# Section 9: Matching without replacement, Genetic matching

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# Matching without replacement

- The standard approach is to minimize the Mahalanobis distance matrix (In GenMatch we use a weighted Mahalanobis distance)
- We can also use other distance functions such as Euclidean distance, however Mahalanobis distance has attractive properties
- Matching with replacement will yield the same results using Euclidean, absolute, or Mahalanobis distance
- Next we consider Matching using Mahalnobis distance without replacement

# Matching without replacement using Mahalanobis distance

- Suppose we have  $X_1, \ldots, X_n$  with each variable  $X_i$  having k components (covariates):  $X_i = (X_{i1}, X_{i2}, \ldots, X_{ik})$
- Suppose we have the  $k \times k$  covariance matrix of these covariates. (either by estimating this matrix, or given to us). Call this covariance matrix S
- The Mahalanobis distance between  $X_i$  and  $X_j$  is,

$$md(X_i, X_j) = \sqrt{[(X_i - X_j)S^{-1}(X_i - X_j)]}$$

- How to implement Mahalanobis distance in *R*:
  - **1** The *Match* function in the package *Matching*.
  - Write your own function that calculates Mahalanobis distance, and creates matches to minimizes the distance matrix between the treated and the control units
- The code for both options is in the next slides using data from Dehejia and Wahba (1999). The results are very similar, the differences result in the minimization of the Mahalanobis distance matrix

Dehejia and Wahba (1999)

```
### Using Jas's "Match" function:
match = Match(Tr=treat1,X=x1,replace=FALSE,Weight=2)
# or using the option "Weight=1", there is not much difference
dm2 = d1[c(match$index.treated,match$index.control),]
```

```
smahal = function(z.X) {
  X = as.matrix(X)
  n = dim(X)[1]
  rownames (X) < -1:n
  k < -dim(X)[2]
  m < -sum(z)
  for (j in 1:k) {
    X[,j] = rank(X[,j])
  cv \leftarrow cov(X)
  vuntied <- var(1:n)
  rat <- sqrt(vuntied/diag(cv))</pre>
  cv <- diag(rat)%*%cv%*%diag(rat)</pre>
  # distance matrix containing: rows are the treatment group and columns are th
  out <- matrix(NA,m,n-m)
  Xc = X[z==0,]; Xt = X[z==1,];
  rownames(out)=rownames(X)[z==1]
  colnames(out)=rownames(X)[z==0]
  library(MASS)
  icov <- ginv(cv)</pre>
  for (i in 1:m) {
    out[i,] = mahalanobis(Xc,Xt[i,],icov,inverted=T)
  }
  out }
```

#### Dehejia and Wahba (1999)

```
source("smahal.R")
### data without the experimental controls with CPS 3
d1 = read.dta("cps1re74.dta")
### CPS-1 vs NSW treatment group ###
x1 = d1[,c("age","ed","black","hisp","married","nodeg","re74","r
treat1 = d1$treat
# Calculating the Mahalanobis distance matrix using the function
distance.matrix = smahal(treat1,x1)
library("optmatch")
pair.match <-pairmatch(distance.matrix, controls=1,data=d1)</pre>
# The matched data set:
dm = d1[is.na(pair.match)=="FALSE",]
```

	Jas's Match function		Code using package optmatch	
	Ave. control	T-test	Ave. control	T-test
age	25.189	0.445	25.330	0.562
ed	10.314	0.878	10.276	0.755
black	0.843	1.000	0.832	0.779
hisp	0.059	1.000	0.059	1.000
married	0.189	1.000	0.189	1.000
nodeg	0.708	1.000	0.708	1.000
re74	2631.729	0.270	2559.173	0.352
re75	2215.897	0.049	2211.832	0.060

- The treated units are not presented as they are the same in both matched data sets.
- The difference in the results follows from the numerical procedure used to minimize the Mahalanobis distance matrix. The Match package and the optmatch package use different algorithms.
- I recommend using the *Match* function, but understand what it does

### What is the Mahalanobis distance doing: A special case

- Suppose covariates are ellipsoidal (in a crude sense, the data looks like an oval).
- Special case of ellipsoidal: Multivariate normal
- Transform the covariates into a circle, where the marginal SD is 1 for any covariate
- Equivalent to "Eliminating the correlation between covariates and scaling all covariates so that they are on the same scale"
- Calculate the distance within this circle
- Mahalanobis distance can be computed for any other original shape of data... but nice properties can be derived if the data is ellipsoidal

# Equal Percent Bias Reduction (EPBR)

- Suppose we are matching treated units to control units
- matching method is Equal Percent Bias Reducing (EPBR) if, for ALL covariates X:

$$\mathbb{E}(X|T=1) - \mathbb{E}(X|\textit{Matched Controls})$$
  
= $\gamma (\mathbb{E}(X|T=1) - \mathbb{E}(X|T=1))$ 

where  $0 \ge \gamma \le 1$ .

- $\bullet$  In words, the average covariate imbalance shrinks by  $\gamma$  for each covariate
- If data is ellipsoidal, then the matching with the Mahalanobis distance is EPBR
- EPBR may not be that attractive of a property; especially if you think that balance over some covariates (height, weight) is more important than others (number of hair follicles)
- EPBR shrinks the bias only on average, the bias in other percentiles is not guaranteed to reduce

### General notes on matching

- EPBR does not mean that your estimated treatment effect will be less biased. It could be more biased if you do not have the right X s. Moreover, balance on covariates that you do not include could get worse
- Measurement error in some of the covariates can bias the matching treatment effect estimations. Is it attenuation bias?
- Is the following suggestion a good placebo test?

"As a placebo test, redo the matching without the previous outcome and test the previous outcome. The two groups should be balanced on the previous outcome"

•

#### How can you tell if matching worked?

- Matching has necessary testable implication, which can be verified in the data. This are only necessary conditions, <u>not sufficient</u> conditions for inference. We cannot test for balance in unobservable...
- The objective of matching is to find people in the control group that look like people that are treated
- If matching works, the distribution of the covariates for the treated group should be the same as the distribution for the matched control group
- If data is quantitative, compare means of covariates between the treatment group and the matched control group (t-test)
- If data is categorical, compare proportions across the covariates. (Fisher exact test)
- These methods compare single points (a mean or proportion) between these two groups
- Kolmogorov-Smirnov test can test if theres discrepancies across the entire distribution

# Example: Kane and Rouse (1995), Labor-Market Returns to Two- and Four-Year College

- We will use the paper, "Labor-Market Returns to Two-Year and Four-Year College" by Kane and Rouse (1995) to demonstrate matching without replacement and Genetic matching
- The data:
  - Outcome: educ86, Years of Education Completed (as of 1986) Treatment: twoyr (1 if started in a two year college, O if started in a four year college.
- We are interested in the effect of the treatment (two-years collage) on the outcome (years of education completed)

# Example: Kane and Rouse (1995)

	Ave. Treat	Ave. control	T-test	Wilcoxon	KS
female	0.501	0.510	0.757	0.757	1.000
black	0.056	0.098	0.002	0.007	0.611
hispanic	0.135	0.096	0.035	0.023	0.709
bytest	59.255	60.919	0.000	0.000	0.000
dadmiss	0.131	0.119	0.525	0.515	1.000
dadvoc	0.093	0.069	0.123	0.097	0.991
dadsome	0.147	0.114	0.091	0.073	0.882
dadcoll	0.210	0.346	0.000	0.000	0.000
mommiss	0.068	0.040	0.035	0.016	0.959
momvoc	0.098	0.088	0.536	0.525	1.000
momsome	0.159	0.159	0.976	0.976	1.000
momcoll	0.145	0.246	0.000	0.000	0.002
fincome	23057.110	26534.917	0.000	0.007	0.026
fincmiss	0.051	0.061	0.426	0.446	1.000
ownhome	0.823	0.844	0.300	0.286	0.998
perwhite	75.956	79.176	0.027	0.002	0.008
urban	0.166	0.221	0.009	0.013	0.264

# Example: Kane and Rouse (1995)

```
female (1 = Female/0 = Male)
black (1 = Black/0 = Not-Black)
hispanic (1 = Hispanic/0 = Not-Hispanic)
bytest (Base Year Composite Test Score)
Father's Highest Education Level:
 dadmiss1 (Education Missing)
 dadvoc (1 = Father attended Vocational/Technical School, 0 if not)
 dadsome (1 = Father attended some college but did not graduate, 0 if not)
 dadcoll (1 = Father college graduate, 0 if not)
Mother's Highest Education Level:
 mommiss (Education Missing)
 momvoc (1= Mother attended Vocational/Technical School, 0 if not)
 momsome (1= Mother attended some college but did not graduate, 0 if not)
momcoll (1= Mother College Graduate, 0 if not)
fincome (Family Income)
fincmiss (Family Income Missing)
ownhome (1= Family Owns Home, 0 if not)
perwhite (Percentage of Student Body White in students high school)
urban (1 if students high school in Urban Area)
cue80 (County unemployment rate in 1980
stypc80 (State Personal Income per capita)
stwmfg80 (State Hourly Wage (mfg), 1980)
```

# Example: Kane and Rouse (1995)

Model 1	Model 2	NAI - I 2
		Model 3
15.30***	11.04***	11.20***
(0.04)	(0.61)	(0.63)
-1.36***	-1.22***	-1.21***
(80.0)	(80.0)	(80.0)
, ,	0.03	0.03
	(0.07)	(0.07)
	-0.23	_0.27 <sup>*</sup>
	(0.13)	(0.14)
		0.00*
		(0.00)
	` ,	0.06***
		(0.01)
	0.03	-0.03
	(0.09)	(0.09)
	()	0.14
		(0.13)
		. ,
		:
0.13	0.16	0.19
0.13	0.16	0.18
1818	1818	1818
	-1.36*** (0.08)	(0.04) (0.61) -1.36*** (0.08) (0.08) (0.08) 0.03 (0.07) -0.23 (0.13) 0.00*** (0.00) 0.07*** (0.01) 0.03 (0.09)

<sup>15/59</sup> 

# Example: Kane and Rouse (1995), matching w/t replacement

	Ave. Treat	Ave. control	T-test	Wilcoxon	KS
female	0.501	0.517	0.633	0.633	1.000
black	0.056	0.054	0.881	0.881	1.000
hispanic	0.135	0.121	0.541	0.540	1.000
bytest	59.255	59.689	0.040	0.019	0.040
dadmiss	0.131	0.119	0.606	0.606	1.000
dadvoc	0.093	0.093	1.000	1.000	1.000
dadsome	0.147	0.138	0.696	0.696	1.000
dadcoll	0.210	0.212	0.933	0.934	1.000
mommiss	0.068	0.068	1.000	1.000	1.000
momvoc	0.098	0.093	0.817	0.817	1.000
momsome	0.159	0.156	0.925	0.926	1.000
momcoll	0.145	0.142	0.923	0.923	1.000
fincome	23057.110	22638.695	0.701	0.894	1.000
fincmiss	0.051	0.051	1.000	1.000	1.000
ownhome	0.823	0.848	0.311	0.311	0.999
perwhite	75.956	79.883	0.028	0.002	0.012
urban	0.166	0.166	1.000	1.000	1.000

# Example: Kane and Rouse (1995), Regressions after matching

	Model 1	Model 2	Model 3		
(Intercept)	15.18***	10.25***	10.33***		
	(80.0)	(1.07)	(1.10)		
twoyr	-1.24***	-1.21***	-1.22***		
	(0.11)	(0.11)	(0.11)		
female		-0.10	-0.09		
		(0.11)	(0.11)		
black		-0.35	-0.45		
		(0.24)	(0.25)		
fincome		0.00***	0.00		
		(0.00)	(0.00)		
bytest		0.08***	0.08***		
		(0.02)	(0.02)		
urban		-0.07	-0.17		
		(0.15)	(0.15)		
hispanic			0.28		
			(0.21)		
			-		
:			<u>:</u>		
$R^2$	0.13	0.16	0.19		
Adj. R <sup>2</sup>	0.13	0.16	0.17		
Num. obs.	858	858	858		
*** <i>p</i> < 0.001, ** <i>p</i> < 0.01, * <i>p</i> < 0.05					

 $<sup>0.001, \ \</sup> p < 0.01, \ \ p < 0.05$ 

- The objective of matching is to create comparable treatment and control units
- Matching on the Propensity score or using Mahalanobis distance, or a mix using both, do we maximize balance? No, they optimize on P-score or on Mahalanobis distance
- Genetic Matching incorporates balancing directly in the algorithm
  - Generalizing Mahalanobis matching
  - Prioritizes differences on certain covariates
  - ullet Optimizes matching for some target: balance on  $X_p$

- Why weight? Some covariates may be more important than others
- Intuitively, we already weight covariates either at 0 or 1 in measuring  $md(X_i, X_j)$
- Define  $X_p$  as a full matrix of covariates and  $X_k$  as the k-dimensional matrix we match on, where k < p
- Let  $\tilde{w}$  be a vector of weights so that,

$$\tilde{w} = \{w_1 = 0, w_2 = 0, \dots, w_k = 0, w_{k+1} = 1, \dots, w_p = 1\}$$

 $\bullet$  Generalize this weighting in a flexible way, Let  $\tilde{w} \in \{\underline{w}, \overline{w}\}$ 

• Incorporate these weights in  $W_{p \times p}$  to rescale unit distance between unit i and unit j,

$$wmd(X_i, X_j) = \sqrt{(X_i - X_j)^T \left(S^{-\frac{1}{2}}\right)^T W\left(S^{-\frac{1}{2}}\right) (X_i - X_j)}$$

- Where W is a  $p \times p$  positive definite weight matrix, with  $\tilde{w}$  the diagonal elements of W, the off diagonal element of W are 0
- "The choice of setting the non-diagonal elements of W to zero is made for reasons of computational power alone", Mebane and Sekhon (2011)
- $\left(S^{\frac{1}{2}}\right)$  is the Cholesky decomposition of the sample covariance matrix S,

$$S = \left(S^{-rac{1}{2}}
ight)^T \left(S^{-rac{1}{2}}
ight)$$

• When  $W = I_{p \times p}$ , then  $wmd(X_i, X_j) = md(X_i, X_j)$ 

Suppose we have,

$$(x_1,x_2) \sim N \left( (0,0), \underbrace{\left( \begin{array}{cc} 1 & 0 \\ 0 & 1 \end{array} \right)}_{S} \right)$$

This is equivalent to  $x_1 \sim N(0,1)$ ,  $x_2 \sim N(0,1)$ , and  $Cov(x_1,x_2)=0$ . Then,

$$wmd(X_{i}, X_{j}) = \sqrt{(X_{i} - X_{j})^{T} \left(S^{-\frac{1}{2}}\right)^{T} W\left(S^{-\frac{1}{2}}\right) (X_{i} - X_{j})}$$

$$= \sqrt{(X_{i} - X_{j})^{T} W(X_{i} - X_{j})}$$

$$= \sqrt{(x_{1i} - x_{1j}, x_{2i} - x_{2j}) \begin{pmatrix} w_{1} & 0 \\ 0 & w_{2} \end{pmatrix} \begin{pmatrix} x_{1i} - x_{1j} \\ x_{2i} - x_{2j} \end{pmatrix}}$$

$$= \sqrt{(x_{1i} - x_{1j})^{2} w_{1} + (x_{2i} - x_{2j})^{2} w_{2}}$$

• If  $w_1 = 2w_2$  then,

$$wmd(X_i, X_j) = \sqrt{(x_{1i} - x_{1j})^2 \underbrace{2w_2}_{w_1} + (x_{1i} - x_{1j})^2 w_2}$$

$$\approx \sqrt{2(x_{1i} - x_{1j})^2 + (x_{2i} - x_{2j})^2}$$

- $x_1$  receives twice as weight as  $x_2$  in the distance calculation between  $X_i$  and  $X_j$ , e.g we penalize more for differences in  $x_1$
- Weighting rescales (upweights) relative differences on some covariates over other differences
- The weighted Mahalanobis distance matching is affine invariant. For example, dividing all weights by  $w_1$ , affine transformation of wmd will not change the matched stratas
- How do we choose the weights  $\tilde{w}$ ? Genetic optimization

### **Evolutionary Algorithm to Select Weights**

- The relevant task is to identify the k weights
- Pick weights to minimizes differences on  $X_p$  across treated and control units after matching This is our objective function
- Genetic Matching
  - Evolutionary algorithm to optimize weights over multiple generations
  - ullet Optimize means reduces imbalance on included X covariates

# Genetic Matching Workflow

Basic argument for the functions GenMatch and Match in the Matching R package

- Tr Treatment indicator
- $X_k$  is covariates for matching (Match Matrix)
- $X_p$  is covariates for balancing("Balance Matrix)
- Note:  $X_p$  can include **any**, **all** or **none** of  $X_k$

The optimization is performed using the package *rgenoud*. This optimization package can be used for other optimization problems, for example *maximum likelihood* 

### Genetic Matching Workflow

- At starting generation  $g_0$ :
  - 1. Select starting  $\tilde{w}_0$
  - 2. Devise *m* weight vectors using evolutionary string operators (called populations)
  - 3. For each weight vector:
    - a. Compute  $wmd(X_k, \tilde{w}_0)$
    - b. Match i treated to j control units using  $wmd(X_{ki}, X_{kj}, \tilde{w}_0)$  distances
    - c. Compute a balance metric  $\mathbb{B}(s)$ , a function of all s matched pairs
  - 4. Select the  $m^{th}$  weights  $(\tilde{w}_0^m)$  that maximizes  $\mathbb{B}(s)$
- At next generation  $g_1$  to  $g_n$ :
  - 5. Retain  $(\tilde{w}_{t-1}^m)$  as the best starting value from  $g_{t1}$
  - 6. Repeat steps 2 to 4 from above, maximizing  $\mathbb{B}_t(s)$  at each step
  - Stop by user control or once no more fitness gains over last few generations

Load the package 'Matching'

Matching includes the functions GenMatch(), Matching(), and MatchBalance().

Use GenMatch() to find the matches

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.func="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE, nboots=0,
ks=TRUE, verbose=FALSE, distance.tolerance=1e-05,
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

Use GenMatch() to find the matches GenMatch(Tr, X,

```
BalanceMatrix=X, estimand="ATT", M=1, weights=NULL,
pop.size = 100, max.generations=100, wait.generations=4,
hard.generation.limit=FALSE, starting.values=rep(1,ncol(X)),
fit.func="pvals", MemoryMatrix=TRUE, exact=NULL.
caliper=NULL, replace=TRUE, ties=TRUE,
CommonSupport=FALSE, nboots=0, ks=TRUE, verbose=FALSE,
distance.tolerance=1e-05, tolerance=sqrt(.Machine$double.eps),
min.weight=0, max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

**Tr** is the vector of treatment assignments.

Use GenMatch() to find the matches

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.func="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE, nboots=0,
ks=TRUE, verbose=FALSE, distance.tolerance=1e-05,
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1.
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

**X** is the vector of variables that we want to match on.

Use GenMatch() to find the matches

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.func="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE, nboots=0.
ks=TRUE, verbose=FALSE, distance.tolerance=1e-05.
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

**BalanceMatrix** is the vector of variables that we to achieve balance on (defaults to  $\mathbf{X}$ ).

"ATE", or "ATC".

Use GenMatch() to find the matches

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.func="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE, nboots=0.
ks=TRUE, verbose=FALSE, distance.tolerance=1e-05,
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
estimand is the parameter of interest. It can be the "ATT",
```

Use GenMatch() to find the matches

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M = 1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.func="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE, nboots=0,
ks=TRUE, verbose=FALSE, distance.tolerance=1e-05.
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

 ${\bf M}$  is the number of control units that you match to each treated unit.  ${\bf M}{=}1$  means one-to-one matching.

Use GenMatch() to find the matches

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.func="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE, nboots=0.
ks=TRUE, verbose=FALSE, distance.tolerance=1e-05,
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

**weights** is a vector with the same length as Y that specifies the weights to be placed on the observations. This is **not** the weights denoting the importance of the covariates.

Use GenMatch() to find the matches

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.fun="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE, nboots=0.
ks=TRUE, verbose=FALSE, distance.tolerance=1e-05,
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

**caliper** is a vector denoting the maximum difference apart each match can be for every covariate. Units with no suitable matches will be dropped.

Use GenMatch() to find the matches

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.fun="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE,
nboots=0, ks=TRUE, verbose=FALSE, distance.tolerance=1e-05,
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

**replace** allows you to choose whether to match with or without replacement.

Use GenMatch() to find the matches

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.fun="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE,
nboots=0, ks=TRUE, verbose=FALSE, distance.tolerance=1e-05,
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

**ties** allows you to deal with cases where there are several equally suitable control units for one treated unit. If **ties** is set at TRUE, GenMatch will average over the control units. If **ties** is FALSE, ties

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.fun="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE, nboots=0,
ks=TRUE, verbose=FALSE, distance.tolerance=1e-05,
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss=1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

paired allows you to choose whether GenMatch uses pairedt-tests.

```
GenMatch(Tr, X, BalanceMatrix=X, estimand="ATT", M=1,
weights=NULL, pop.size = 100, max.generations=100,
wait.generations=4, hard.generation.limit=FALSE,
starting.values=rep(1,ncol(X)), fit.fun="pvals",
MemoryMatrix=TRUE, exact=NULL, caliper=NULL,
replace=TRUE, ties=TRUE, CommonSupport=FALSE, nboots=0,
ks=TRUE, verbose=FALSE, distance.tolerance=1e-05.
tolerance=sqrt(.Machine$double.eps), min.weight=0,
max.weight=1000, Domains=NULL, print.level=2,
project.path=NULL, paired=TRUE, loss = 1,
data.type.integer=FALSE, restrict=NULL, cluster=FALSE,
balance=TRUE, ...)
```

**loss** is the loss function. It should take the vector of p-values for the covariates and return some value that GenMatch will try to maximize. The default is to sort the p-values and minimize the maximum discrepancy.

**Y** is the vector of outcomes.

```
\label{eq:match} \begin{split} &\text{Match}(\boldsymbol{Y}{=}\text{NULL}, \, \text{Tr}, \, X, \, Z = X, \, V = \text{rep}(1, \, \text{length}(Y)), \, \text{estimand} \\ &= \text{"ATT"}, \, M = 1, \, \text{BiasAdjust} = \text{FALSE}, \, \text{exact} = \text{NULL}, \, \text{caliper} = \\ &\text{NULL}, \, \text{replace}{=}\text{TRUE}, \, \text{ties}{=}\text{TRUE}, \\ &\text{CommonSupport}{=}\text{FALSE}, \text{Weight} = 1, \, \text{Weight.matrix} = \text{NULL}, \\ &\text{weights} = \text{NULL}, \, \text{Var.calc} = 0, \, \text{sample} = \text{FALSE}, \, \text{restrict}{=}\text{NULL}, \\ &\text{match.out} = \text{NULL}, \, \text{distance.tolerance} = 1\text{e-}05, \\ &\text{tolerance}{=}\text{sqrt}(.\text{Machine} \text{\$double.eps}), \, \text{version}{=}\text{"standard"}) \end{split}
```

```
\label{eq:match} \begin{split} &\text{Match}(Y {=} \text{NULL}, \ \textbf{Tr}, \ X, \ Z = X, \ V = \text{rep}(1, \text{length}(Y)), \text{ estimand} \\ &= \text{"ATT"}, \ M = 1, \ \text{BiasAdjust} = \text{FALSE}, \text{ exact} = \text{NULL}, \text{ caliper} = \\ &\text{NULL}, \ \text{replace} {=} \text{TRUE}, \ \text{ties} {=} \text{TRUE}, \\ &\text{CommonSupport} {=} \text{FALSE}, \text{Weight} = 1, \ \text{Weight.matrix} = \text{NULL}, \\ &\text{weights} = \text{NULL}, \ \text{Var.calc} = 0, \ \text{sample} = \text{FALSE}, \ \text{restrict} {=} \text{NULL}, \\ &\text{match.out} = \text{NULL}, \ \text{distance.tolerance} = 1e{-}05, \\ &\text{tolerance} {=} \text{sqrt}(.\text{Machine} \text{\$double.eps}), \ \text{version} {=} \text{"standard"}) \end{split}
```

**Tr** is the treatment assignment vector.

```
\label{eq:match} \begin{split} &\text{Match}(Y=\text{NULL}, \, \text{Tr}, \, \textbf{X}, \, Z=X, \, V=\text{rep}(1, \, \text{length}(Y)), \, \text{estimand} \\ &=\text{"ATT"}, \, M=1, \, \text{BiasAdjust}=\text{FALSE}, \, \text{exact}=\text{NULL}, \, \text{caliper}=\text{NULL}, \, \text{replace}=\text{TRUE}, \, \text{ties}=\text{TRUE}, \\ &\text{CommonSupport}=\text{FALSE}, \text{Weight}=1, \, \text{Weight.matrix}=\text{NULL}, \, \text{weights}=\text{NULL}, \, \text{Var.calc}=0, \, \text{sample}=\text{FALSE}, \, \text{restrict}=\text{NULL}, \, \text{match.out}=\text{NULL}, \, \text{distance.tolerance}=1\text{e-}05, \, \text{tolerance}=\text{sqrt}(.\text{Machine}\text{\$double.eps}), \, \text{version}=\text{"standard"}) \end{split}
```

**X** is the matrix of control variables.

Match(Y=NULL, Tr, X, Z = X, V = rep(1, length(Y)), estimand = "ATT", M = 1, BiasAdjust = FALSE, exact = NULL, caliper = NULL, replace=TRUE, ties=TRUE, CommonSupport=FALSE, Weight = 1, Weight.matrix = NULL, weights = NULL, Var.calc = 0, sample = FALSE, restrict=NULL, match.out = NULL, distance.tolerance = 1e-05, tolerance=sqrt(.Machine\$double.eps), version="standard")

**Weight** can be set at 2 to do Mahalanobis Distance matching. Leave this argument blank if you want to do Genetic matching.

```
Match(Y=NULL, Tr, X, Z = X, V = rep(1, length(Y)), estimand = "ATT", M = 1, BiasAdjust = FALSE, exact = NULL, caliper = NULL, replace=TRUE, ties=TRUE, CommonSupport=FALSE,Weight = 1, Weight.matrix = NULL, weights = NULL, Var.calc = 0, sample = FALSE, restrict=NULL, match.out = NULL, distance.tolerance = 1e-05, tolerance=sqrt(.Machine$double.eps), version="standard")
```

**Weight.matrix** takes the output from the GenMatch() function.

```
$est
      [,1]
[1,] 20933
$se
[1] 28768.51
$est.noadj
[1] 20933
$se.standard
[1] 23754.81
```

\$est is the estimated treatment effect (adjusted for bias if you BiasAdjust=TRUE)

\$se is the estimated standard error (adjusted for the uncertainty in the matching procedure)

\$est.noadj is the estimated treatment effect (not adjusted for bias)

\$se.standard is the standard error estimated the normal way

```
$index.treated
[1] 6 8 11 28 30 34 35 42 43 45 50 51 52 56 60

$index.control
[1] 48 3 48 48 17 36 40 46 25 10 48 64 15 64 19
```

```
MatchBalance(formul, data = NULL, match.out = NULL, ks = TRUE, nboots=500, weights=NULL, digits=5, paired=TRUE, print.level=1)
```

```
MatchBalance(formul, data = NULL, match.out = NULL, ks = TRUE, nboots=500, weights=NULL, digits=5, paired=TRUE, print.level=1)
```

**formul** is not the regular formula. It should be written as Treat  $\sim$  Control 1+ Control 2+...

```
MatchBalance(formul, data = NULL, match.out = NULL, ks = TRUE, nboots=500, weights=NULL, digits=5, paired=TRUE, print.level=1)
```

data is the data frame.

MatchBalance(formul, data = NULL, match.out = NULL, ks = TRUE, nboots=500, weights=NULL, digits=5, paired=TRUE, print.level=1)

match.out is the output of Match(). If you do not include this, MatchBalance will only return the balance before matching.

```
MatchBalance(formul, data = NULL, match.out = NULL, ks = TRUE, nboots=500, weights=NULL, digits=5, paired=TRUE, print.level=1)
```

paired determines if the t.tests on the matched data is paired.

```
> mat=Match(Y=data$b2004, Tr=data$etouch, X=with(data,cbind(income, b2000, hispanic, size)), Weight.matrix=gen)
> 
> MatchBalance(etouch ~ income + b2000 + hispanic + size, data=data, match.out=mat)
```

***** (V1) income *****			
	Before Matching	After	Matching
mean treatment	39282	39282	
mean control	34261	39358	
std mean diff	105.39	-1.5994	
mean raw eQQ diff	5574.9	1128.3	
med raw eQQ diff	6222	1045	
max raw eQQ diff	7100	4158	
mean eCDF diff	0.26883	0.066667	
med eCDF diff	0.25513	0.066667	
max eCDF diff	0.52949	0.2	
var ratio (Tr/Co)	0.56619	1.0204	
T-test p-value	0.0023728	0.88572	
KS Bootstrap p-value	< 2.22e-16	0.85	
KS Naive p-value	0.0016245	0.92509	
KS Statistic	0.52949	0.2	

- Your custom loss function should take a vector containing all t-tests p-values, followed by all the ks-test p-values, for the covariates in the BalanceMatrix.
- Therefore, the length of the vector it will receive is twice the length of the number of covariates in the BalanceMatrix
- Your loss function should return either a number or a vector
- If it returns a number, then GenMatch will try to maximize that number
- If it returns a vector, then GenMatch will try to maximize the first number in that vector. In cases where multiple matching schemes result in the same first number, then GenMatch will look at the second number, and then the third number, and so on

#### What does this loss function do?

```
loss.function=function(x) {
  p.vals = x
  return(sort(p.vals))
}
```

Answer: Maximize the minimum p-value, with ties are broken by looking at the second lowest p-value. This is the default loss function.

#### What does this loss function do?

```
loss.function=function(x) {
  p.vals = x
  return(min(p.vals))
}
```

Answer: Maximize the minimum p-value. This is the loss function if you set loss=2

#### What does this loss function do?

```
loss.function=function(x) {
  p.vals = x
  return(mean(p.vals))
}
```

Answer: Maximize the mean of the p-values

#### What does this loss function do?

```
loss.function=function(x) {
  p.vals = x
  p.vals[1]=p.vals[1]*0.5
  p.vals[length(p.vals)/
2+1]=p.vals[length(p.vals)/2+1]*0.5
  return(sort(p.vals))
}
```

Answer: Maximize the minimum p-value, giving more weight to the first covariate (we will treat the p-values for the first covariate as half their actual value

#### What does this loss function do?

```
loss.function = function(x) {
  p.vals = x
  if(sum(x < initial) > 0) {
    p.vals = 0.1*p.vals
  }
  return(sort(p.vals))
}
```

Answer: Restricts the search to only the weighting schemes where balance on every covariate is better than before matching. (Note: If GenMatch cannot find any such balancing scheme, the loss function defaults to maximize the minimum balance