

# Package: OmicsPLS (via r-universe)

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**Type** Package

**Title** Data Integration with Two-Way Orthogonal Partial Least Squares

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**Description** Performs the O2PLS data integration method for two datasets, yielding joint and data-specific parts for each dataset. The algorithm automatically switches to a memory-efficient approach to fit O2PLS to high dimensional data. It provides a rigorous and a faster alternative cross-validation method to select the number of components, as well as functions to report proportions of explained variation and to construct plots of the results. See the software article by el Bouhaddani et al (2018) <[doi:10.1186/s12859-018-2371-3](https://doi.org/10.1186/s12859-018-2371-3)>, and Trygg and Wold (2003) <[doi:10.1002/cem.775](https://doi.org/10.1002/cem.775)>. It also performs Sparse Group (Penalized) O2PLS, see Gu et al (2020) <[doi:10.1186/s12859-021-03958-3](https://doi.org/10.1186/s12859-021-03958-3)> and cross-validation for the degree of sparsity.

**License** GPL-3

**Encoding** UTF-8

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adjR2

*Gridwise adjusted R2 for O2PLS*

---

### Description

For (a grid of) values for  $a$ ,  $n_x$  and  $n_y$ , `loocv` calculates the  $R^2$  of the joint part. Parallel computing is supported on Windows with package `parallel`.

**Usage**

```
adjR2(
  X,
  Y,
  a = 1:2,
  a2 = 1,
  b2 = 1,
  func = o2m,
  parall = F,
  cl = NULL,
  stripped = TRUE,
  p_thresh = 3000,
  q_thresh = p_thresh,
  tol = 1e-10,
  max_iterations = 100
)
```

**Arguments**

X	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
Y	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
a	Vector of integers. Contains the numbers of joint components.
a2	Vector of integers. Contains the numbers of orthogonal components in <i>X</i> .
b2	Vector of integers. Contains the numbers of orthogonal components in <i>Y</i> .
func	Function to fit the O2PLS model with. Only <code>o2m</code> and <code>o2m_stripped</code> are supported.
parall	Integer. Should a parallel cluster be set up using package <code>parallel</code> (Windows)? Best is to leave it to <code>FALSE</code> .
cl	Object of class <code>'cluster'</code> . If <code>parall</code> is <code>TRUE</code> and <code>cl</code> is not <code>NULL</code> , calculations are parallelized over workers in <code>cl</code> .
stripped	Logical. Use the stripped version of <code>o2m</code> (usually when cross-validating)?
p_thresh	Integer. If <i>X</i> has more than <code>p_thresh</code> columns, a power method optimization is used, see <a href="#">o2m2</a>
q_thresh	Integer. If <i>Y</i> has more than <code>q_thresh</code> columns, a power method optimization is used, see <a href="#">o2m2</a>
tol	Double. Threshold for which the NIPALS method is deemed converged. Must be positive.
max_iterations	Integer. Maximum number of iterations for the NIPALS method.

**Details**

The use of this function is to calculate the R2 of the joint part, while varying the number of orthogonal components. Adding more joint components will increase the R2!

A parallelized version is built in -tested on windows-, use package `parallel` and set `parall=TRUE` to activate this. There should not be already a cluster object with the name `c1`. In case of some error, don't forget to invoke `stopCluster(c1)` to end the cluster. See Task Manager (Windows) to verify that the workers are spanned/ended.

See [loocv](#) for more intuition.

### Value

Matrix with two rows:

<code>adjR2X</code>	Contains the joint R2 in X
<code>adjR2Y</code>	Contains the joint R2 in Y

---

<code>crossval_o2m</code>	<i>Cross-validate procedure for O2PLS</i>
---------------------------	---

---

### Description

Cross-validate procedure for O2PLS

### Usage

```
crossval_o2m(
  X,
  Y,
  a,
  ax,
  ay,
  nr_folds,
  nr_cores = 1,
  stripped = TRUE,
  p_thresh = 3000,
  q_thresh = p_thresh,
  tol = 1e-10,
  max_iterations = 100
)
```

### Arguments

<code>X</code>	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
<code>Y</code>	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
<code>a</code>	Vector of positive integers. Denotes the numbers of joint components to consider.
<code>ax</code>	Vector of non-negative integers. Denotes the numbers of X-specific components to consider.

ay	Vector of non-negative integers. Denotes the numbers of Y-specific components to consider.
nr_folds	Positive integer. Number of folds to consider. Note: kcv=N gives leave-one-out CV. Note that CV with less than two folds does not make sense.
nr_cores	Positive integer. Number of cores to use for CV. You might want to use <a href="#">detectCores()</a> . Defaults to 1.
stripped	Logical. Use the stripped version of o2m (usually when cross-validating)?
p_thresh	Integer. If X has more than p_thresh columns, a power method optimization is used, see <a href="#">o2m2</a>
q_thresh	Integer. If Y has more than q_thresh columns, a power method optimization is used, see <a href="#">o2m2</a>
tol	Double. Threshold for which the NIPALS method is deemed converged. Must be positive.
max_iterations	Integer. Maximum number of iterations for the NIPALS method.

### Details

This is the standard CV approach. It minimizes the sum of the prediction errors of X and Y over a three-dimensional grid of integers. Parallelization is possible on all platforms. On Windows it uses [makePSOCKcluster](#), then exports all necessary objects to the workers, and then calls [parLapply](#). On OSX and Linux the more friendly [mclapply](#) is used, which uses forking. A print method is defined, printing the minimizers and minimum in a readable way. Also the elapsed time is tracked and reported.

### Value

List of class "cvo2m" with the original and sorted Prediction errors and the number of folds used.

### Examples

```
local({
  X = scale(jitter(tcrossprod(rnorm(100),runif(10))))
  Y = scale(jitter(tcrossprod(rnorm(100),runif(10))))
  crossval_o2m(X, Y, a = 1:4, ax = 1:2, ay = 1:2,
              nr_folds = 5, nr_cores = 1)
})
```

---

crossval\_o2m\_adjR2      *Adjusted Cross-validate procedure for O2PLS*

---

### Description

Combines CV with R2 optimization

**Usage**

```
crossval_o2m_adjR2(
  X,
  Y,
  a,
  ax,
  ay,
  nr_folds,
  nr_cores = 1,
  stripped = TRUE,
  p_thresh = 3000,
  q_thresh = p_thresh,
  tol = 1e-10,
  max_iterations = 100
)
```

**Arguments**

X	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
Y	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
a	Vector of positive integers. Denotes the numbers of joint components to consider.
ax	Vector of non-negative integers. Denotes the numbers of X-specific components to consider.
ay	Vector of non-negative integers. Denotes the numbers of Y-specific components to consider.
nr_folds	Positive integer. Number of folds to consider. Note: <code>kcv=N</code> gives leave-one-out CV. Note that CV with less than two folds does not make sense.
nr_cores	Positive integer. Number of cores to use for CV. You might want to use <code>detectCores()</code> . Defaults to 1.
stripped	Logical. Use the stripped version of <code>o2m</code> (usually when cross-validating)?
p_thresh	Integer. If X has more than <code>p_thresh</code> columns, a power method optimization is used, see <a href="#">o2m2</a>
q_thresh	Integer. If Y has more than <code>q_thresh</code> columns, a power method optimization is used, see <a href="#">o2m2</a>
tol	Double. Threshold for which the NIPALS method is deemed converged. Must be positive.
max_iterations	Integer. Maximum number of iterations for the NIPALS method.

**Details**

This is an alternative way of cross-validating. It is proposed in [citation\(OmicsPLS\)](#). This approach is (much) faster than the standard `crossval_o2m` approach and works fine even with two

folders. For each element in  $n$  it looks for  $n_x$  and  $n_y$  that maximize the  $R^2$  between  $T$  and  $U$  in the O2PLS model. This approach often yields similar integer as the standard approach. We however suggest to use the standard approach to minimize the prediction error around the found integers.

### Value

data.frame with four columns: MSE,  $n$ ,  $n_x$  and  $n_y$ . Each row corresponds to an element in  $a$ .

### Examples

```
local({
  X = scale(jitter(tcrossprod(rnorm(100),runif(10))))
  Y = scale(jitter(tcrossprod(rnorm(100),runif(10))))
  crossval_o2m_adjR2(X, Y, a = 1:4, ax = 1:2, ay = 1:2,
                    nr_folds = 5, nr_cores = 1)
})
```

---

crossval_sparsity	<i>Perform cross-validation to find the optimal number of variables/groups to keep for each joint component</i>
-------------------	---

---

### Description

Perform cross-validation to find the optimal number of variables/groups to keep for each joint component

### Usage

```
crossval_sparsity(
  X,
  Y,
  n,
  nx,
  ny,
  nr_folds,
  keepx_seq = NULL,
  keepy_seq = NULL,
  groupx = NULL,
  groupy = NULL,
  tol = 1e-10,
  max_iterations = 100
)
```

### Arguments

$X$	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
-----	---

Y	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
n	Integer. Number of joint PLS components. Must be positive.
nx	Integer. Number of orthogonal components in <i>X</i> . Negative values are interpreted as 0
ny	Integer. Number of orthogonal components in <i>Y</i> . Negative values are interpreted as 0
nr_folds	Integer. Number of folds of CV
keepx_seq	Numeric vector. A vector indicating how many variables/groups to keep for CV in each of the joint component of <i>X</i> . Sparsity of each joint component will be selected sequentially.
keepy_seq	Numeric vector. A vector indicating how many variables/groups to keep for CV in each of the joint component of <i>Y</i> . Sparsity of each joint component will be selected sequentially.
groupx	Vector. Used when <code>sparse = TRUE</code> . A vector of strings indicating group names of each <i>X</i> -variable. Its length must be equal to the number of variables in <i>X</i> . The order of group names must corresponds to the order of the variables.
groupy	Vector. Used when <code>sparse = TRUE</code> . A vector of strings indicating group names of each <i>Y</i> -variable. The length must be equal to the number of variables in <i>Y</i> . The order of group names must corresponds to the order of the variables.
tol	Double. Threshold for which the NIPALS method is deemed converged. Must be positive.
max_iterations	Integer. Maximum number of iterations for the NIPALS method.

### Value

A list containing

x_1sd	A vector with length n, giving the optimal number of variables/groups to keep for each X-joint component. One standard error rule is applied
y_1sd	A vector with length n, giving the optimal number of variables/groups to keep for each Y-joint component. One standard error rule is applied
x	A vector with length n, giving the optimal number of variables/groups to keep for each X-joint component, without applying the one standard error rule
y	A vector with length n, giving the optimal number of variables/groups to keep for each Y-joint component, without applying the one standard error rule



---

impute_matrix	<i>Impute missing values in a matrix</i>
---------------	--

---

**Description**

Impute missing values in a matrix

**Usage**

```
impute_matrix(X, ...)
```

**Arguments**

X	A matrix with missing values in some entries.
...	Further arguments for <code>softimpute</code> .

**Details**

This function is based on the [softImpute](#) function in its eponymous package.

**Value**

An imputed version of matrix *X*

**Examples**

```
X <- matrix(rnorm(20*100),20)
Xmis <- X
Xmis[sample(length(Xmis),length(Xmis)/10)] <- NA
anyNA(X)
anyNA(impute_matrix(Xmis))
```

---

loadings	<i>Extract the loadings from an O2PLS fit</i>
----------	---

---

**Description**

This function extracts loading parameters from an O2PLS fit

**Usage**

```
loadings(x, ...)

## S3 method for class 'o2m'
loadings(
  x,
  loading_name = c("Xjoint", "Yjoint", "gr_Xjoint", "gr_Yjoint", "Xorth", "Yorth"),
  subset = 0,
  sorted = FALSE,
  ...
)
```

**Arguments**

x	Object of class o2m
...	For consistency
loading_name	character string. One of the following: 'Xjoint', 'Yjoint', 'gr_Xjoint', 'gr_Yjoint', 'Xorth' or 'Yorth'.
subset	subset of loading vectors to be extracted.
sorted	Logical. Should the rows of the loadings be sorted according to the absolute magnitude of the first column?

**Value**

Loading matrix

**See Also**

[scores.o2m](#)

**Examples**

```
loadings(o2m(scale(-2:2), scale(-2:2*4), 1, 0, 0))
```

**Description**

For (a grid of) values for a, nx and ny, loocv estimates the prediction error using k-fold CV.

**Usage**

```
loocv(
  X,
  Y,
  a = 1:2,
  a2 = 1,
  b2 = 1,
  fitted_model = NULL,
  func = o2m,
  app_err = F,
  kcv,
  stripped = TRUE,
  p_thresh = 3000,
  q_thresh = p_thresh,
  tol = 1e-10,
  max_iterations = 100
)
```

**Arguments**

<code>X</code>	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
<code>Y</code>	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
<code>a</code>	Vector of integers. Contains the numbers of joint components.
<code>a2</code>	Vector of integers. Contains the numbers of orthogonal components in $X$ .
<code>b2</code>	Vector of integers. Contains the numbers of orthogonal components in $Y$ .
<code>fitted_model</code>	List. Deprecated. O2PLS model fit with <code>o2m</code> . Is used to calculate the apparent error without recalculating this fit.
<code>func</code>	Function to fit the O2PLS model with. Only <code>o2m</code> and <code>o2m_stripped</code> are supported.
<code>app_err</code>	Logical. Deprecated. Should the apparent error also be computed?
<code>kcv</code>	Integer. The value of $k$ , i.e. the number of folds. Choose $N$ for LOO-CV.
<code>stripped</code>	Logical. Use the stripped version of <code>o2m</code> (usually when cross-validating)?
<code>p_thresh</code>	Integer. If $X$ has more than <code>p_thresh</code> columns, a power method optimization is used, see <code>o2m2</code>
<code>q_thresh</code>	Integer. If $Y$ has more than <code>q_thresh</code> columns, a power method optimization is used, see <code>o2m2</code>
<code>tol</code>	Double. Threshold for which the NIPALS method is deemed converged. Must be positive.
<code>max_iterations</code>	Integer. Maximum number of iterations for the NIPALS method.

**Details**

Note that this function can be easily parallelized (on Windows e.g. with the `parallel` package.).

The parameters `a`, `a2` and `b2` can be integers or vectors of integers. A for loop is used to loop over all combinations. The resulting output is a list, which is more easy to interpret if you use `array(unlist(output_of_loocv$CVerr))` as in the example below. The array will have varying `a` along the first dimension and `a2` and `b2` along the second and third respectively. Typing `example(loocv)` (hopefully) clarifies this function.

**Value**

List with two numeric vectors:

<code>CVerr</code>	Contains the k-fold CV estimated RMSEP
<code>Fiterr</code>	Contains the apparent error

---

<code>loocv_combi</code>	<i>K-fold CV based on symmetrized prediction error</i>
--------------------------	--

---

**Description**

The prediction error of both  $X \sim \hat{X}$  and  $Y \sim \hat{Y}$  are summed. This provides a symmetrized version of [loocv](#).

**Usage**

```
loocv_combi(
  X,
  Y,
  a = 1:2,
  a2 = 1,
  b2 = 1,
  fitted_model = NULL,
  func = o2m,
  app_err = F,
  kcv,
  stripped = TRUE,
  p_thresh = 3000,
  q_thresh = p_thresh,
  tol = 1e-10,
  max_iterations = 100
)
```

**Arguments**

<code>X</code>	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
<code>Y</code>	Numeric matrix. Vectors will be coerced to matrix with <code>as.matrix</code> (if this is possible)
<code>a</code>	Vector of integers. Contains the numbers of joint components.
<code>a2</code>	Vector of integers. Contains the numbers of orthogonal components in $X$ .
<code>b2</code>	Vector of integers. Contains the numbers of orthogonal components in $Y$ .
<code>fitted_model</code>	List. Deprecated. O2PLS model fit with <code>o2m</code> . Is used to calculate the apparent error without recalculating this fit.
<code>func</code>	Function to fit the O2PLS model with. Only <code>o2m</code> and <code>o2m_stripped</code> are supported.
<code>app_err</code>	Logical. Deprecated. Should the apparent error also be computed?
<code>kcv</code>	Integer. The value of $k$ , i.e. the number of folds. Choose $N$ for LOO-CV.
<code>stripped</code>	Logical. Use the stripped version of <code>o2m</code> (usually when cross-validating)?
<code>p_thresh</code>	Integer. If $X$ has more than <code>p_thresh</code> columns, a power method optimization is used, see <code>o2m2</code>
<code>q_thresh</code>	Integer. If $Y$ has more than <code>q_thresh</code> columns, a power method optimization is used, see <code>o2m2</code>
<code>tol</code>	Double. Threshold for which the NIPALS method is deemed converged. Must be positive.
<code>max_iterations</code>	Integer. Maximum number of iterations for the NIPALS method.

**Details**

Note that this function can be easily parallelized (on Windows e.g. with the `parallel` package.). If there are NAs in the `CVerr` component, this is due to an error in the fitting.

**Value**

List with two numeric vectors:

<code>CVerr</code>	Contains the $k$ -fold CV estimated RMSEP
<code>Fiterr</code>	Contains the apparent error

---

mse	<i>Calculate mean squared difference</i>
-----	--

---

**Description**

Calculate mean squared difference

**Usage**

```
mse(x, y = 0, na.rm = FALSE)
```

**Arguments**

x	Numeric vector or matrix.
y	Numeric vector or matrix. Defaults to 0.
na.rm	Remove NA's?

**Details**

Is equal to  $\text{ssq}(x-y)/\text{length}(c(x))$ . If x and y are of unequal length, the invoked minus-operator will try to make the best out of it by recycling elements of the shorter object (usually you don't want that). In particular if x is an N x p matrix and y an N x 1 vector, y is subtracted from each column of x, and if y=0 (default) you get the mean of  $\text{vec}(x^2)$

**Value**

The mean of the squared differences elementwise.

**Examples**

```
mse(2)
mse(1:10,2:11) == 1
mse(matrix(rnorm(500),100,5),matrix(rnorm(500),100,5))
```

---

norm_vec	<i>Norm of a vector</i>
----------	-------------------------

---

**Description**

Norm of a vector

**Usage**

```
norm_vec(x)
```

**Arguments**

x Numerical vector

**Value**

L2 norm of a vector

---

o2m *Perform O2PLS data integration with two-way orthogonal corrections*

---

**Description**

NOTE THAT THIS FUNCTION DOES NOT CENTER NOR SCALE THE MATRICES! Any normalization you will have to do yourself. It is best practice to at least center the variables though.

**Usage**

```
o2m(
  X,
  Y,
  n,
  nx,
  ny,
  stripped = FALSE,
  p_thresh = 3000,
  q_thresh = p_thresh,
  tol = 1e-10,
  max_iterations = 1000,
  sparse = F,
  groupx = NULL,
  groupy = NULL,
  keepx = NULL,
  keepy = NULL,
  max_iterations_sparsity = 1000
)
```

**Arguments**

X Numeric matrix. Vectors will be coerced to matrix with `as.matrix` (if this is possible)

Y Numeric matrix. Vectors will be coerced to matrix with `as.matrix` (if this is possible)

n Integer. Number of joint PLS components. Must be positive.

nx Integer. Number of orthogonal components in *X*. Negative values are interpreted as 0

<code>ny</code>	Integer. Number of orthogonal components in $Y$ . Negative values are interpreted as 0
<code>stripped</code>	Logical. Use the stripped version of o2m (usually when cross-validating)?
<code>p_thresh</code>	Integer. If $X$ has more than <code>p_thresh</code> columns, a power method optimization is used, see <a href="#">o2m2</a>
<code>q_thresh</code>	Integer. If $Y$ has more than <code>q_thresh</code> columns, a power method optimization is used, see <a href="#">o2m2</a>
<code>tol</code>	Double. Threshold for which the NIPALS method is deemed converged. Must be positive.
<code>max_iterations</code>	Integer. Maximum number of iterations for the NIPALS method.
<code>sparse</code>	Boolean. Default value is FALSE, in which case O2PLS will be fitted. Set to TRUE for GO2PLS.
<code>groupx</code>	Vector. Used when <code>sparse = TRUE</code> . A vector of strings indicating group names of each $X$ -variable. Its length must be equal to the number of variables in $X$ . The order of group names must corresponds to the order of the variables.
<code>groupy</code>	Vector. Used when <code>sparse = TRUE</code> . A vector of strings indicating group names of each $Y$ -variable. The length must be equal to the number of variables in $Y$ . The order of group names must corresponds to the order of the variables.
<code>keepx</code>	Vector. Used when <code>sparse = TRUE</code> . A vector of length $n$ indicating how many variables (or groups if <code>groupx</code> is provided) to keep in each of the joint component of $X$ . If the input is an integer, all the components will have the same amount of variables or groups retained.
<code>keepy</code>	Vector. Used when <code>sparse = TRUE</code> . A vector of length $n$ indicating how many variables (or groups if <code>groupx</code> is provided) to keep in each of the joint component of $Y$ . If the input is an integer, all the components will have the same amount of variables or groups retained.
<code>max_iterations_sparsity</code>	Integer. Used when <code>sparse = TRUE</code> . Maximum number of iterations for the NIPALS method for GO2PLS.

### Details

If both `nx` and `ny` are zero, `o2m` is equivalent to PLS2 with orthonormal loadings. This is a ‘slower’ (in terms of memory) implementation of O2PLS, and is using [svd](#), use `stripped=T` for a stripped version with less output. If either `ncol(X) > p_thresh` or `ncol(Y) > q_thresh`, the NIPALS method is used which does not store the entire covariance matrix. The squared error between iterands in the NIPALS approach can be adjusted with `tol`. The maximum number of iterations in the NIPALS approach is tuned by `max_iterations`.

### Value

A list containing

<code>Tt</code>	Joint $X$ scores
<code>W.</code>	Joint $X$ loadings
<code>U</code>	Joint $Y$ scores



C.	Joint $Y$ loadings
E	Residuals in $X$
Ff	Residuals in $Y$
T_Yosc	Orthogonal $X$ scores
P_Yosc.	Orthogonal $X$ loadings
W_Yosc	Orthogonal $X$ weights
U_Xosc	Orthogonal $Y$ scores
P_Xosc.	Orthogonal $Y$ loadings
C_Xosc	Orthogonal $Y$ weights
B_U	Regression coefficient in $Tt \sim U$
B_T.	Regression coefficient in $U \sim Tt$
H_TU	Residuals in $Tt$ in $Tt \sim U$
H_UT	Residuals in $U$ in $U \sim Tt$
X_hat	Prediction of $X$ with $Y$
Y_hat	Prediction of $Y$ with $X$
R2X	Variation (measured with <a href="#">ssq</a> ) of the modeled part in $X$ (defined by joint + orthogonal variation) as proportion of variation in $X$
R2Y	Variation (measured with <a href="#">ssq</a> ) of the modeled part in $Y$ (defined by joint + orthogonal variation) as proportion of variation in $Y$
R2Xcorr	Variation (measured with <a href="#">ssq</a> ) of the joint part in $X$ as proportion of variation in $X$
R2Ycorr	Variation (measured with <a href="#">ssq</a> ) of the joint part in $Y$ as proportion of variation in $Y$
R2X_Y0	Variation (measured with <a href="#">ssq</a> ) of the orthogonal part in $X$ as proportion of variation in $X$
R2Y_X0	Variation (measured with <a href="#">ssq</a> ) of the orthogonal part in $Y$ as proportion of variation in $Y$
R2Xhat	Variation (measured with <a href="#">ssq</a> ) of the predicted $X$ as proportion of variation in $X$
R2Yhat	Variation (measured with <a href="#">ssq</a> ) of the predicted $Y$ as proportion of variation in $Y$
W_gr	Joint loadings of $X$ at group level (only available when GO2PLS is used)
C_gr	Joint loadings of $Y$ at group level (only available when GO2PLS is used)

**See Also**

[summary.o2m](#), [plot.o2m](#), [crossval\\_o2m\\_adjR2](#), [crossval\\_sparsity](#)

## Examples

```

test_X <- scale(matrix(rnorm(100*10),100,10))
test_Y <- scale(matrix(rnorm(100*11),100,11))
# ----- Default run -----
o2m(test_X, test_Y, 3, 2, 1)
# ----- Stripped version -----
o2m(test_X, test_Y, 3, 2, 1, stripped = TRUE)
# ----- High dimensional version -----
o2m(test_X, test_Y, 3, 2, 1, p_thresh = 1)
# ----- High D and stripped version -----
o2m(test_X, test_Y, 3, 2, 1, stripped = TRUE, p_thresh = 1)
# ----- Now with more iterations -----
o2m(test_X, test_Y, 3, 2, 1, stripped = TRUE, p_thresh = 1, max_iterations = 1e6)
# -----

```

---

OmicsPLS

*Data integration with O2PLS: Two-Way Orthogonal Partial Least Squares*

---

## Description

The OmicsPLS package is an R package for penalized integration of heterogeneous omics data. The software articles are published in (el Bouhaddani et al, 2018, doi: [10.1186/s1285901823713](https://doi.org/10.1186/s1285901823713)) and (Gu et al, 2020, doi: [10.1186/s12859021039583](https://doi.org/10.1186/s12859021039583)). OmicsPLS includes the O2PLS fit, the GO2PLS fit, cross-validation tools and some misc functions.

## Model and assumptions

**Note that the rows of  $X$  and  $Y$  are the subjects and columns are variables.** The number of columns may be different, but the subjects should be the same in both datasets.

The O2PLS model (Trygg & Wold, 2003) decomposes two datasets  $X$  and  $Y$  into three parts.

- 1. A joint part, representing the relationship between  $X$  and  $Y$
- 2. An orthogonal part, representing the unrelated latent variation in  $X$  and  $Y$  separately.
- 3. A noise part capturing all residual variation.

See also the corresponding paper (el Bouhaddani et al, 2018).

## Fitting

The O2PLS fit is done with `o2m`. For data  $X$  and  $Y$  you can run `o2m(X, Y, n, nx, ny)` for an O2PLS fit with  $n$  joint and  $nx$ ,  $ny$  orthogonal components. See the help page of `o2m` for more information on parameters. There are four ways to obtain an O2PLS fit, depending on the dimensionality.

- For the not-too-high dimensional case, you may use `o2m` with default parameters. E.g. `o2m(X, Y, n, nx, ny)`.
- In case you only want the parameters, you may add `stripped = TRUE` to obtain a stripped version of `o2m` which avoids calculating and storing some matrices. E.g. `o2m(X, Y, n, nx, ny, stripped=TRUE)`.

- For high dimensional cases, defined by  $\text{ncol}(X) > p\_thresh$  and  $\text{ncol}(Y) > q\_thresh$ , a NI-PALS approach is used which avoids storing large matrices. E.g. `o2m(X, Y, n, nx, ny, p_thresh=3000, q_thresh=3000)`. The thresholds are by default both at 3000 variables.
- If you want a stripped version in the high dimensional case, add `stripped = TRUE`. E.g. `o2m(X, Y, n, nx, ny, stripped=TRUE, p_thresh=3000, q_thresh=3000)`.
- For GO2PLS, add `sparsity = TRUE` and specify how many variables or groups to retain. E.g. `o2m(X, Y, n, nx, ny, sparse=TRUE, keepx, keepy)`.

### Obtaining results

After fitting an O2PLS model, by running e.g. `fit = o2m(X, Y, n, nx, ny)`, the results can be visualised. Use `plot(fit, ...)` to plot the desired loadings with/without `ggplot2`. Use `summary(fit, ...)` to see the relative explained variances in the joint/orthogonal parts. Also plotting the joint scores `fit$Tt`, `fit$U` and orthogonal scores `fit$T_Yosc`, `fit$U_Xosc` are of help.

### Cross-validating

Determining the number of components  $n, nx, ny$  is an important task. For this we have two methods. See `citation("OmicsPLS")` for our proposed approach for determining the number of components, implemented in `crossval_o2m_adjR2`!

- Cross-validation (CV) is done with `crossval_o2m` and `crossval_o2m_adjR2`, both have built in parallelization which relies on the `parallel` package. Usage is something like `crossval_o2m(X, Y, a, ax, ay, nr_folds)` where `a, ax, ay` are vectors of integers. See the help pages. `nr_folds` is the number of folds, with `nr_folds = nrow(X)` for Leave-One-Out CV.
- For `crossval_o2m_adjR2` the same parameters are to be specified. This way of cross-validating is (potentially much) faster than the standard approach. It is also recommended over the standard CV.
- To cross-validate the number of variables to keep, use `crossval_sparsity`.

### S3 methods

There are S3 methods implemented for a fit obtained with `o2m`, i.e. `fit <- o2m(X, Y, n, nx, ny)`

- Use `plot(fit)` to plot the loadings, see above.
- Use `loadings(fit)` to extract a matrix with loading values
- Use `scores(fit)` to extract the scores
- Use `print` and `summary` to print and summarize the fit object

### Imputation

When the data contains missing values, one should impute them prior to using O2PLS. There are many sophisticated approaches available, such as MICE and MissForest, and no one approach is the best for all situations. To still allow users to quickly impute missing values in their data matrix, the `impute_matrix` function is implemented. It relies on the `softImpute` function+package and imputes based on the singular value decomposition.

**Misc**

Also some handy tools are available

- `orth(X)` is a function to obtain an orthogonalized version of a matrix or vector  $X$ .
- `ssq(X)` is a function to calculate the sum of squares (or squared Frobenius norm) of  $X$ . See also `vnorm` for calculating the norm of each column in  $X$ .
- `mse(x, y)` returns the mean squared difference between two matrices/vectors.

**Citation**

If you use the OmicsPLS R package in your research, please cite the corresponding software paper:

**el Bouhaddani, S., Uh, H.-W., Jongbloed, G., Hayward, C., Klarić, L., Kielbasa, S. M., & Houwing-Duistermaat, J.** (2018). *Integrating omics datasets with the OmicsPLS package*. BMC Bioinformatics, 19(1). doi: [10.1186/s1285901823713](https://doi.org/10.1186/s1285901823713)

The bibtex entry can be obtained with command `citation("OmicsPLS")`. Thank you!

The original paper proposing O2PLS is

**Trygg, J., & Wold, S.** (2003). *O2-PLS, a two-block (X-Y) latent variable regression (LVR) method with an integral OSC filter*. Journal of Chemometrics, 17(1), 53-64. doi: [10.1002/cem.775](https://doi.org/10.1002/cem.775)

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Maintainer: Said el Bouhaddani (<[s.elbouhaddani@umcutrecht.nl](mailto:s.elbouhaddani@umcutrecht.nl)>).

---

orth	<i>Orthogonalize a matrix</i>
------	-------------------------------

---

**Description**

Orthogonalize a matrix

**Usage**

```
orth(X, X_true = NULL, type = c("QR", "SVD"))
```

**Arguments**

X	Numeric vector or matrix.
X_true	(optional) A 'true' matrix/vector. Used to correct the sign of the orthonormalized X if QR is used. Only the first column is corrected.
type	A character or numeric. Should be one of "QR" or "SVD".

**Details**

Choosing `type='QR'` uses a QR decomposition of  $X$  to produce orthonormal columns. For `type=='SVD'` it uses an SVD decomposition. The columns are corrected for sign.

**Value**

An orthogonalized representation of  $X$

**Examples**

```
orth(c(3,4))
round(crossprod(orth(matrix(rnorm(500),100,5))),4)
orth(matrix(1:9,3,3),type='QR')[,1] - orth(1:3); orth(matrix(1:9,3,3),type='SVD')[,1] - orth(1:3);
```

---

 orth\_vec

---

*Orthogonalize a sparse loading vector with regard to a matrix*


---

**Description**

Orthogonalize a sparse loading vector with regard to a matrix

**Usage**

```
orth_vec(x, W)
```

**Arguments**

`x` sparse loading vector to be orthogonalized  
`W` sparse loading matrix of the previous loading vectors

**Value**

A sparse loading vector

---

plot.o2m

---

*Plot one or two loading vectors for class o2m*


---

### Description

This function plots one or two loading vectors, by default with ggplot2.

### Usage

```
## S3 method for class 'o2m'
plot(
  x,
  loading_name = c("Xjoint", "Yjoint", "gr_Xjoint", "gr_Yjoint", "Xorth", "Yorth"),
  i = 1,
  j = NULL,
  use_ggplot2 = TRUE,
  label = c("number", "colnames"),
  ...
)
```

### Arguments

x	An O2PLS fit, with class 'o2m'
loading_name	character string. One of the following: 'Xjoint', 'Yjoint', 'gr_Xjoint', 'gr_Yjoint', 'Xorth' or 'Yorth'.
i	Integer. First component to be plotted.
j	NULL (default) or Integer. Second component to be plotted.
use_ggplot2	Logical. Default is TRUE. If FALSE, the usual plot device will be used.
label	Character, either 'number' or 'colnames'. The first option prints numbers, the second prints the colnames
...	Further arguments to geom_text, such as size, col, alpha, etc.

### Value

If use\_ggplot2 is TRUE a ggplot2 object. Else NULL.

### See Also

[summary.o2m](#)

---

predict.o2m	<i>Predicts X or Y</i>
-------------	------------------------

---

**Description**

Predicts X or Y based on new data on Y or X

**Usage**

```
## S3 method for class 'o2m'  
predict(object, newdata, XorY = c("X", "Y"), ...)
```

**Arguments**

object	List. Should be of class <code>o2m</code> .
newdata	New data, which one of X or Y is specified in XorY.
XorY	Character specifying whether newdata is X or Y.
...	For compatibility

**Details**

Prediction is done after correcting for orthogonal parts.

**Value**

Predicted Data

**Examples**

```
predict(o2m(scale(1:10), scale(1:10), 1, 0, 0), newdata = scale(1:5), XorY = "X")
```

---

print.cvo2m	<i>Cross-validate procedure for O2PLS</i>
-------------	---

---

**Description**

Cross-validate procedure for O2PLS

**Usage**

```
## S3 method for class 'cvo2m'  
print(x, include_matrix = FALSE, ...)
```

**Arguments**

x List of class "cvo2m", produced by [crossval\\_o2m](#).  
 include\_matrix Logical. Should the 3-d array with Prediction errors also be printed.  
 ... For consistency.

---

print.o2m *Print function for O2PLS.*

---

**Description**

This function is the print method for an O2PLS fit

**Usage**

```
## S3 method for class 'o2m'
print(x, ...)
```

**Arguments**

x An O2PLS fit (an object of class o2m)  
 ... For consistency

---

print.pre.o2m *Print function for O2PLS.*

---

**Description**

This function is the print method for an O2PLS fit

**Usage**

```
## S3 method for class 'pre.o2m'
print(x, ...)
```

**Arguments**

x An O2PLS fit (an object of class o2m)  
 ... For consistency



---

rmsep	<i>Root MSE of Prediction</i>
-------	-------------------------------

---

**Description**

Calculates the Root MSE of prediction on test data. Only tested to work inside [loocv](#).

**Usage**

```
rmsep(Xtst, Ytst, fit, combi = FALSE)
```

**Arguments**

Xtst	Numeric vector or matrix.
Ytst	Numeric vector or matrix.
fit	<a href="#">o2m</a> fit (on data without Xtst and Ytst).
combi	Logical. Should the symmetrized MSE be used, i.e. add both MSEs. Not yet implemented, but see <a href="#">rmsep_combi</a>

**Details**

This function is the building block for [loocv](#), as it produced the prediction error for test (left out) data.

**Value**

Mean squares difference between predicted Y and true Y

---

rmsep_combi	<i>Symmetrized root MSE of Prediction</i>
-------------	---

---

**Description**

Calculates the symmetrized root MSE of prediction on test data. \*Expected\* to work in combination with [loocv](#).

**Usage**

```
rmsep_combi(Xtst, Ytst, fit)
```

**Arguments**

Xtst	Numeric vector or matrix.
Ytst	Numeric vector or matrix.
fit	<a href="#">o2m</a> fit (on data without Xtst and Ytst).

**Details**

This function is the building block for `loocv`, as it produced the prediction error for test (left out) data.

This is a symmetrized version of `rmsep`, and is used by `loocv`. The prediction error of both `Xtst` and `Ytst` are calculated and summed. Whether this is a good idea depends: If  $X$  and  $Y$  have similar meanings (LC-MS versus MALDI) this is a good thing to do. If the two matrices do not have similar meanings, (Metabolomics versus Transcriptomics) then you may want to not sum up the two prediction errors or include weights in the sum.

**Value**

Mean squares difference between predicted  $Y$  and true  $Y$

---

scores	<i>Extract the scores from an O2PLS fit</i>
--------	---

---

**Description**

This function extracts score matrices from an O2PLS fit

**Usage**

```
scores(x, ...)  
  
## S3 method for class 'o2m'  
scores(  
  x,  
  which_part = c("Xjoint", "Yjoint", "Xorth", "Yorth"),  
  subset = 0,  
  ...  
)
```

**Arguments**

<code>x</code>	Object of class <code>o2m</code>
<code>...</code>	For consistency
<code>which_part</code>	character string. One of the following: 'Xjoint', 'Yjoint', 'Xorth' or 'Yorth'.
<code>subset</code>	subset of scores vectors to be extracted.

**Value**

Scores matrix

**See Also**

[loadings.o2m](#)

**Examples**

```
scores(o2m(scale(-2:2),scale(-2:2*4),1,0,0))
```

---

ssq

*Calculate Sum of Squares*


---

**Description**

Calculate Sum of Squares

**Usage**

```
ssq(X)
```

**Arguments**

*X*                    Numeric vector or matrix.

**Details**

This is the Frobenius norm of *X*.

**Value**

The sum of squared elements of *X*

**Examples**

```
ssq(tcrossprod(1:5))
ssq(rnorm(1e5))/1e5
```

---

summary.o2m

*Summary of an O2PLS fit*


---

**Description**

Until now only variational summary given by the R2's is outputted

**Usage**

```
## S3 method for class 'o2m'
summary(object, digits = 3, ...)
```

**Arguments**

object	List. Should be of class <code>o2m</code> .
digits	Integer, number of digits.
...	For compatibility

**Value**

List with R2 values.

**See Also**

[plot.o2m](#)

**Examples**

```
summary(o2m(scale(-2:2), scale(-2:2*4), 1, 0, 0))
```

---

thresh_n	<i>Soft thresholding a vector with respect to a number of variables</i>
----------	---

---

**Description**

Soft thresholding a vector with respect to a number of variables

**Usage**

```
thresh_n(x, keepx)
```

**Arguments**

x	Numerical vector
keepx	How many non-zero

**Value**

Soft-thresholded vector

---

thresh_n_gr	<i>Soft thresholding a vector with respect to a number of groups</i>
-------------	--

---

**Description**

Soft thresholding a vector with respect to a number of groups

**Usage**

```
thresh_n_gr(w, keep_gr, index_gr)
```

**Arguments**

w	Numerical loading vector
keep_gr	How many groups to retain
index_gr	List of index and size. index are the index of variables belongs to the group in the original vector, size is the group size

**Value**

A list containing sparse loading vector and names of the selected groups

---

vnorm	<i>Norm of a vector or columns of a matrix</i>
-------	--

---

**Description**

Norm of a vector or columns of a matrix

**Usage**

```
vnorm(x)
```

**Arguments**

x	Numeric vector or matrix.
---	---------------------------

**Value**

(columnwise) Euclidian norm of  $x$

**Examples**

```
vnorm(orth(1:5))
vnorm(matrix(1:9,3,3))^2 - colSums(matrix(1:9,3)^2)
```

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