# DeepAffinity Usage Manual

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### 1 DeepAffinity download link and its prerequisites

Source code of DeepAffinity is available at https://github.com/Shen-Lab/DeepAffinity.

DeepAffinity has been developed and tested in:

- Tensorflow-gpu v1.1
- Python 3.6
- TFlearn v0.3
- Scikit-learn v0.19
- Anaconda 3/5.0.0.1

To install Anaconda 3: Please refer to https://www.anaconda.com/distribution/ to install Anaconda 3 according to your operating system.

We have already provided our environment list as environment.yml at https://github.com/Shen-Lab/DeepAffinity/environment.yml. After installing Anaconda, you can replace "envname" with any environment name you like and create your own environment by:

```
$ conda env create -n envname -f environment.yml
```

Then please activate your environment by:

\$ source activate envname

## 2 Input Data Preparation

• Protein: SPS format.

```
e.g "CETM, ANGS, CETL, BETM, AEKM, CETL, AEKM, CEKL, AEDM, CEDL.."
```

We have provided the SPS dictionary we have generated for now. We will keep updating it and you are more than welcome to send us the new SPS sequence you generate. The dictionary can be found at:

https://github.com/Shen-Lab/DeepAffinity/data/dataset/protein\_grouped\_finalPresentation

For each two lines in this dictionary, the first line is original protein sequence from Uniprot and the second line is corresponding SPS format.

To generate SPS format for your own protein, please refer to DeepAffinity/data/script/split\_data\_script/README

• Compound: Canonical SMILE from Pubchem.

```
e.g CCC(C(=0)C)(C(=0)0)0
```

Please make sure your SMILE format is the same as Pubchem Canonical SMILE. To generate canonical SMILE based on CID, please refer to

DeepAffinity/data/script/split\_data\_script/README

• Label:

e.g 5.2

Here we use  $pIC50 = -log_{10}(IC50)$ , in which the unit of IC50 is in Mol. Please notice that you can calculate other labels, like EC50, Ki or Kd, in the same way.

# 3 Run the DeepAffinity and Output Analysis

We will introduce how to train our model and load checkpoints to test your data directly. Then we will give an example of result.

#### 3.1 Run the DeepAffinity by script

We have provided automatic script for users to run our code in one command. The options of "DeepAffinity\_inference.sh" are introduced below:

- -h: display this help and exit
- -o|--output OUTPUT\_FILE\_NAME: specify output filename
- -p|--prot PROTEIN\_SPS\_FILE: replace test\_sps with user's protein sps format file
- -c|--comp COMPOUND\_SMILE\_FILE: replace test\_smile with user's compound SIMLE format file
- -l|--label LABEL\_FILE: replace test\_ic50 with user's label file
- -t|--labeltype LABEL\_TYPE: specify the label type you are using, it can be either IC50(default), EC50, Ki or Kd

#### 3.1.1 Run the sample code

To simply run our model with IC50 dataset and test case, you can directly run

\$ ./DeepAffinity\_inference.sh

The default output will be at "./Joint\_models/joint\_attention/joint\_warm\_start/output". You can also specify the output name by

\$ ./DeepAffinity\_inference.sh -o OUTPUT\_FILE\_NAME

#### 3.1.2 Test your own data from scratch

Also, our script will allow you to test your own data with our training set by:

\$ ./DeepAffinity\_inference.sh -o OUTPUT\_FILE\_NAME -p PROTEIN\_SPS\_FILE
-c COMPOUND\_SMILE\_FILE -l LABEL\_FILE

If your label is not IC50, you can also choose your label type by adding -t LABEL\_TYPE:

\$ ./DeepAffinity\_inference.sh -o OUTPUT\_FILE\_NAME -p PROTEIN\_SPS\_FILE
-c COMPOUND\_SMILE\_FILE -1 LABEL\_FILE -t LABEL\_TYPE

#### 3.1.3 Test your own data with our checkpoints

If you would like to use our checkpoints, we have provided two versions as follows:

- Download the checkpoints trained based on training set of IC50 from https://drive.google.com/drive/folders/1 Pwn8uTyHNig4G2JDy0TErzH9hVacSadt?usp=sharing
- Download the checkpoints trained based on the whole dataset of IC50 from

```
https://drive.google.com/drive/folders/1
XAnXHSRnr08DGA1drW3YnmaBaCihdiP5?usp=sharing
```

Please note that both checkpoints we provided are for the IC50 label. To use our checkpoint to test your data, you can run it by:

```
$ ./DeepAffinity_inference.sh -o OUTPUT_FILE_NAME -p PROTEIN_SPS_FILE
    -c COMPOUND_SMILE_FILE -1 LABEL_FILE -m CHECKPOINT_FILE
```

#### 3.2 Output

You should be able to see your result at the end below "error on test" line. Here the "error" means the root mean squared error(RMSE). One example is stated as below. 48928 should be the number of cases in test set and 0.6411142 is the RMSE. There are other results from OLS regression for your reference:

```
error on test
48928
[0.6411142]
```

OLS Regression Results

Dep. Variable:	у	R-squared:	0.698
Model:	OLS	Adj. R-squared:	0.698
Method:	Least Squares	F-statistic:	1.129e+05
Date:	Thu, 23 Jan 2020	Prob (F-statistic):	0.00
Time:	16:12:46	Log-Likelihood:	-51554.

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No. Observations:	48928	AIC:	1.031e+05
Df Residuals:	48926	BIC:	1.031e+05

Df Model: 1

Covariance Type:		nonrobust				
	coef	std err	t	P> t	[0.025	0.975]
const	1.8087	0.014	126.433	0.000	1.781	1.837
x1	0.7264	0.002	336.061	0.000	0.722	0.731
Omnibus: Prob(Omnibus)	· · · · · · · · · · · · · · · · · · ·	1354. 0.		bin-Watson: que-Bera (JB):		1.750 3473.420
Skew: Kurtosis:				b(JB): d. No.		0.00 30.8