User manual for PANDA: Preferential Attachment based common Neighbor Distribution derived functional Associations

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2014-10-30

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1 Introduction

This file gives a brief introduction on the functions used in the R package of PANDA (preferential-attachment based common neighbor distribution derived associations). PANDA is designed to perform the following tasks in protein-protein interaction (PPI) networks: (1) identify significantly functionally associated protein pairs, (2) predict GO terms and KEGG pathways for proteins, (3) make a cluster of proteins based on the significant protein pairs, (4) identify subclusters whose members are enriched in KEGG pathways. For other types of biological networks, (1) and (3) can still be performed. For more details on PANDA, please refer to "PAND: a distribution to identify functional linkage from networks with preferential attachment property", or consult Hua Li (kaixinsjtu@hotmail.com).

2 A Quick Example

The first step is to load the package from the library.

> library(PANDA)

Then we load the example data shipped with this package.

- > data(dfPPI)
- > data(GENE2GOtopLite)

- > data(GENE2KEGG)
- > data(KEGGID2NAME)

The "dfPPI" is a PPI network consisting of 2360 proteins and 5355 interactions that is used as an example to demonstrate the capability of PANDA. The "GENE2GOtopLite" and "GENE2KEGG" are examples of GO and KEGG annotations of proteins. "KEGGID2NAME" maps KEGG pathway ID to KEGG pathway names.

We first apply PAND to the PPI network to derive functional links between protein by using the function "SignificantPairs" (protein pairs will be ranked by p-values if "pvalue=TRUE" is specified, otherwise by probabilities):

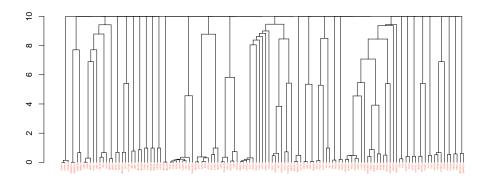
> OrderAll=SignificantPairs(PPIdb=dfPPI)

> head(OrderAll)

	Sym_A	Sym_B	${\tt Probability}$	${\tt CommonNeighbor}$
10234	CPNE1	CPNE4	-69.52455	18
8343	SMARCA4	SMARCA2	-54.93328	26
26211	LDB1	LDB2	-47.30440	16
1	SMAD2	SMAD3	-39.21016	20
4255	PER1	PER2	-39.00400	10
21882	JARID2	MTF2	-38.85149	10

Based on the p-values (or probabilities) of the significant protein pairs obtained above, we can perform agglomerative hierarchical clustering (using the unweighted group average) for proteins of all significant pairs. This function returns an object in the class "dendrogram". If "Plot=TRUE" is specified, it will also plot the dendrogram.

> dendMap=ProteinCluster(Pfile=OrderAll, Plot=TRUE, TextScaler=50)

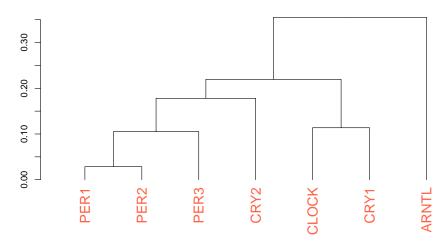


The significant protein pairs generated by the function "SignificantPairs" constituted a new network, with which we can make further functional predictions. We use the functions "GOpredict" and "KEGGpredict" to perform functional enrichment analysis (p-values were calculated with Fisher's exact test) among a protein's significant partners to predict GO terms and KEGG pathways for the protein:

```
> GP=GOpredict(Pfile=OrderAll, PPIdb=dfPPI, Gene2Annotation=GENE2GOtopLite, p_value=0.001)
> head(GP)
   Symbol
                GOTD
  TGFBR1 G0:0045669
2
     CBX2 GO:0071535
3
    MTA1 GO:0016581
4
    PBX1 G0:0007387
    THBS2 GO:0048603
6 GATAD2A GD:0072092
                                              GOterm Ratio
1 positive regulation of osteoblast differentiation
                                                       2/2
2
               RING-like zinc finger domain binding
                                                       2/2
3
                                        NuRD complex
                                                       4/5
             anterior compartment pattern formation
4
                                                       1/1
5
                   fibroblast growth factor binding
                                                       1/1
                              ureteric bud invasion
                                                       1/2
        Pvalue
1 5.496440e-05
2 3.592444e-07
3 2.764488e-09
4 4.237288e-04
5 4.237288e-04
6 8.474576e-04
> KP=KEGGpredict(Pfile=OrderAll, PPIdb=dfPPI, Gene2Annotation=GENE2KEGG,
    p_value=0.001, IDtoNAME=KEGGID2NAME)
> head(KP)
             KEGGID
    Symbol
                                      PathName Ratio
                                                           Pvalue
     SAP18 hsa05217
                         Basal cell carcinoma
                                                 2/2 5.334780e-04
2 TIMELESS hsa04710 Circadian rhythm - mammal
                                                 4/4 5.545925e-10
```

We use the following function to identify subclusters (from the cluster generated by "ProteinCluster") whose members are significantly enriched in any KEGG pathway (if KGremove=TURE, "hsa05200" and "hsa01100" will be excluded from this analysis as they are too broad):

hsa04710



3 File Location and Session Info

> sessionInfo()

R version 3.1.2 (2014-10-31)

Platform: x86_64-apple-darwin13.4.0 (64-bit)

locale:

[1] C/en_US.UTF-8/en_US.UTF-8/C/en_US.UTF-8/en_US.UTF-8

attached base packages:

- [1] parallel stats graphics grDevices utils datasets
- [7] methods base

other attached packages:

- [1] PANDA_0.9.9 AnnotationDbi_1.26.1 GenomeInfoDb_1.0.2
- [4] Biobase_2.24.0 BiocGenerics_0.10.0 RSQLite_1.0.0
- [7] DBI_0.3.1

loaded via a namespace (and not attached):

- [1] GO.db_2.14.0 IRanges_1.22.10 cluster_1.15.3 stats4_3.1.2
- [5] tools_3.1.2