

# Package ‘abundant’

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**Type** Package

**Title** High-Dimensional Principal Fitted Components and Abundant Regression

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**Depends** R (>= 2.10), glasso

**Description** Fit and predict with the high-dimensional principal fitted components model. This model is described by Cook, Forzani, and Rothman (2012) <[doi:10.1214/11-AOS962](https://doi.org/10.1214/11-AOS962)>.

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abundant-package	<i>Abundant regression and high-dimensional principal fitted components</i>
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## Description

Fit and predict with the high-dimensional principal fitted components model.

## Details

The main functions are `fit.pfc`, `pred.response`.

## Author(s)

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## References

Cook, R. D., Forzani, L., and Rothman, A. J. (2012). Estimating sufficient reductions of the predictors in abundant high-dimensional regressions. *Annals of Statistics* 40(1), 353-384.

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fit.pfc

*Fit a high-dimensional principal fitted components model using the method of Cook, Forzani, and Rothman (2012).*

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## Description

Let  $(x_1, y_1), \dots, (x_n, y_n)$  denote the  $n$  measurements of the predictor and response, where  $x_i \in R^p$  and  $y_i \in R$ . The model assumes that these measurements are a realization of  $n$  independent copies of the random vector  $(X, Y)'$ , where

$$X = \mu_X + \Gamma\beta\{f(Y) - \mu_f\} + \epsilon,$$

$\mu_X \in R^p$ ;  $\Gamma \in R^{p \times d}$  with rank  $d$ ;  $\beta \in R^{d \times r}$  with rank  $d$ ;  $f : R \rightarrow R^r$  is a known vector valued function;  $\mu_f = E\{f(Y)\}$ ;  $\epsilon \sim N_p(0, \Delta)$ ; and  $Y$  is independent of  $\epsilon$ . The central subspace is  $\Delta^{-1}\text{span}(\Gamma)$ .

This function computes estimates of these model parameters by imposing constraints for identifiability. The mean parameters  $\mu_X$  and  $\mu_f$  are estimated with  $\bar{x} = n^{-1} \sum_{i=1}^n x_i$  and  $\bar{f} = n^{-1} \sum_{i=1}^n f(y_i)$ . Let  $\hat{\Phi} = n^{-1} \sum_{i=1}^n \{f(y_i) - \bar{f}\}\{f(y_i) - \bar{f}\}'$ , which we require to be positive definite. Given a user-specified weight matrix  $\hat{W}$ , let

$$(\hat{\Gamma}, \hat{\beta}) = \arg \min_{G \in R^{p \times d}, B \in R^{d \times r}} \sum_{i=1}^n [x_i - \bar{x} - GB\{f(y_i) - \bar{f}\}]' \hat{W} [x_i - \bar{x} - GB\{f(y_i) - \bar{f}\}],$$

subject to the constraints that  $G' \hat{W} G$  is diagonal and  $B \hat{\Phi} B' = I$ . The sufficient reduction estimate  $\hat{R} : R^p \rightarrow R^d$  is defined by

$$\hat{R}(x) = (\hat{\Gamma}' \hat{W} \hat{\Gamma})^{-1} \hat{\Gamma}' \hat{W} (x - \bar{x}).$$

## Usage

```
fit.pfc(X, y, r=4, d=NULL, F.user=NULL, weight.type=c("sample", "diag", "L1"),
        lam.vec=NULL, kfold=5, silent=TRUE, qrtol=1e-10, cov.tol=1e-4,
        cov.maxit=1e3, NPERM=1e3, level=0.01)
```

**Arguments**

<code>X</code>	The predictor matrix with $n$ rows and $p$ columns. The $i$ th row is $x_i$ defined above.
<code>y</code>	The vector of measured responses with $n$ entries. The $i$ th entry is $y_i$ defined above.
<code>r</code>	When polynomial basis functions are used (which is the case when <code>F.user=NULL</code> ), $r$ is the polynomial order, i.e., $f(y) = (y, y^2, \dots, y^r)'$ . The default is $r=4$ . This argument is not used when <code>F.user</code> is specified.
<code>d</code>	The dimension of the central subspace defined above. This must be specified by the user when <code>weight.type="L1"</code> . If unspecified by the user this function will use the sequential permutation testing procedure, described in Section 8.2 of Cook, Forzani, and Rothman (2012), to select $d$ .
<code>F.user</code>	A matrix with $n$ rows and $r$ columns, where the $i$ th row is $f(y_i)$ defined above. This argument is optional, and will typically be used when polynomial basis functions are not desired.
<code>weight.type</code>	The type of weight matrix estimate $\widehat{W}$ to use. Let $\widehat{\Delta}$ be the observed residual sample covariance matrix for the multivariate regression of $X$ on $f(Y)$ with $n - r - 1$ scaling. There are three options for $\widehat{W}$ : <ul style="list-style-type: none"> <li>• <code>weight.type="sample"</code> uses a Moore-Penrose generalized inverse of <math>\widehat{\Delta}</math> for <math>\widehat{W}</math>, when <math>p \leq n - r - 1</math> this becomes the inverse of <math>\widehat{\Delta}</math>;</li> <li>• <code>weight.type="diag"</code> uses the inverse of the diagonal matrix with the same diagonal as <math>\widehat{\Delta}</math> for <math>\widehat{W}</math>;</li> <li>• <code>weight.type="L1"</code> uses the L1-penalized inverse of <math>\widehat{\Delta}</math> described in equation (5.4) of Cook, Forzani, and Rothman (2012). In this case, <code>lam.vec</code> and <code>d</code> must be specified by the user. The glasso algorithm of Friedman et al. (2008) is used through the R package <code>glasso</code>.</li> </ul>
<code>lam.vec</code>	A vector of candidate tuning parameter values to use when <code>weight.type="L1"</code> . If this vector has more than one entry, then <code>kfold</code> cross validation will be performed to select the optimal tuning parameter value.
<code>kfold</code>	The number of folds to use in cross-validation to select the optimal tuning parameter when <code>weight.type="L1"</code> . Only used if <code>lam.vec</code> has more than one entry.
<code>silent</code>	Logical. When <code>silent=FALSE</code> , progress updates are printed.
<code>qrtol</code>	The tolerance for calls to <code>qr.solve()</code> .
<code>cov.tol</code>	The convergence tolerance for the QUIC algorithm used when <code>weight.type="L1"</code> .
<code>cov.maxit</code>	The maximum number of iterations allowed for the QUIC algorithm used when <code>weight.type="L1"</code> .
<code>NPERM</code>	The number of permutations to used in the sequential permutation testing procedure to select $d$ . Only used when <code>d</code> is unspecified.
<code>level</code>	The significance level to use to terminate the sequential permutation testing procedure to select $d$ .

**Details**

See Cook, Forzani, and Rothman (2012) more information.

**Value**

A list with

Gamhat	this is $\widehat{\Gamma}$ described above.
bhat	this is $\widehat{\beta}$ described above.
Rmat	this is $\widehat{W}\widehat{\Gamma}(\widehat{\Gamma}'\widehat{W}\widehat{\Gamma})^{-1}$ .
What	this is $\widehat{W}$ described above.
d	this is $d$ described above.
r	this is $r$ described above.
GWG	this is $\widehat{\Gamma}'\widehat{W}\widehat{\Gamma}$
fc	a matrix with $n$ rows and $r$ columns where the $i$ th row is $f(y_i) - \bar{f}$ .
Xc	a matrix with $n$ rows and $p$ columns where the $i$ th row is $x_i - \bar{x}$ .
y	the vector of $n$ response measurements.
mx	this is $\bar{x}$ described above.
mf	this is $\bar{f}$ described above.
best.lam	this is selected tuning parameter value used when <code>weight.type="L1"</code> , will be NULL otherwise.
lam.vec	this is the vector of candidate tuning parameter values used when <code>weight.type="L1"</code> , will be NULL otherwise.
err.vec	this is the vector of validation errors from cross validation, one error for each entry in <code>lam.vec</code> . Will be NULL unless <code>weight.type="L1"</code> and <code>lam.vec</code> has more than one entry.
test.info	a dataframe that summarizes the results from the sequential testing procedure. Will be NULL unless <code>d</code> is unspecified.

**Author(s)**

Adam J. Rothman

**References**

- Cook, R. D., Forzani, L., and Rothman, A. J. (2012). Estimating sufficient reductions of the predictors in abundant high-dimensional regressions. *Annals of Statistics* 40(1), 353-384.
- Friedman, J., Hastie, T., and Tibshirani R. (2008). Sparse inverse covariance estimation with the lasso. *Biostatistics* 9(3), 432-441.

**See Also**

[pred.response](#)

**Examples**

```

set.seed(1)
n=20
p=30
d=2
y=sqrt(12)*runif(n)
Gam=matrix(rnorm(p*d), nrow=p, ncol=d)
beta=diag(2)
E=matrix(0.5*rnorm(n*p), nrow=n, ncol=p)
V=matrix(c(1, sqrt(12), sqrt(12), 12.8), nrow=2, ncol=2)
tmp=eigen(V, symmetric=TRUE)
V.msqr=tcrossprod(tmp$vec*rep(tmp$val^(-0.5), each=2), tmp$vec)
Fyc=cbind(y-sqrt(3), y^2-4)%*%V.msqr
X=0+Fyc%*%t(beta)%*%t(Gam) + E

fit=fit.pfc(X=X, y=y, r=3, weight.type="sample")
## display hypothesis testing information for selecting d
fit$test.info
## make a response versus fitted values plot
plot(pred.response(fit), y)

```

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pred.response	<i>Predict the response with the fitted high-dimensional principal fitted components model</i>
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**Description**

Let  $x \in R^p$  denote the values of the  $p$  predictors. This function computes  $\hat{E}(Y|X = x)$  using equation (8.1) of Cook, Forzani, and Rothman (2012).

**Usage**

```
pred.response(fit, newx=NULL)
```

**Arguments**

fit	The object returned by <code>fit.pfc()</code> .
newx	A matrix with $N$ rows and $p$ columns where each row is an instance of $x$ described above. If this argument is unspecified, then the fitted values are returned, i.e, <code>newx=X</code> , where $X$ was the predictor matrix used in the call to <code>fit.pfc()</code> .

**Details**

See Cook, Forzani, and Rothman (2012) for more information.

**Value**

A vector of response prediction with `nrow(newx)` entries.

**Author(s)**

Adam J. Rothman

**References**

Cook, R. D., Forzani, L., and Rothman, A. J. (2012). Estimating sufficient reductions of the predictors in abundant high-dimensional regressions. *Annals of Statistics* 40(1), 353-384.

**See Also**

[fit.pfc](#)

**Examples**

```
set.seed(1)
n=25
p=50
d=1
true.G = matrix(rnorm(p*d), nrow=p, ncol=d)
y=rnorm(n)
fy = y
E=matrix(rnorm(n*p), nrow=n, ncol=p)
X=fy%*%t(true.G) + E
fit=fit.pfc(X=X, r=4, d=d, y=y, weight.type="diag")
fitted.values=pred.response(fit)
mean((y-fitted.values)^2)
plot(fitted.values, y)

n.new=100
y.new=rnorm(n.new)
fy.new=y.new
E.new=matrix(rnorm(n.new*p), nrow=n.new, ncol=p)
X.new = fy.new%*%t(true.G) + E.new
mean((y.new - pred.response(fit, newx=X.new))^2)
```

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