

MetaboAnalystを用いた PCA, PLS-DA

<https://www.metaboanalyst.ca/>

データファイルの作り方

各サンプルが各行になるようにする。行と列を入れ替えることも可能

1列目にID、2列目にサンプルの分類情報、3列目以降に変数データを記入する

id	class	Q1	Q2	Q3	Q4	Q5	Q6	Q7	Q8
1	M	100	100	50	70	10	100	100	
2	F	80	100	90	100	0	90	100	
3	F	50	100	100	100	0	70	80	
4	F	100	100	50	50	0	70	100	
5	F	94	80	98	30	60	98	99	
6	F	99	100	80	70	0	10	60	
7	M	50	40	80	50	0	80	40	
8	F	55	50	69	100	0	0	15	
9	F	25	0	100	100	0	25	75	
10	F	30	75	80	60	0	90	80	
11	M	85	100	85	90	0	100	40	
12	M	23	74	87	72	0	68	82	
13	F	70	100	80	100	10	60	60	
14	F	55	90	100	80	30	50	42	
15	F	50	80	100	80	0	30	100	
16	F	90	50	70	70	0	0	40	
17	M	80	30	100	100	2	30	95	

転置可能



データファイルの作り方

カンマ区切りファイル(csv)にしたいデータだけを記載した、エクセルシートを作る。

そのシートを閲覧している状態で、
「名前を付けて保存」から、カンマ区切り
(csv)を選んで保存する。

MetaboAnalystへアクセス

click here to startをクリック



MetaboAnalyst - statistical, functional and integrative analysis of metabolomics data

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Welcome [>> click here to start <<](#)

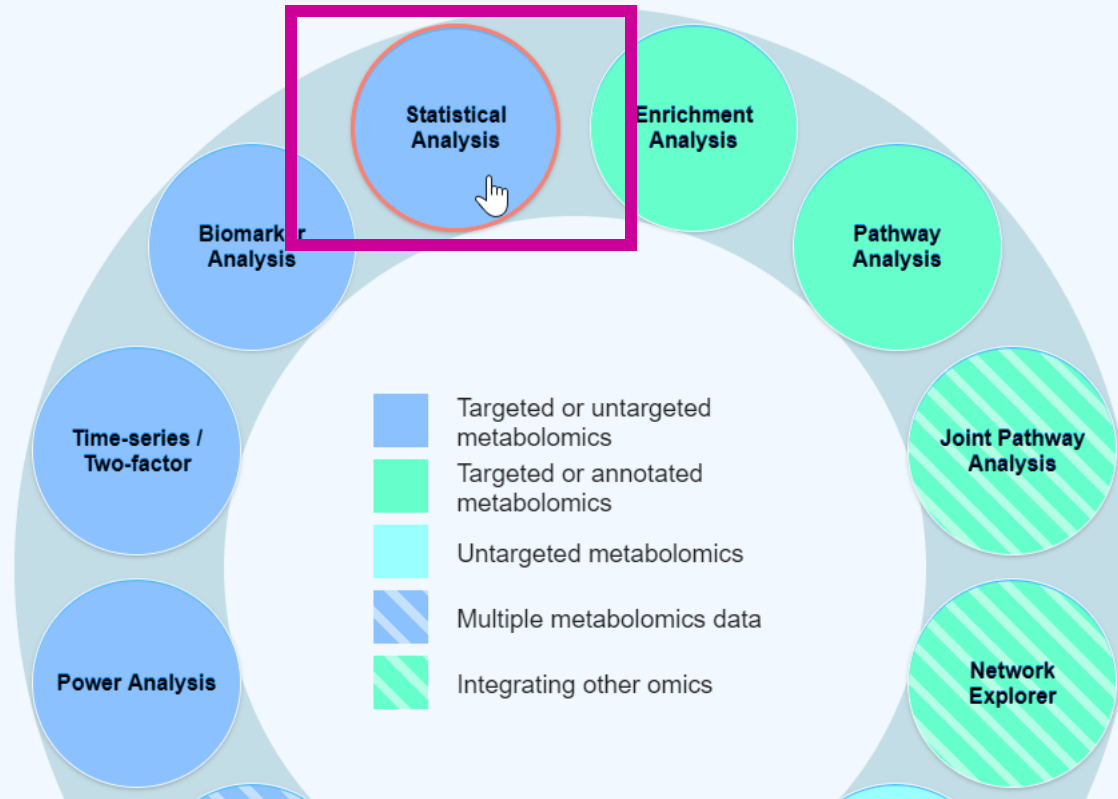
News & Updates

- Check out [OmicsNet](#) for multi-omics data integration via 3D network visual analytics;
- Check out [MicrobiomeAnalyst](#) for comprehensive analysis of microbiome data;
- Check out [NetworkAnalyst](#) for comprehensive gene expression profiling & meta-analysis;
- Users can upload MS peak intensity tables and test enriched pathways for any peak cluster of interest (MS Peaks to Pathways) (09/13/2019); **NEW**
- Enhanced enrichment network view for metabolite set enrichment analysis (06/25/2019); **NEW**
- Added mixed mode support to allow combining peaks from both positive and negative ion modes (MS peaks to pathways) (06/10/2019); **NEW**
- Upgraded the web framework (05/27/2019);
- Improved box plot graphics based on user feedback (05/06/2019);
- Enhanced the summary plot for individual metabolite set in Enrichment Analysis module (04/23/2019);
- Enhanced support for various input formats and added interactive result summary for MS Peaks to Pathways module (04/10/2019);
- Added support for synchronized 3D visual exploration for PCA/PLS-DA scores and loading plots (04/07/2019);
- Added support for the GSEA algorithm (MS Peaks to Pathways) (03/22/2019);
- The download table in compound ID conversion now contains SMILES (03/15/2019);
- Enhanced ID mapping for KEGG compounds (03/11/2019);
- Enhanced KEGG pathway generation to address the occasional failure issue during peak time (01/24/2019);
- Fixed the issue for compound view in Pathway Analysis module (01/07/2019);
- Enhanced support for parsing data input and data editing (01/02/2019);

[Read more](#)

Statistical Analysisをクリック

Click a module to proceed, or **scroll down** for more details:



準備したCSVファイルをアップロード

必要に応じて、Formatなどを選択する

Submitボタンを押す

1) Upload your data

Tab-delimited text (.txt) or comma-separated values (.csv) file:

Data Type: ☒ Concentrations ☐ Spectral bins ☐ Peak intensity table

Format: ▼

Data File: okonomi.csv

Submit

データチェックが行われる 問題なければskipで先に進む

Data Integrity Check:

1. Checking the class labels - at least three replicates are required in each class.
2. If the samples are paired, the pair labels must conform to the specified format.
3. The data (except class labels) must not contain non-numeric values.
4. The presence of missing values or features with constant values (i.e. all zeros)

Data processing information:

Checking data content ...passed

Samples are in rows and features in columns

The uploaded file is in comma separated values (.csv) format.

The uploaded data file contains 33 (samples) by 29 (compounds) data matrix.

Samples are not paired.

2 groups were detected in samples.

Only English letters, numbers, underscore, hyphen and forward slash (/) are allowed.

Other special characters or punctuations (if any) will be stripped off.

All data values are numeric.

A total of 0 (0%) missing values were detected.

By default, these values will be replaced by a small value.

Click **Skip** button if you accept the default practice

Or click **Missing value imputation** to use other methods

Missing value estimation

Skip

必要に応じて、データの標準化（normalization）、変換（transformation）、スケーリング方法を選択する

Normalizeボタンを押す

Normalization overview:

The normalization procedures are grouped into three categories. The sample normalization allows general-purpose adjustment for differences among your sample; data transformation and scaling are two different approaches to make individual features more comparable. You can use one or combine them to achieve better results.

Sample normalization

☒ None

☐ Sample-specific normalization (i.e. weight, volume) [Specify](#)

☐ Normalization by sum

☐ Normalization by median

☐ Normalization by reference sample (PQN) [Specify](#)

☐ Normalization by a pooled sample from group [Specify](#)

☐ Normalization by reference feature [Specify](#)

☐ Quantile normalization

Data transformation

☒ None

☐ Log transformation (generalized logarithm transformation or log)

☐ Cube root transformation (takes the cube root of data values)

Data scaling

☒ None

☐ Mean centering (mean-centered only)

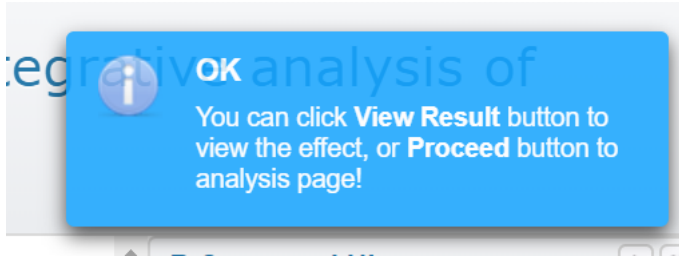
☐ Auto scaling (mean-centered and divided by the standard deviation of each variable)

☐ Pareto scaling (mean-centered and divided by the square root of the standard deviation of each variable)

☐ Range scaling (mean-centered and divided by the range of each variable)

Normalize **View Result** **Proceed**

右上にこのウィンドウが出ればOK



アクティブになったProceedボタンを押す

Data scaling

☒ **None**

☐ **Mean centering** (mean-centered only)

☐ **Auto scaling** (mean-centered and divided by the standard deviation of each variable)

☐ **Pareto scaling** (mean-centered and divided by the square root of the standard deviation of each variable)

☐ **Range scaling** (mean-centered and divided by the range of each variable)

Normalize **View Result** **Proceed**

主成分分析

Principal Component Analysis (PCA) をクリック

Select an analysis path to explore :

Univariate Analysis

[Fold Change Analysis](#) [T-tests](#) [Volcano plot](#)

[One-way Analysis of Variance \(ANOVA\)](#)

[Correlation Analysis](#) [Pattern Searching](#)

Chemometrics Analysis

[Principal Component Analysis \(PCA\)](#)

[Partial Least Squares - Discriminant Analysis \(PLS-DA\)](#)

[Sparse Partial Least Squares - Discriminant Analysis \(sPLS-DA\)](#)

[Orthogonal Partial Least Squares - Discriminant Analysis \(orthoPLS-DA\)](#)

Feature Identification

[Significance Analysis of Microarray \(and Metabolites\) \(SAM\)](#)

[Empirical Bayesian Analysis of Microarray \(and Metabolites\) \(EBAM\)](#)

結果画面が表示される

Overview

Scree Plot

2D Scores Plot

Loadings Plot

Synchronized 3D Plots

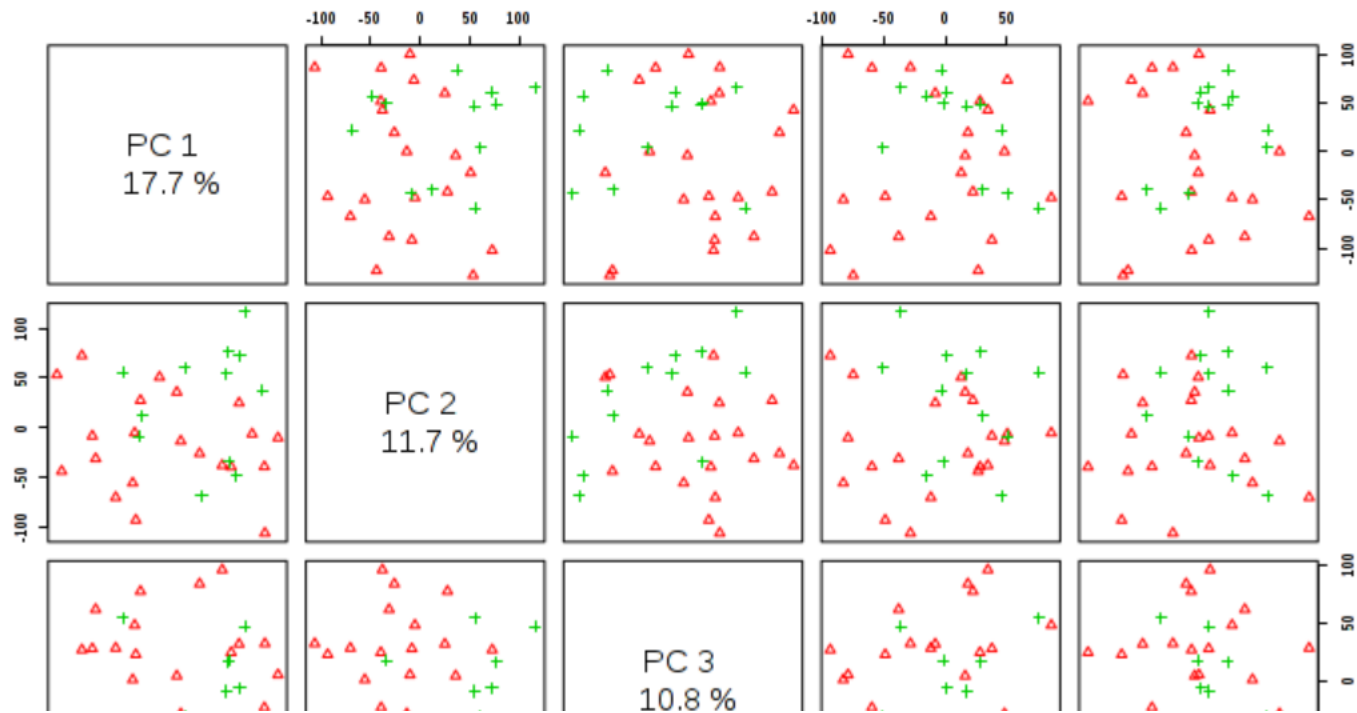
Biplot

Display pairwise score plot for top

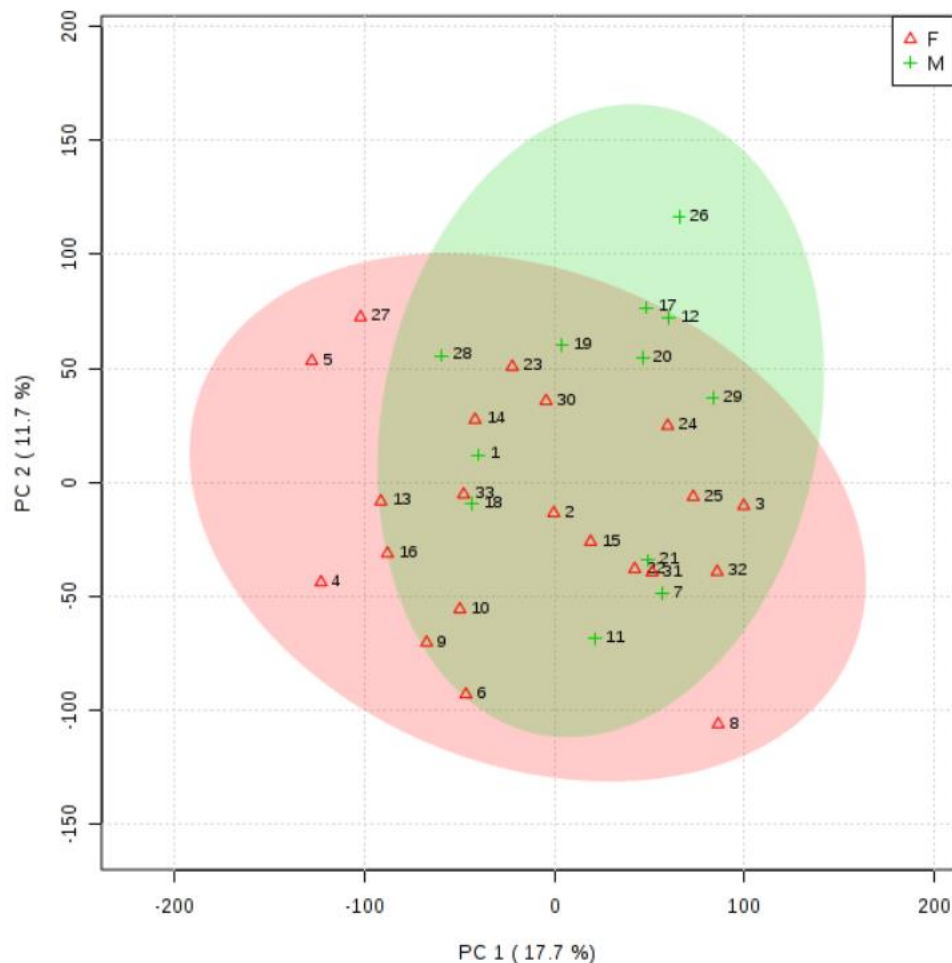
5

PCs

Update



2D Scores Plot



Specify PC on X-axis:

Specify PC on Y-axis:

Display 95% confidence regions:

☒

Display sample names:

☒

Use grey-scale colors:

☐

Flip Image

☐

X axis

☒

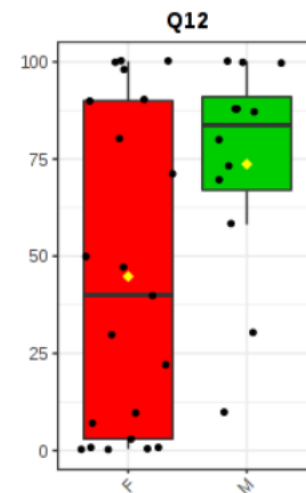
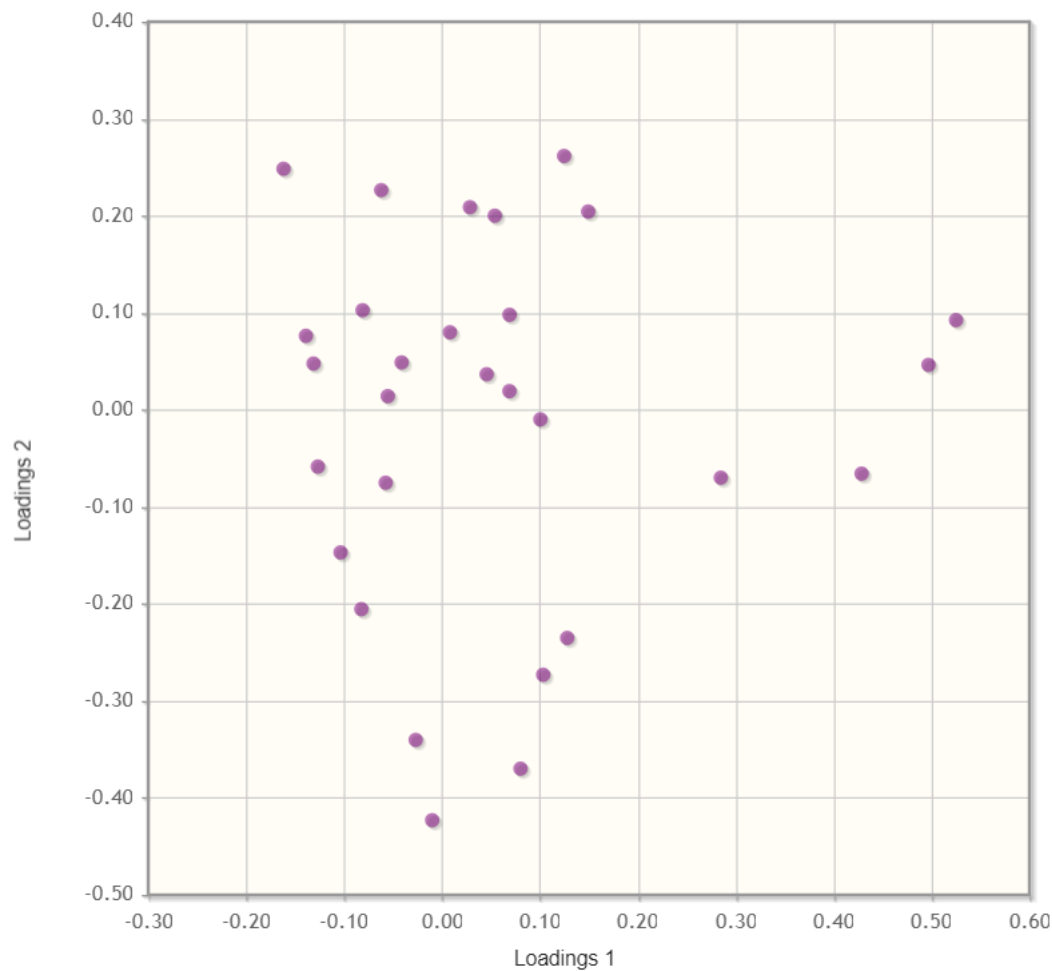
Y axis

☐

All

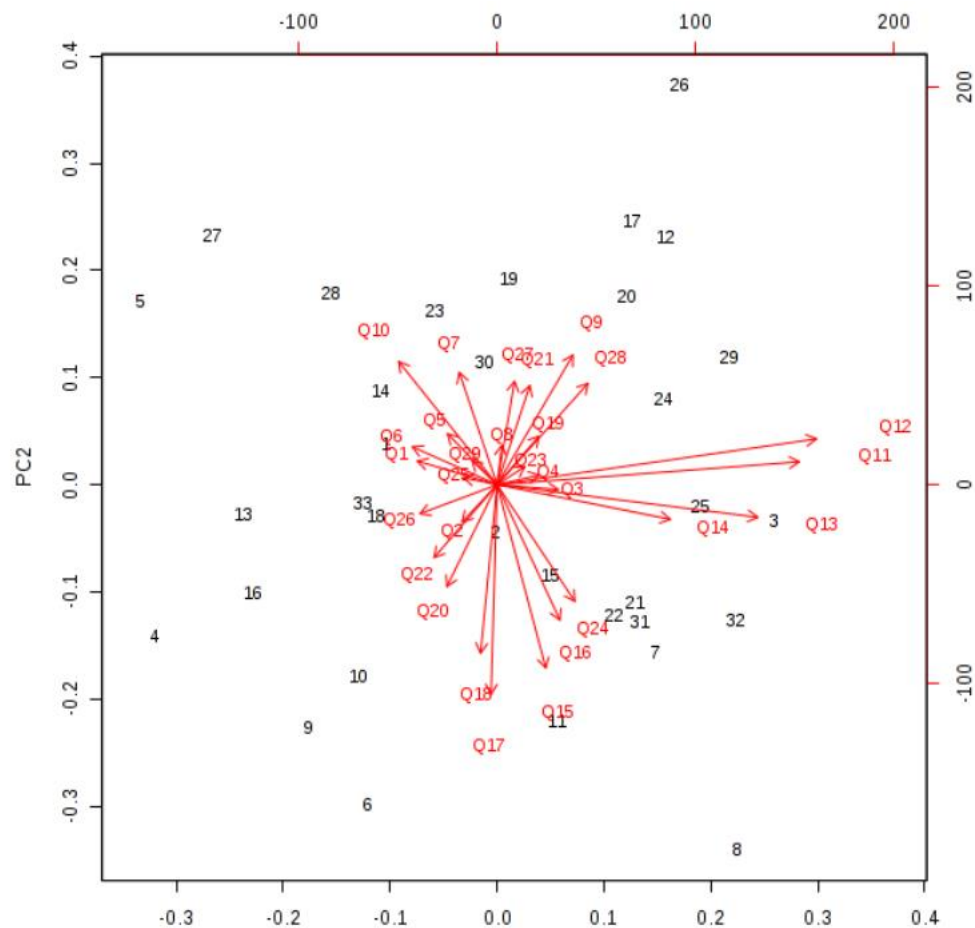
Display sample
nameにチェックが
入っていると、サン
プル名が一緒に表示
される

Loadings Plot



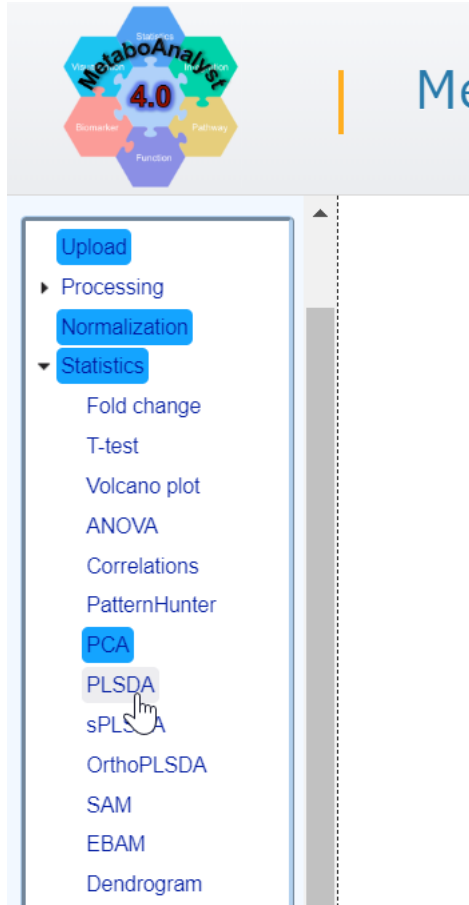
プロットをクリックすると、
その変数の特徴が表示される。

Biplot

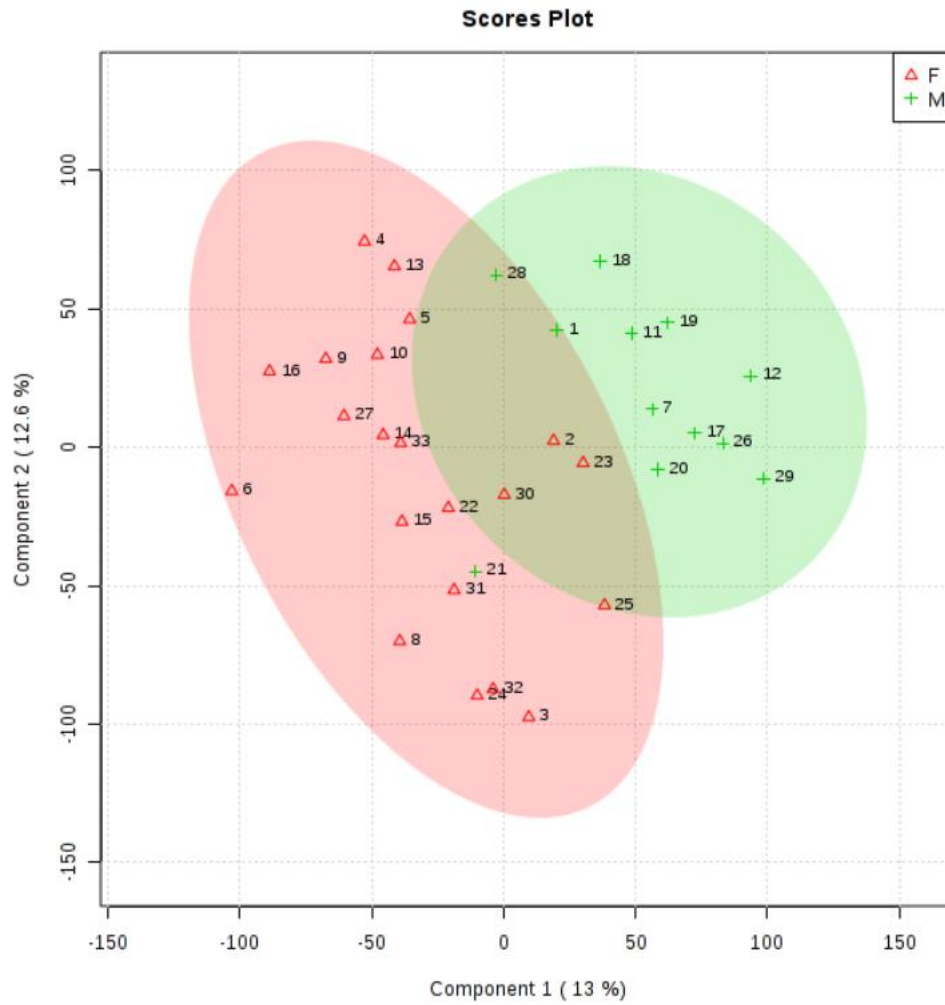


スコアプロットとローディングプロットが一緒に描かれたもの。

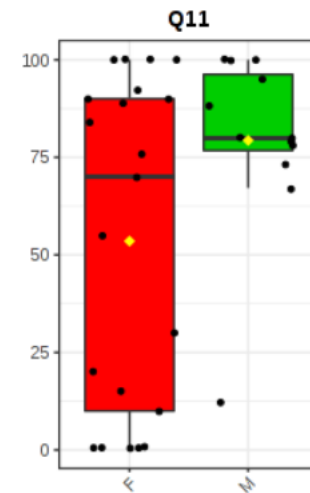
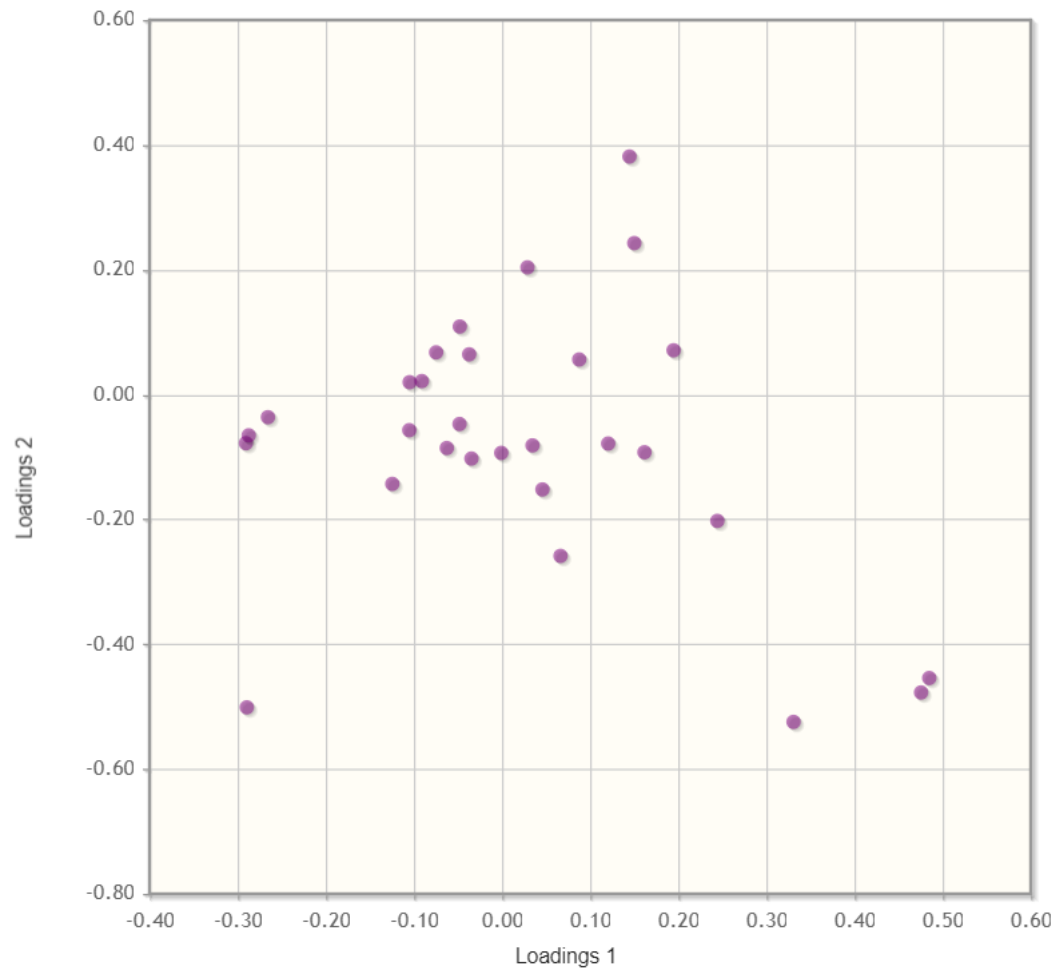
PLSDA



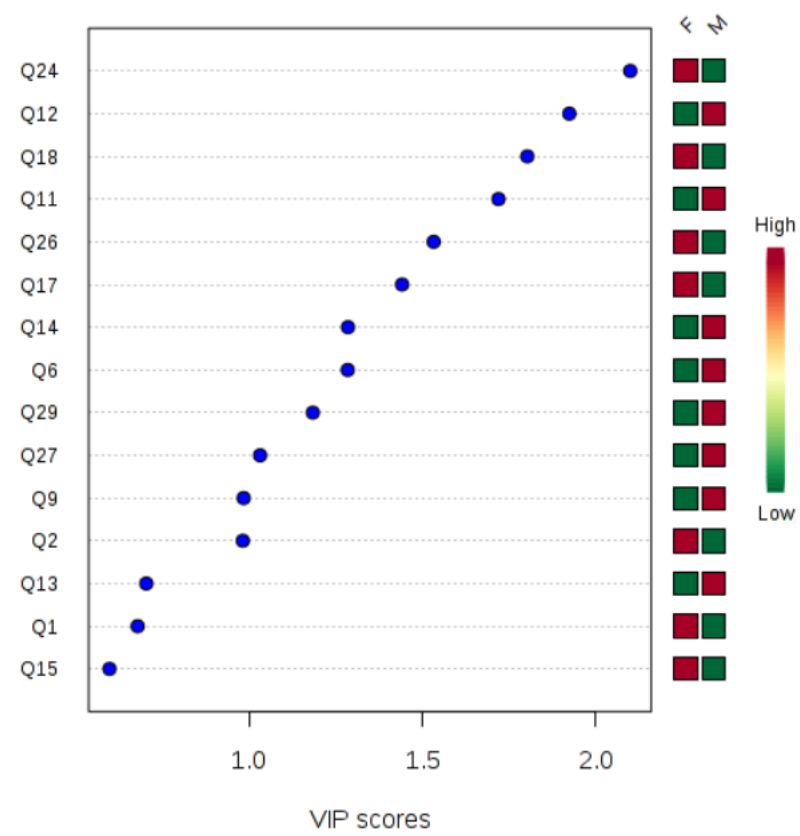
2D Scores Plot



Loadings Plot



VIP值



累積寄与率と構築したモデルの精度(Q)

