', 'target_organism',
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      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL445740&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182226&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL251280&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
```

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s',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1911053&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
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mbl_value',
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ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
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d', 'type',
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228374&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL117&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL242383&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
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ion', 'src_id',
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ue',
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e',
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d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL31574&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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ion', 'src_id',
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ue',
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e',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
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      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL210635&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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ion', 'src_id',
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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL379064&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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ion', 'src_id',
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ue',
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e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL28626&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
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ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
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ue',
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e',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL413552&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL55415&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
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e',
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       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
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```
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL247484&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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ion', 'src_id',
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ue',
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e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL30707&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
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ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
```

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mbl_value',
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ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL8145&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL491174&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
```

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s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
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ion', 'src_id',
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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2139332&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL472877&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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ao_endpoint',
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idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
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ion', 'src_id',
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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182514&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2106238&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2252747&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2229622&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1093743&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL454759&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2269080&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL448058&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL450072&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
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       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4074177&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL507518&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL443949&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL451328&limit=5&standard_type=IC50

These are the available columns for the Activities API:
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Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228452&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183866&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2105350&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184538&limit=5&standard_type=IC50
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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL25424&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1364260&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL324794&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
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ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228463&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL186141&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
```

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ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL121548&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228465&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1172198&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2270062&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL53493&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL250421&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3355912&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183445&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187552&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3184730&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1728549&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1594090&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3187344&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL561014&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186526&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182724&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
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https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186514&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4062155&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL206566&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2106289&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2251454&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL4076957&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL256368&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2269088&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228466&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2228467&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3188009&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL21404&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL116268&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3273404&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3733030&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL390773&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185276&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3182851&limit=5&standard_type=IC50
```

```
These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL184827&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186926&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2268552&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3183603&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL401203&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1488393&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL66926&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL506678&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
```

```
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1453070&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3560760&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186423&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3185994&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2252948&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2289233&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
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This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL477900&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2252947&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL285376&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
```

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       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL425677&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL558557&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
```

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       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1420783&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL486208&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3586094&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL448502&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
```

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       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL430091&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2048654&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3186909&limit=5&standard_type=IC50
```

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These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1077088&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL263094&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL486795&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
```

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       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL496447&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL505813&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3120650&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3120652&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
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initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2115552&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL3277900&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1222273&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL506247&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
```

```
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL1797128&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL273977&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2106930&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL430341&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
```

```
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL196&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index(['activity_comment', 'activity_id', 'activity_propertie
s',
       'assay_chembl_id', 'assay_description', 'assay_type', 'b
ao_endpoint',
       'bao_format', 'bao_label', 'canonical_smiles', 'data_val
idity_comment',
       'data_validity_description', 'document_chembl_id', 'docu
ment_journal',
       'document_year', 'ligand_efficiency', 'molecule_chembl_i
d',
       'molecule_pref_name', 'parent_molecule_chembl_id', 'pche
mbl_value',
       'potential_duplicate', 'qudt_units', 'record_id', 'relat
ion', 'src_id',
       'standard_flag', 'standard_relation', 'standard_text_val
ue',
       'standard_type', 'standard_units', 'standard_upper_valu
e',
       'standard_value', 'target_chembl_id', 'target_organism',
       'target_pref_name', 'target_tax_id', 'text_value', 'toi
d', 'type',
       'units', 'uo_units', 'upper_value', 'value'],
      dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL40274&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL486293&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
This is the url string that calls the 'Activities' API with the
initial query specification:
```

```
https://www.ebi.ac.uk/chembl/api/data/activity.json?molecule_ch
embl_id__in=CHEMBL2269081&limit=5&standard_type=IC50

These are the available columns for the Activities API:
Index([], dtype='object')
```

In [14]:
```
df.to_csv('till_now_bioactivity_3.csv', index=False)
```

Scraped Bioactivity Data and saved it as 'till_now_bioactivity_3.csv'.

In [1]:
```
import pandas as pd
df=pd.read_csv('till_now_bioactivity_3.csv')
```

In [2]:
```
df.info()
```

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 13055 entries, 0 to 13054
Data columns (total 11 columns):
 #   Column              Non-Null Count  Dtype  
---  ------              --------------  -----  
 0   target_chembl_id    13055 non-null  object 
 1   target_organism     9733 non-null   object 
 2   target_pref_name    13055 non-null  object 
 3   molecule_chembl_id  13055 non-null  object 
 4   molecule_pref_name  12845 non-null  object 
 5   pchembl_value       4591 non-null   float64
 6   standard_type       13055 non-null  object 
 7   standard_relation   7283 non-null   object 
 8   standard_value      7364 non-null   float64
 9   standard_units      7329 non-null   object 
 10  assay_chembl_id     13055 non-null  object 
dtypes: float64(2), object(9)
memory usage: 1.1+ MB
```

Scraped Bioactivity Data and saved it as 'till_now_bioactivity_3.csv'. The standard bioactivity type and units were kept same to maintain uniformity.

In [3]:
```
df.standard_type.unique()
df_ic50=df.loc[(df['standard_type']=='IC50')&(df['standard_units']=
='nM')]
```

```
df_ic50
#standard value is the potency of the drug , lower the number bette
r the potency
df_ic50.head()
```

Out[3]:

	target_chembl_id	target_organism	target_pref_name	molecule_chembl_id
0	CHEMBL364	Plasmodium falciparum	Plasmodium falciparum	CHEMBL326602
1	CHEMBL612545	NaN	Unchecked	CHEMBL37537
2	CHEMBL3879801	NaN	NON-PROTEIN TARGET	CHEMBL37537
3	CHEMBL2362975	NaN	No relevant target	CHEMBL37537
4	CHEMBL612545	NaN	Unchecked	CHEMBL37537

In [5]:

```
df_pic=df_ic50.drop(['standard_relation','pchembl_value','assay_che
mbl_id'], axis=1)
df_pic.head()
```

Out[5]:

	target_chembl_id	target_organism	target_pref_name	molecule_chembl_id
0	CHEMBL364	Plasmodium falciparum	Plasmodium falciparum	CHEMBL326602
1	CHEMBL612545	NaN	Unchecked	CHEMBL37537
2	CHEMBL3879801	NaN	NON-PROTEIN TARGET	CHEMBL37537
3	CHEMBL2362975	NaN	No relevant target	CHEMBL37537
4	CHEMBL612545	NaN	Unchecked	CHEMBL37537

In [8]:

```
#df2.info()
df.info()
df_ic50.info()
```

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 13055 entries, 0 to 13054
Data columns (total 11 columns):
 #   Column              Non-Null Count  Dtype
---  ------              --------------  -----
 0   target_chembl_id    13055 non-null  object
```

```
 1   target_organism     9733 non-null   object
 2   target_pref_name    13055 non-null  object
 3   molecule_chembl_id  13055 non-null  object
 4   molecule_pref_name  12845 non-null  object
 5   pchembl_value       4591 non-null   float64
 6   standard_type       13055 non-null  object
 7   standard_relation   7283 non-null   object
 8   standard_value      7364 non-null   float64
 9   standard_units      7329 non-null   object
 10  assay_chembl_id     13055 non-null  object
dtypes: float64(2), object(9)
memory usage: 1.1+ MB
<class 'pandas.core.frame.DataFrame'>
Int64Index: 6553 entries, 0 to 13054
Data columns (total 11 columns):
 #   Column              Non-Null Count  Dtype
---  ------              --------------  -----
 0   target_chembl_id    6553 non-null   object
 1   target_organism     4735 non-null   object
 2   target_pref_name    6553 non-null   object
 3   molecule_chembl_id  6553 non-null   object
 4   molecule_pref_name  6379 non-null   object
 5   pchembl_value       4591 non-null   float64
 6   standard_type       6553 non-null   object
 7   standard_relation   6472 non-null   object
 8   standard_value      6553 non-null   float64
 9   standard_units      6553 non-null   object
 10  assay_chembl_id     6553 non-null   object
dtypes: float64(2), object(9)
memory usage: 614.3+ KB
```

In [9]:

```
df4=df_ic50[df_ic50.standard_value.notna()]
df4.info()
df4.to_csv('bioactivity_data_flavormolecules_actual_ouput22.csv', i
ndex=False)
```

```
<class 'pandas.core.frame.DataFrame'>
Int64Index: 6553 entries, 0 to 13054
Data columns (total 11 columns):
 #   Column              Non-Null Count  Dtype
---  ------              --------------  -----
 0   target_chembl_id    6553 non-null   object
 1   target_organism     4735 non-null   object
 2   target_pref_name    6553 non-null   object
 3   molecule_chembl_id  6553 non-null   object
 4   molecule_pref_name  6379 non-null   object
 5   pchembl_value       4591 non-null   float64
```

```
 6   standard_type       6553 non-null   object
 7   standard_relation   6472 non-null   object
 8   standard_value      6553 non-null   float64
 9   standard_units      6553 non-null   object
 10  assay_chembl_id     6553 non-null   object
dtypes: float64(2), object(9)
memory usage: 614.3+ KB
```

In [13]:

```
selection = ['molecule_chembl_id','molecule_pref_name','standard_va
lue','target_chembl_id','target_pref_name','target_organism']
df5 = df4[selection]
rslt_df = df5.sort_values(by = 'standard_value', ascending = True)
rslt_df.drop_duplicates(subset='molecule_chembl_id', keep='first',
inplace=True)
rslt_df.reset_index(drop=True, inplace=True)
rslt_df.info()
rslt_df.head()
```

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 365 entries, 0 to 364
Data columns (total 6 columns):
 #   Column              Non-Null Count  Dtype
---  ------              --------------  -----
 0   molecule_chembl_id  365 non-null    object
 1   molecule_pref_name  306 non-null    object
 2   standard_value      365 non-null    float64
 3   target_chembl_id    365 non-null    object
 4   target_pref_name    365 non-null    object
 5   target_organism     293 non-null    object
dtypes: float64(1), object(5)
memory usage: 17.2+ KB
```

Out[13]:

	molecule_chembl_id	molecule_pref_name	standard_value	target_chembl_id
0	CHEMBL165	RESVERATROL	0.0225	CHEMBL399
1	CHEMBL19224	PAPAVERINE	0.0590	CHEMBL613633
2	CHEMBL50	QUERCETIN	0.1100	CHEMBL2362975
3	CHEMBL107	COLCHICINE	0.3600	CHEMBL3879801
4	CHEMBL441687	GLYCYRRHIZIN	0.4000	CHEMBL3746

In [15]:

```
rslt_df.to_csv("bioactivity_data_flavormolecules_actual_datapreproc
essed_output22.csv", index=False)
```