

# An Introduction to *CCPROMISE*

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## 1 Introduction

CCPROMISE, Canonical correlation with PROjection onto the Most Interesting Statistical Evidence, is a general procedure to integrate two forms of genomic features that exhibit a specific biologically interesting pattern of association with multiple phenotypic endpoint variables. In biology, one type of genomic feature tends to regulate the other types. For example, DNA methylation regulates gene expression. Biological knowledge of the endpoint variables is used to define a vector that represents the biologically most interesting values for a set of association statistics. The CCPROMISE performs one hypothesis test for each gene, and is flexible to accommodate two type of genomic features with various types of endpoints.

In this document, we describe how to perform CCPROMISE procedure using hypothetical example data sets provided with the package.

## 2 Requirements

The CCPROMISE package extends our former PROMISE package to integrate two forms of molecular data with multiple biologically related endpoints in gene level or probe set level. The understanding of *ExpressionSet* is a prerequisite to perform the CCPROMISE procedure. Due to the internal handling of multiple endpoints, the consistency of *ExpressionSet* is assumed. The detailed requirements are illustrated below.

Load the CCPROMISE package and the example data sets: *exmplESet*, *exmplMSet*, *exmplGeneSet*, and *exmplPat* into R.

```
> library(CCPROMISE)
> data(exmplESet)
> data(exmplMSet)
> data(exmplGeneSet)
> data(exmplPat)
```

The *ExpressionSet* should contain at least two components: *exprs* (array data) and *phenoData* (endpoint data). The subject id and order of *ESet* and

*MSet* should be same. *exprs* is a data frame with column names representing the array identifiers (IDs) and row names representing the probe (genomic feature) IDs. *phenoData* is an *AnnotatedDataFrame* with column names representing the endpoint variables and row names representing array. The array IDs of *phenoData* and *exprs* should be matched.

The association pattern definition is critical. The prior biological knowledge is required to define the vector that represents the biologically most interesting values for statistics. In this hypothetical example, we are interested in identifying genomic features that are negatively associated with drug level to kill 50% cells, negatively associated with disease, and negatively associated with rate of events. The three endpoints are represented in three rows as shown below:

```
> exmplPat

  stat.coef      stat.func      endpt.vars
1   -0.33 spearman.rstat      LC50
2   -0.33 spearman.rstat      MRD22
3   -0.33      jung.rstat EFSTIME,EFSCENSOR
```

### 3 CCPROMISE Analysis

As mentioned in section 2, the *ExpressionSet* of two forms of genomic data and pattern definition are required by CCPROMISE procedure. The code below performs a CCPROMISE analysis at gene level with fast permutation based on negative binomial.

```
> test1 <- CCPROMISE(geneSet=exmplGeneSet,
+                   ESet=exmplESet,
+                   MSet=exmplMSet,
+                   promise.pattern=exmplPat,
+                   strat.var=NULL,
+                   prlbl=c('LC50', 'MRD22', 'EFS', 'PR3'),
+                   EMLbl=c("Expr", "Methyl"),
+                   nbperm=TRUE,
+                   max.ntail=10,
+                   nperms=100,
+                   seed=13)
```

Gene level result:

```
> head(test1$PRres)

  Gene Expr_LC50.Stat Expr_MRD22.Stat Expr_EFS.Stat
1  DDR1  0.02537225    0.06515531    0.096304751
2  RFC2 -0.19860902    0.11057641    0.194535613
3  PAX8 -0.17035650   -0.05050530   -0.014700892
4  GUCA1A -0.11857641    0.03538941    0.002563084
```

```

5  UBE1L      0.06880280      -0.05714667      -0.041141660
6  THRA      -0.05640794      -0.05223717      -0.071381381
Expr_PR3.Stat Methyl_LC50.Stat Methyl_MRD22.Stat
1  -0.062277435      -0.006796137      0.02987477
2  -0.035501001      0.049870701      0.07273351
3  0.078520899      -0.187767560      -0.06772779
4  0.026874639      0.034627937      0.15896498
5  0.009828512      NaN      NaN
6  0.060008831      0.011165082      -0.04654276
Methyl_EFS.Stat Methyl_PR3.Stat PR3.Stat
1  0.01013990      -0.011072845      -0.036675140
2  0.05561824      -0.059407483      -0.047454242
3  -0.10627160      0.120588984      0.099554942
4  0.09232005      -0.095304321      -0.034214841
5  NaN      NaN      0.009828512
6  0.01189162      0.007828685      0.033918758
Expr_LC50.Pval Expr_MRD22.Pval Expr_EFS.Pval
1  0.9000000      0.6500000      0.4000000
2  0.1666667      0.1333333      0.0000000
3  0.2100000      0.5300000      0.8700000
4  0.4615385      0.7307692      0.9230769
5  0.6363636      0.4545455      0.6363636
6  0.8333333      0.5833333      0.4166667
Expr_PR3.Pval Methyl_LC50.Pval Methyl_MRD22.Pval
1  0.4500000      0.9500000      0.9000000
2  0.5666667      0.7000000      0.3000000
3  0.2200000      0.1500000      0.4100000
4  0.6153846      0.7307692      0.0000000
5  0.9090909      NA      NA
6  0.6666667      1.0000000      0.5833333
Methyl_EFS.Pval Methyl_PR3.Pval PR3.Pval nperm
1  0.8500000      0.9000000      0.5000000      20
2  0.4666667      0.3666667      0.3333333      30
3  0.1700000      0.0900000      0.1000000      100
4  0.1923077      0.07692308      0.3846154      26
5  NA      NA      0.9090909      11
6  0.6666667      1.0000000      0.8333333      12

```

The code below performs a prbPROMISE analysis at probe pair level with fast permutation.

```

> test2 <- PrbPROMISE(geneSet=exmplGeneSet,
+                     ESet=exmplESet,
+                     MSet=exmplMSet,
+                     promise.pattern=exmplPat,
+                     strat.var=NULL,
+                     prlbl=c('LC50', 'MRD22', 'EFS', 'PR3'),

```

```

+           EMIbl=c("Expr", "Methyl"),
+           nbperm=TRUE,
+           max.ntail=10,
+           nperms=100,
+           seed=13)

```

Probe pair level correlation result at p value cut off 0.05:

```
> head(test2$CORres)
```

	Gene	Expr	Methyl
1007_s_at*cg00466425	DDR1	"1007_s_at"	"cg00466425"
1007_s_at*cg01386080	DDR1	"1007_s_at"	"cg01386080"
1007_s_at*cg01936707	DDR1	"1007_s_at"	"cg01936707"
1007_s_at*cg02313535	DDR1	"1007_s_at"	"cg02313535"
1007_s_at*cg02376496	DDR1	"1007_s_at"	"cg02376496"
1007_s_at*cg03270204	DDR1	"1007_s_at"	"cg03270204"
	Spearman.rstat	Spearman.p	
1007_s_at*cg00466425	"0.3467"	"1.46642e-05"	
1007_s_at*cg01386080	"0.2624"	"0.0011758053"	
1007_s_at*cg01936707	"0.4245"	"7.56e-08"	
1007_s_at*cg02313535	"0.1978"	"0.0150390162"	
1007_s_at*cg02376496	"0.3069"	"0.0001359269"	
1007_s_at*cg03270204	"-0.2394"	"0.0031454459"	

Probe pair level PROMISE result of probe pair at p value cut off 0.05 as above:

```
> head(test2$PRres)
```

	Gene	Expr_LC50.Stat	Expr_MRD22.Stat
1	1007_s_at*cg00466425	-0.02436901	-0.1044313
2	1007_s_at*cg01386080	-0.02436901	-0.1044313
3	1007_s_at*cg01936707	-0.02436901	-0.1044313
4	1007_s_at*cg02313535	-0.02436901	-0.1044313
5	1007_s_at*cg02376496	-0.02436901	-0.1044313
6	1007_s_at*cg03270204	-0.02436901	-0.1044313
	Expr_EFS.Stat	Expr_PR3.Stat	Methyl_LC50.Stat
1	-0.1031647	0.07732165	-0.03802601
2	-0.1031647	0.07732165	-0.12352788
3	-0.1031647	0.07732165	-0.14957974
4	-0.1031647	0.07732165	0.11877059
5	-0.1031647	0.07732165	-0.02394829
6	-0.1031647	0.07732165	0.05799370
	Methyl_MRD22.Stat	Methyl_EFS.Stat	Methyl_PR3.Stat
1	0.092249971	0.11770019	-0.05730805
2	-0.009996237	0.03141073	0.03403780
3	0.069168489	-0.02279546	0.03440224
4	0.165563303	0.05556954	-0.11330114

5	0.167249598	0.17349992	-0.10560041	
6	-0.133609780	-0.14570372	0.07377326	
	PR3.Stat	Expr_LC50.Pval	Expr_MRD22.Pval	Expr_EFS.Pval
1	0.01000680	1.0000000	0.4000000	0.3000000
2	0.05567973	0.8947368	0.1842105	0.2631579
3	0.05586195	0.9000000	0.3500000	0.3000000
4	-0.01798974	0.9166667	0.3333333	0.2500000
5	-0.01413938	1.0000000	0.4000000	0.3000000
6	0.07554746	0.8730159	0.1904762	0.2222222
	Expr_PR3.Pval	Methyl_LC50.Pval	Methyl_MRD22.Pval	
1	0.5000000	0.6000000	0.3000000	
2	0.3157895	0.4210526	0.8684211	
3	0.5000000	0.3500000	0.4500000	
4	0.5000000	0.1666667	0.0000000	
5	0.5000000	1.0000000	0.0000000	
6	0.2539683	0.7301587	0.1428571	
	Methyl_EFS.Pval	Methyl_PR3.Pval	PR3.Pval	nperm
1	0.2000000	0.5000000	1.0000000	10
2	0.76315789	0.68421053	0.2631579	38
3	0.8500000	0.6500000	0.5000000	20
4	0.33333333	0.08333333	0.8333333	12
5	0.0000000	0.1000000	1.0000000	10
6	0.06349206	0.28571429	0.1587302	63