Documentation of SVMcon 1.0

SVMcon is a software to predict protein residue-residue contacts in 8 Angstrom distance threshold. It is one of the most accurate contact predictors in the 7th and 8th editions of Critical Assessment of Techniques for Protein Structure Prediction (CASP7&8). It is the only free contact map prediction software (including source code) available so far.

Reference

J. Cheng and P. Baldi. Improved Residue Contact Prediction Using Support Vector Machines and a Large Feature Set. BMC Bioinformatics. 8113, 2007.

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Please do not re-distribute the software without the permission of the author.

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Third-party software used by SVMcon

SVMcon softwre uses svm_classify program in SVM-light (svmlight.joachims.org) to classify data points.

Reference to SVM-light

T. Joachims, 11 in Making large-Scale SVM Learning Practical. Advances in Kernel Methods - Support Vector Learning, B. Schölkopf and C. Burges and A. Smola (ed.), MIT Press, 1999.

I thank the author of SVM-light (Dr. Thorsten Joachims) for the permit to include svm_classify in the SVMcon package.

Here is the license quoted from the SVM-light website:

SVM-light is free only for non-commercial use. It must not be distributed without prior permission of the author. The author is not responsible for implications from the use of this software.

Installation

1. install secondary structure and solvent accessibility predictor PSpro (PSpro1.1 or PSpro2.0)

(download it from http://sysbio.rnet.missouri.edu/multicom_toolbox/. easy to install)

- 2. unzip svmcon.tar.gz
- 3. do configuration

cd svmcon

open configure.pl

set \$install_dir to svmcon directory

set \$pspro_dir to PSpro directory (full path)

save configure.pl

run .configure.pl to configure.

installation is done.

Test Installation

cd test

..binpredict_map.sh T0288.fasta test.map

The output test.map should be the same as T0288.map.final

Run the program

predict_map.sh fasta_file output_file

fasta_file protein sequence in the simple FASTA foramt

(line1 name; line 2 a plain protein sequence).

Output file includes the predicted residue contacts in CASP format.

Notice

Since there are lot of support vectors in the SVM model, it takes from tens of minutes to several hours to predict contacts for a protein, depending on sequence length.