Tutorial of ligand similarity search or comparsion by MolAICal

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1. Introduction

Sometimes, the similarity search of ligands can help scientists to find potential compounds fastly. In this tutorial, the fingerprint similarity and 3D structural similarity searches are introduced for molecular comparison or simple virtual screening based on appointed ligand structure.

2. Materials

2.1. Software requirement

1) MolAICal: https://molaical.github.io

2.2. Example files

1) All the necessary tutorial files are downloaded from:
https://github.com/MolAICal/tutorials/tree/master/011-similaritySearch

3. Procedure

3.1. Fingerprint similarity search

Go to folder "011-similaritySearch/fingerprint"

```
#> cd 011-similaritySearch/fingerprint
```

1. Using the command "ls > listfiles.dat" in Linux console, or "dir /b > listfiles.dat" in DOS console of Window. Open generated file "listfiles.dat" and delete no useful characters. Make sure the file "listfiles.dat" only contains the ligand names.

2. Run command as below:

```
#> molaical.exe -tool finger -i listfiles.dat -o result.dat
```

It will generate the file "result.dat" which contains the values of structural similarities that correspond to names in "listfiles.dat". **Note:** the first line is the compared ligand in file "result.dat" except for the comment line which starts with "#".

3.2. 3D structural similarity search or comparsion

Go to folder "011-similaritySearch/3Dsimilarity"

```
#> cd 011-similaritySearch/3Dsimilarity
```

1. Using the command "ls > listfiles.dat" in Linux console, or "dir /b > listfiles.dat" in DOS console of Window. Open generated file "listfiles.dat" and delete no useful characters. Make sure the file "listfiles.dat" only contains the ligand names.

2. Run command as below:

```
#> molaical.exe -tool 3Dcompare -i listfiles.dat -s result.dat -f mol2list
```

It will generate the file "result.dat" which contains the values of structural similarities that

correspond to names in “listfiles.dat” except the first line. **Note:** the first line is not the compared ligand in file “result.dat”. The first line of “result.dat” corresponds to the second line of “listfiles.dat”, the second line of “result.dat” corresponds to the third line of “listfiles.dat”, and so on.

In addition, MolAICal supports the 3D structural similarity comparsion between two ligands:

```
#> molaical.exe -tool 3Dcompare -i ligand.mol2 -s 1_10023_out.mol2
```