Package 'RSurveillance'

October 6, 2014

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adj.risk calculate adjusted risk

Description

calculate adjusted risk

Usage

adj.risk(rr, ppr)

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Arguments

rr relative risk values (vector)

ppr population proportions corresponding to rr values (vector of equal length)

Value

vector of adjusted risk values (in order corresponding to rr)

Examples

```
# examples for adj.risk
adj.risk(c(5, 1), c(0.1, 0.9))
adj.risk(c(5, 3, 1), c(0.1, 0.1, 0.8))
```

ар

Apparent prevalence

Description

Estimates apparent prevalence and confidence limits for given sample size and result

Usage

```
ap(x, n, type = "wilson", conf = 0.95)
```

Arguments

x number of positives in sample

n sample size, note: either x or n can be a vector, but at least one must be scalar

type method for estimating CI, one of c("normal", "exact", "wilson", "jeffreys", "agresticoull", "all"), default = "wilson"

conf level of confidence required, default = 0.95

Value

either 1) if type = "all", a list with 5 elements, each element a matrix with 6 columns, x, n, proportion, lower confidence limit, upper confidence limit, confidence level and CI method; or 2) a matrix of results for the chosen method

```
# examples for ap function
n<- 200
x<- 25
conf<- 0.95
ap(x, n)
ap(seq(10, 100, 10), 200, type = "agresti")
ap(seq(10, 100, 10), 200, type = "all")</pre>
```

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binom.a	gresti
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Agresti-Coull confidence limits

Description

Calculates Agresti-Coull confidence limits for a simple proportion (apparent prevalence)

Usage

```
binom.agresti(x, n, conf = 0.95)
```

Arguments

x number of positives in sample

n sample size, note: either x or n can be a vector, but at least one must be scalar

conf level of confidence required, default 0.95

Value

a dataframe with 6 columns, x, n, proportion, lower confidence limit, upper confidence limit, confidence level and CI method

Examples

```
# test binom.agresti
binom.agresti(25, 200)
binom.agresti(seq(10, 100, 10), 200)
binom.agresti(50, seq(100, 1000, 100))
```

binom.cp

Clopper-Pearson exact confidence limits

Description

Calculates Clopper-Pearson exact binomial confidence limits for a simple proportion (apparent prevalence)

Usage

```
binom.cp(x, n, conf = 0.95)
```

Arguments

x number of positives in sample

n sample size, note: either x or n can be a vector, but at least one must be scalar

conf level of confidence required, default = 0.95

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Value

a dataframe with 6 columns, x, n, proportion, lower confidence limit, upper confidence limit, confidence level and CI method

Examples

```
# test binom.cp
binom.cp(25, 200)
binom.cp(seq(10, 100, 10), 200)
binom.cp(50, seq(100, 1000, 100))
```

binom.jeffreys

Jeffreys confidence limits

Description

Calculates Jeffreys confidence limits for a simple proportion (apparent prevalence)

Usage

```
binom.jeffreys(x, n, conf = 0.95)
```

Arguments

n number of positives in sample
 n sample size, note: either x or n can be a vector, but at least one must be scalar
 conf level of confidence required, default = 0.95

Value

a dataframe with 6 columns, x, n, proportion, lower confidence limit, upper confidence limit, confidence level and CI method

```
# test binom.jeffreys
binom.jeffreys(25, 200)
binom.jeffreys(seq(10, 100, 10), 200)
binom.jeffreys(50, seq(100, 1000, 100))
```

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disc.prior

calculate discounted prior

Description

calculate discounted prior

Usage

```
disc.prior(prior, p.intro)
```

Arguments

prior prior probability of freedom before surveillance

p.intro probability of introduction for time period (scalar or vactor equal length to sep)

Value

vector of discounted prior probability of freedom

Examples

```
# examples for disc.prior
disc.prior(0.5, 0.01)
disc.prior(0.95, c(0.001, 0.005, 0.01, 0.02, 0.05))
disc.prior(c(0.5, 0.6, 0.7, 0.8, 0.9, 0.95), 0.01)
```

epi.calc

calculate effective probability of infection

Description

calculate effective probability of infection

Usage

```
epi.calc(pstar, rr, ppr)
```

Arguments

pstar design prevalence

rr relative risk values (vector)

ppr population proportions corresponding to rr values (vector of equal length)

Value

list of 2 elements, a vector of EPI values and a vector of corresponding adjusted risks (in corresponding order to rr)

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Examples

```
# examples for epi.calc
epi.calc(0.1, c(5, 1), c(0.1, 0.9))
epi.calc(0.02, c(5, 3, 1), c(0.1, 0.1, 0.8))
```

n.2stage

calculate 2-stage sample size

Description

calculate 2-stage sample size

Usage

```
n.2stage(H = NA, N = NA, sep.sys = 0.95, sep.c, pstar.c, pstar.u, se = 1)
```

Arguments

Н	population size = number of clusters or NA if not known
N	populaton size if known, scalar or vector
sep.sys	desired population sensitivity
sep.c	desired cluster-level sensitivity
pstar.c	specified cluster-level design prevalence (proportion or integer)
pstar.u	specified population-level design prevalence (proportion or integer)
se	unit sensitivity

Value

list of number of clusters to sample and sample size per cluster

```
# examples of n.2stage - checked
n.2stage(NA, NA, 0.95, 0.5, 0.01, 0.1, 0.95)
n.2stage(500, NA, 0.95, 0.5, 10, 0.1, 0.95)
n.2stage(1000, c(50, 100, 200, 500, 1000, 5000, NA), 0.95, 0.5, 0.01, 0.05, 0.8)
n.2stage(1000, c(50, 100, 200, 500, 1000, 5000, NA), 0.95, 0.5, 0.01, 1, 0.8)
n.2stage(1000, c(50, 100, 200, 500, 1000, 5000, NA), 0.9, 0.95, 1, 0.1, 0.8)
```

n.binom

n.ap

Sample size for apparent prevalence

Description

Calculates sample size for estimating apparent prevalence (simple proportion)

Usage

```
n.ap(p, precision, conf = 0.95)
```

Arguments

p expected proportion, scalar or vector of values

precision absolute precision, +/- proportion equivalent to half the width of the desired con-

fidence interval, scalar or vector of values, note: at least one of p and precision

must be a scalar

conf level of confidence required, default = 0.95

Value

a vector of sample sizes

Examples

```
# examples of n.ap
n.ap(0.5, 0.1)
n.ap(0.5, 0.1, conf=0.99)
n.ap(seq(0.1, 0.5, by = 0.1), 0.05)
n.ap(0.2, c(0.01, 0.02, 0.05, 0.1))
```

 ${\sf n.binom}$

calculate sample size assuming sampling with replacement (binomial)

Description

calculate sample size assuming sampling with replacement (binomial)

Usage

```
n.binom(sep, pstar, se = 1)
```

Arguments

sep desired population sensitivity pstar specified design prevalence

se unit sensitivity

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Value

vector of sample sizes

Examples

```
# examples for n.binom - checked
n.binom(sep=0.95, pstar=c(0.01, 0.02, 0.05, 0.1, 0.2))
n.binom(c(0.5, 0.8, 0.9, 0.95), 0.01)
```

n.c.freecalc

Freecalc optimum sample size and specified cut-point number of reac-

Description

Calculates optimum sample size and cut-point reactors to achieve specified population sensitivity, for given population size and other parameters, using freecalc algorithm, all paramaters must be scalars

Usage

```
n.c.freecalc(N, sep = 0.95, c = 1, se, sp = 1, pstar, minSpH = 0.95)
```

Arguments

N	population size
sep	target population sensitivity
С	The cut-point number of reactors to classify a herd/flock as positive, default=1, if reactors $<$ c result is negative, $>=$ c is positive
se	test unit sensitivity
sp	test unit specificity, default=1
pstar	design prevalence
minSpH	minimium desired population specificity

Value

a list of 3 elements, a dataframe with 1 row and six columns for the recommended sample size and corresponding values for population sensitivity (SeP), population specificity (SpP), N, c and pstar, a vector of SeP values and a vector of SpP values, for n = 1:N

```
# examples for n.c.hp
n.c.freecalc(120,0.95,c=5,se=0.9,sp=0.99,pstar=0.1, minSpH=0.9)[[1]]
n.c.freecalc(65,0.95,c=5,se=0.95,sp=0.99,pstar=0.05, minSpH=0.9)
```

n.freecalc

n.c.hp	Hypergeometric (HerdPlus) optimum sample size and specified cut- point number of reactors
	point number of reactors

Description

Calculates optimum sample size and cut-point reactors to achieve specified population sensitivity, for given population size and other parameters, all paramaters must be scalars

Usage

```
n.c.hp(N, sep = 0.95, c = 1, se, sp = 1, pstar, minSpH = 0.95)
```

Arguments

N	population size
sep	target population sensitivity
С	The cut-point number of reactors to classify a herd/flock as positive, default=1, if reactors < c result is negative, >= c is positive
se	test unit sensitivity
sp	test unit specificity, default=1
pstar	design prevalence
minSpH	minimium desired population specificity

Value

a list of 3 elements, a dataframe with 1 row and six columns for the recommended sample size and corresponding values for population sensitivity (SeP), population specificity (SpP), N, c and pstar, a vector of SeP values and a vector of SpP values, for n = 1:N

Examples

```
# examples for n.c.hp
n.c.hp(65,0.95,c=5,se=0.95,sp=0.99,pstar=0.05, minSpH=0.9)[[1]]
tmp<- n.c.hp(120,0.95,c=5,se=0.9,sp=0.99,pstar=0.1, minSpH=0.9)</pre>
```

n.freecalc	Freecalc sample size for a finite population and specified cut-point
	number of reactors

Description

Calculates sample size required for a specified population sensitivity, for a given population size, cut-point number of reactors and other parameters, using Freecalc algorithm. All paramaters must be scalars

Usage

```
n.freecalc(N, sep = 0.95, c = 1, se, sp = 1, pstar, minSpH = 0.95)
```

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Arguments

N	population size
sep	target population sensitivity
С	The cut-point number of reactors to classify a herd/flock as positive, default=1, if reactors $<$ c result is negative, $>=$ c is positive
se	test unit sensitivity
sp	test unit specificity, default=1
pstar	design prevalence
minSpH	minimium desired population specificity

Value

a list of 2 elements, a dataframe with 1 row and six columns for the recommended sample size and corresponding values for population sensitivity (SeP), population specificity (SpP), N, c and pstar and a dataframe of n rows with SeP and SpP values for each value of n up to the recommended value

Examples

```
# examples for n.freecalc
n.freecalc(65,0.95,c=1,se=0.95,sp=0.99,pstar=0.05, minSpH=0.9)[[1]]
n.freecalc(65,0.95,c=2,se=0.95,sp=0.99,pstar=0.05, minSpH=0.9)[[1]]
n.freecalc(65,0.95,c=3,se=0.95,sp=0.99,pstar=0.05, minSpH=0.9)
```

n.freedom

calculate sample size usng appropriate method for data provided

Description

calculate sample size usng appropriate method for data provided

Usage

```
n.freedom(N = NA, sep = 0.95, pstar, se = 1)
```

Arguments

N	populaton size if known, scalar or vector
sep	desired population sensitivity
pstar	specified design prevalence (proportion or integer)
se	unit sensitivity

Value

vector of sample sizes, NA if N is specified and n>N

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Examples

```
# examples for n.freedom - checked
n.freedom(NA, sep=0.95, pstar=0.01, se=1)
n.freedom(500, sep=0.95, pstar=0.01, se=1)
n.freedom(N=c(100, 500, 1000, 5000, 10000, 100000, NA), sep=0.95, pstar=0.01, se=1)
n.freedom(500, sep=0.95, pstar=0.01, se=c(0.5, 0.6, 0.7, 0.8, 0.9, 0.99, 1))
```

n.hp

Hypergeometric (HerdPlus) sample size for finite population and specified cut-point number of reactors

Description

Calculates sample size to achieve specified population sensitivity with population specificity >= specified minimum value, for given population size, cut-point number of reactors and other parameters, all paramaters must be scalars

Usage

```
n.hp(N, sep = 0.95, c = 1, se, sp = 1, pstar, minSpH = 0.95)
```

Arguments

N	population size
sep	target population sensitivity
С	The cut-point number of reactors to classify a herd/flock as positive, default=1, if reactors $<$ c result is negative, $>=$ c is positive
se	test unit sensitivity
sp	test unit specificity, default=1
pstar	design prevalence
minSpH	minimium desired population specificity

Value

A list of 2 elements, a dataframe with 1 row and six columns for the recommended sample size and corresponding values for population sensitivity (SeP), population specificity (SpP), N, c and pstar and a dataframe of n rows with SeP and SpP values for each value of n up to the recommended value. Returns sample size for maximum achievable sep if it is not possible to achieve target sep AND SpP>= minSpH.

```
# examples for n.hp
n.hp(65,0.95,c=1,se=0.95,sp=0.99,pstar=0.05, minSpH=0.9)[[1]]
n.hp(65,0.95,c=2,se=0.95,sp=0.99,pstar=0.05, minSpH=0.9)
```

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n.hypergeo	calculate sample size assuming sampling without replacement (hypergeometric)
n.hypergeo	

Description

calculate sample size assuming sampling without replacement (hypergeometric)

Usage

```
n.hypergeo(sep, N, d, se = 1)
```

Arguments

sep	desired population sensitivity
N	population size
d	expected number of infected units in population (=pstar*N rounded to next integer)
se	unit sensitivity

Value

vector of sample sizes, NA if n>N

Examples

```
# examples for n.hypergeo - checked
n.hypergeo(0.95, N=100, d=1, se = 0.95)
n.hypergeo(sep=0.95, N=c(100, 200, 500, 1000, 10000), d=ceiling(0.01*c(100, 200, 500, 1000, 10000)))
n.hypergeo(c(0.5, 0.8, 0.9, 0.95), N=100, d=5)
n.hypergeo(0.95, N=80, d=c(1, 2, 5, 10))
n.hypergeo(0.95, N=80, d=c(1, 2, 5, 10), se = 0.8)
```

n.pfree

calculate sample size rqued to achieve target confidence of freedom

Description

calculate sample size rqued to achieve target confidence of freedom

Usage

```
n.pfree(pfree, prior, p.intro, pstar, se, N = NA)
```

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Arguments

pfree desired probability of freedom (scalar or vector) prior prior probability of freedom before surveillance probability of introduction for time period (scalar or vactor equal length to sep) p.intro

pstar design prevalence

unit sensitivity (scalar or vector) se

Ν population size

Value

vector of sample sizes

Examples

```
# examples for n.pfree
n.pfree(0.95, 0.5, 0.01, 0.05, 0.9)
n.pfree(0.95, 0.5, 0.01, 0.05, 0.9, N=300)
n.pfree(pfree = c(0.9, 0.95, 0.98, 0.99), prior = 0.7, 0.01, 0.01, 0.8, 1000)
n.pfree(0.95, 0.7, 0.01, 0.1, 0.96)
```

n.pooled

sample size for pooled testing for freedom

Description

sample size for pooled testing for freedom

Usage

```
n.pooled(sep, k, pstar, pse, psp = 1)
```

Arguments

desired population sensitivity (scalar or vector) sep

k pool size (constant) (scalar or vector)

pstar design prevalence pool-level sensitivity pse pool-level specificity psp

Value

vector of sample sizes

```
# examples for n.pooled
n.pooled(0.95, 5, 0.01, 1, 1)
n.pooled(0.95, 10, 0.1, 0.9, 1)
n.pooled(0.95, c(2, 5, 10, 20), 0.1, c(0.99, 0.98, 0.97, 0.95), 1)
```

n.rb

n.rb	calculate sample size for risk-based sampling \- binomial

Description

calculate sample size for risk-based sampling \- binomial

Usage

```
n.rb(pstar, rr, ppr, spr, se, sep)
```

Arguments

pstar	design prevalence (scalar)
rr	relative risk values (vector)
ppr	population proportions corresponding to rr values (vector of equal length)
spr	surveillance proportion for each risk group (vector equal length to rr, ppr)
se	unit sensitivity (fixed or vector same length as rr, ppr, n)
sep	required population sensitivity

Value

list of 2 elements, a vector of sample sizes for each risk group a scalar of total sample size, a vector of EPI values and a vector of adjusted risks

Examples

```
# examples for n.rb
n.rb(0.1, c(5, 3, 1), c(0.1, 0.10, 0.80), c(0.5, 0.3, 0.2), 0.9, 0.95)
n.rb(0.01, c(5, 1), c(0.1, 0.9), c(0.8, 0.2), c(0.9, 0.95), 0.95)
```

n.rb.varse	sample size for varying sensitivity	

Description

sample size for varying sensitivity

Usage

```
n.rb.varse(pstar, rr, ppr, spr, se, spr.rg, sep)
```

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Arguments

pstar	design prevalence
rr	vector of relative risk values
ppr	vector of population proportions for each risk group - same length as rr
spr	vector of surveillance proportions for each risk group - same length as rr
se	vector of sensitivity values
spr.rg	matrix of proportions of samples for each sensitivity value in each risk group (rows - risk groups, columns = sensitivity values) row sums must equal 1
sep	required population sensitivity

Value

list of 3 elements, a matrix of sample sizes for each risk and sensitivity group, a vector of EPI values and a vector of mean sensitivity for each risk group

Examples

```
# examples for n.rb.varse
m<- rbind(c(0.8, 0.2), c(0.5, 0.5), c(0.7, 0.3))
n.rb.varse(0.01, c(5, 3, 1), c(0.1, 0.1, 0.8), c(0.4, 0.4, 0.2), c(0.92, 0.8), m, 0.95)

m<- rbind(c(0.8, 0.2), c(0.6, 0.4))
n.rb.varse(0.05, c(3, 1), c(0.2, 0.8), c(0.7, 0.3), c(0.95, 0.8), m, 0.95)

m<- rbind(c(1), c(1))
n.rb.varse(0.05, c(3, 1), c(0.2, 0.8), c(0.7, 0.3), c(0.95), m, 0.99)</pre>
```

n.tp

Sample size for true prevalence

Description

Calculates sample size for estimating true prevalence using normal approximation

Usage

```
n.tp(p, se, sp, precision, conf = 0.95)
```

Arguments

p	estimated true prevalence
se	test sensitivity
sp	test specificity
precision	absolute precision, $+$ /- proportion equal to half the width of the desired confidence interval
conf	desired level of confidence for CI, default = 0.95

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Value

a vector of sample sizes

Examples

```
# examples for n.tp
n.tp(0.1, 0.9, 0.99, 0.05)
n.tp(0.1, 0.9, 0.99, 0.05, conf = 0.99)
n.tp(c(0.05, 0.1, 0.2, 0.3, 0.4, 0.5), 0.9, 0.99, 0.05)
n.tp(0.5, 0.9, 0.99, c(0.01, 0.02, 0.05, 0.1, 0.2))
```

pfree.1

calculate confidence of freedom for a single time period

Description

calculate confidence of freedom for a single time period

Usage

```
pfree.1(sep, p.intro, prior = 0.5)
```

Arguments

population sensitivity for time period (scalar or vector)

p.intro probability of introduction for time period (scalar or vactor equal length to sep)

prior prior probability of freedom before surveillance

Value

data. frame with columns for sep, p.intro, discounted prior, pfree, pfree.equ and prior.equ

```
# examples for pfree.1
pfree.1(0.8, 0.01, 0.5)
pfree.1(0.6, c(0.001, 0.005, 0.01, 0.02, 0.05), 0.5)
pfree.1(runif(10, 0.4, 0.6), 0.01, 0.5)
pfree.1(runif(10, 0.4, 0.6), runif(10, 0.005, 0.015), 0.5)
```

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pfree.calc	calculate probability (confidence) of freedom for given prior, sep and p.intro over 1 or more time periods

Description

calculate probability (confidence) of freedom for given prior, sep and p.intro over 1 or more time periods

Usage

```
pfree.calc(sep, p.intro, prior = 0.5)
```

Arguments

sep population sensitivity for time period (scalar or vector)

p.intro probability of introduction for time period (scalar or vactor equal length to sep)

prior prior probability of freedom before surveillance

Value

data. frame with columns for sep, p.intro, discounted prior, pfree, pfree.equ and prior.equ

Examples

```
# examples for pfree.calc
pfree.calc(0.8, 0.01, 0.5)
pfree.calc(rep(0.6,24), 0.01, 0.5)
pfree.calc(runif(10, 0.4, 0.6), 0.01, 0.5)
pfree.calc(runif(10, 0.4, 0.6), runif(10, 0.005, 0.015), 0.5)
```

pfree.equ

calculate equilibrium pfree and equilibrium prior Pfree

Description

calculate equilibrium pfree and equilibrium prior Pfree

Usage

```
pfree.equ(sep, p.intro)
```

Arguments

population sensitivity for time period (scalar or vector)p.introprobability of introduction for time period (scalar or vactor equal length to sep)

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Value

list of 2 vectors, equilibrium posterior probability of freedom and equilibrium prior (discounted) probability of freedom

Examples

```
# examples of pfree.equ
pfree.equ(runif(10, 0.4, 0.6), 0.01)
pfree.equ(0.8, 0.05)
pfree.equ(rep(0.9, 6), c(0.0001, 0.0005, 0.001, 0.005, 0.01, 0.05))
```

pstar.calc

calculate design prevalence for given sample size anddesired population-level sensitivity

Description

calculate design prevalence for given sample size anddesired population-level sensitivity

Usage

```
pstar.calc(N = NA, n, sep, se)
```

Arguments

N populaton size if known, scalar or vector

n sample size

sep desired population sensitivity

se unit sensitivity

Value

vector of design prevalence values

```
# examples of pstar.calc- checked
pstar.calc(NA, 280, 0.95, 0.98)
pstar.calc(500, 250, sep=0.95, se=1)
pstar.calc(N=c(100, 500, 1000, 5000, 10000, 100000, NA), n=30, sep=0.95, se=1)
pstar.calc(500, n=30, sep=0.95, se=c(0.5, 0.6, 0.7, 0.8, 0.9, 0.99, 1))
```

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sd.tp

Standard deviation of true prevalence estimate

Description

Calculates the standard deviation of true prevalence estimate assuming se and sp known exactly, used to calculate normal approximation CI for estimate

Usage

```
sd.tp(x, n, se, sp)
```

Arguments

X	number of positive results in sample
n	sample size
se	test sensitivity

sp test specificity

Value

vector of standard deviation values for true prevalence estimates

Examples

```
# example of sd.tp
sd.tp(1:10, 20, 0.9, 0.99)
```

se.parallel

calculate combined sensitivity for multiple tests in parallel (assuming independence)

Description

calculate combined sensitivity for multiple tests in parallel (assuming independence)

Usage

```
se.parallel(se)
```

Arguments

se

vector of unit sensitivity values

Value

scalar of combined sensitivity, assuming independence

```
# examples for se.parallel
se.parallel(c(0.99, 0.95, 0.8))
```

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se.series calculate combined sensitivity for multiple tests in series (assuming independence)	se.series	
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Description

calculate combined sensitivity for multiple tests in series (assuming independence)

Usage

```
se.series(se)
```

Arguments

se

vector of unit sensitivity values

Value

scalar of combined sensitivity, assuming independence

Examples

```
# examples for se.series
se.series(c(0.99, 0.95, 0.8))
```

sep

Population sensitivity function to calculate seh using most appropriate option depending on whether or not N provided

Description

Population sensitivity function to calculate seh using most appropriate option depending on whether or not N provided

Usage

```
sep(N = NA, n, pstar, se = 1)
```

Arguments

N vector of population sizes: NA or vector of same length as n

n vector of sample sizes

pstar design prevalence: single value as a proportion or integer se unit sensitivity: single value or vector same lengthh as n

Value

value vector of population-level sensitivities

sep.binom

Examples

```
# examples for sep - checked
sep(n=300, pstar=0.01, se=1)
sep(NA, 300, 0.01, 1)
sep(10000, 150, 0.02, 1)
sep(n=1:100, pstar = 0.05, se=0.95)
N<- seq(30, 100, by = 5)
se<- 0.95
pstar<- 0.1
n<- rep(30, length(N))
sep(N, n, pstar, se = se)
sep(rep(100, 10), seq(10, 100, by = 10), pstar = 1, se=0.99)
N<- c(55, 134, NA, 44, 256)
n<- c(15, 30, 28, 15, 33)
sep(N, n, 0.1, 0.95)</pre>
```

sep.binom

Population sensitivity assuming sampling with replacement (binomial)

Description

Population sensitivity assuming sampling with replacement (binomial)

Usage

```
sep.binom(n, pstar, se = 1, sp = 1)
```

Arguments

n	integer scalar or vector of number tested (sample size)
pstar	scalar or vector of design prevalence as proportion
se	unit sensitivity of test (proportion)
sp	unit specificity of test (proportion)

Value

vector of population-level sensitivities

```
# examples for sep.binom - checked
sep.binom(n=300, pstar = 0.02, se = 0.92)
tested<- seq(10,100, by=10)
prev<- 0.05
sens<- 0.9
sep.binom(tested, prev, sens)</pre>
```

sep.binom.imperfect 23

sep.binom.imperfect

Binomial population sensitivity for imperfect test

Description

Calculates population sensitivity for a large or unknown population and allowing for imperfect test sensitivity and specificity, using Binomial distribution

Usage

```
sep.binom.imperfect(n, c = 1, se, sp = 1, pstar)
```

Arguments

n	sample size
С	The cut-point number of reactors to classify a herd/flock as positive, default=1, if reactors $<$ c result is negative, $>=$ c is positive
se	test unit sensitivity
sp	test unit specificity, default=1
pstar	design prevalence

Value

a vector of population-level sensitivities

Examples

```
# examples for sep.imperfect.binom
sep.binom.imperfect(1:10*5, 2, 0.95, 0.98, 0.1)
sep.binom.imperfect(50, 1:5, 0.95, 0.98, 0.1)
sep.binom.imperfect(30, 2, 0.9, 0.98, 0.1)
sep.binom.imperfect(30, 1, 0.9, 0.98, 0.1)
```

sep.exact

Population sensitivity assuming census (all units tested)

Description

Population sensitivity assuming census (all units tested)

Usage

```
sep.exact(d = 1, se = 1)
```

Arguments

d	expected number of infected units in population (=pstar*N rounded to next in-
	teger)

se unit sensitivity of test (proportion)

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Value

vector of population-level sensitivities

Examples

```
# examples for sep.exact - checked
sep.exact(d=1, se = 0.92)
inf<- 1:5
sens<- 0.8
sep.exact(d=inf, se=sens)
sep.exact(se=0.8, d = ceiling(0.01*c(10, 50, 100, 250, 500)))</pre>
```

sep.freecalc

FreeCalc population sensitivity for imperfect test

Description

Calculates population sensitivity for a finite population and allowing for imperfect test sensitivity and specificity, using Freecalc method

Usage

```
sep.freecalc(N, n, c = 1, se, sp = 1, pstar)
```

Arguments

N	population size
n	sample size
С	The cut-point number of reactors to classify a herd/flock as positive, default=1, if reactors < c result is negative, >= c is positive
se	test unit sensitivity
sp	test unit specificity, default=1
pstar	design prevalence - assumed or target prevalence for detection of disease in the population

Value

population-level sensitivity

```
# examples of sep.freecalc
sep.freecalc(150, 30, 2, 0.9, 0.98, 0.1)
sep.freecalc(150, 30, 1, 0.9, 0.98, 0.1)
```

sep.hp 25

Hypergeometric (HerdPlus) population sensitivity for imperfect test

Description

Calculates population sensitivity for a finite population and allowing for imperfect test sensitivity and specificity, using Hypergeometric distribution

Usage

```
sep.hp(N, n, c = 1, se, sp = 1, pstar)
```

Arguments

N	population size
n	sample size
С	The cut-point number of reactors to classify a herd/flock as positive, default=1, if reactors $<$ c result is negative, $>=$ c is positive
se	test unit sensitivity
sp	test unit specificity, default=1
pstar	design prevalence

Value

a vector of population-level sensitivities

Examples

```
# examples of sep.hp
sep.hp(150, 1:5*10, 2, 0.9, 0.98, 0.1)
sep.hp(150, 30, 2, 0.9, 0.98, 15)
sep.hp(150, 30, 1, 0.9, 0.98, 15)
sep.hp(150, 30, 1, 0.9, 0.98, 0.1)
```

sep.hypergeo

Population sensitivity assuming sampling without replacement (hypergeometric approximation)

Description

Population sensitivity assuming sampling without replacement (hypergeometric approximation)

Usage

```
sep.hypergeo(N, n, d, se = 1)
```

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Arguments

N	population size
n	sample size (tested)
d	expected number of infected units in population (=pstar*N rounded to next integer) $\frac{1}{2}$
se	unit sensitivity of test (proportion)

Value

value vector of population-level sensitivities

Examples

```
# examples for sep.hypergeo - checked
sep.hypergeo(N=100, n=50, d=1, se = 0.92)
inf<- 1:5
sens<- 0.8
sep.hypergeo(N=100, n=50, d=inf, se=sens)
N<- c(10, 50, 100, 250, 500)
sep.hypergeo(se=0.8, N=N, n=c(5, 25, 50, 125, 250), d = ceiling(0.01*N))</pre>
```

sep.pfree

calcuate population sensitivity for given PFree

Description

calcuate population sensitivity for given PFree

Usage

```
sep.pfree(prior, pfree)
```

Arguments

prior prior probability of freedom before surveillance (scalar or vector)

pfree desired probability of freedom (scalar or vector)

Value

vector of population-level sensitivities

```
# examples of sep.pfree
sep.pfree(0.5, 0.95)
sep.pfree(c(0.5, 0.6, 0.7, 0.8, 0.9, 0.95), 0.99)
sep.pfree(0.5, c(0.8, 0.9, 0.95, 0.99))
```

sep.pooled 27

sep.pooled

population sensitivity with pooled sampling

Description

population sensitivity with pooled sampling

Usage

```
sep.pooled(r, k, pstar, pse, psp = 1)
```

Arguments

r number of pools sampled (scalar or vector)
k pool size (constant) (scalar or vector)
pstar design prevalence
pse pool-level sensitivity
psp pool-level specificity

Value

list of 2 elements, vector of sep values and vector of spp values

Examples

```
# examples for sep.pooled
sep.pooled(60, 5, 0.01, 1, 1)
sep.pooled(4, 10, 0.1, 0.9, 1)
sep.pooled(1:10*5, 5, 0.02, 0.9, 0.99)
sep.pooled(10, 5, 0.05, c(0.8, 0.9, 0.95, 0.99), 1)
```

sep.prior

calcuate population sensitivity for given prior pfree

Description

calcuate population sensitivity for given prior pfree

Usage

```
sep.prior(prior, p.intro)
```

Arguments

```
prior prior probability of freedom before surveillance (scalar or vector)

p.intro probability of introduction for time period (scalar or vactor equal length to sep)
```

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Value

vector of population-level sensitivities

Examples

```
# examples of sep.prior
sep.prior(0.95, 0.01)
sep.prior(c(0.9, 0.95, 0.98, 0.99), 0.01)
sep.prior(0.95, c(0.001, 0.005, 0.01, 0.02, 0.05))
```

sep.rb.bin

calculate popuation sensitivity for single risk factor - binomial

Description

calculate popuation sensitivity for single risk factor - binomial

Usage

```
sep.rb.bin(pstar, rr, ppr, n, se)
```

Arguments

pstar	design prevalence (scalar)
rr	relative risk values (vector)
ppr	population proportions corresponding to rr values (vector of equal length)
n	sample size per risk category (vector same length as rr and ppr)
se	unit sensitivity (fixed or vector same length as rr, ppr, n)

Value

list of 3 elements, a scalar of population-level sensitivity a vector of EPI values and a vector of corresponding Adjusted risks

```
# examples for sep.rb.bin
sep.rb.bin(0.1, c(5, 3, 1), c(0.1, 0.1, 0.8), c(5, 5, 5), 0.9)
sep.rb.bin(0.1, c(5, 1), c(0.1, 0.9), c(10, 5), c(0.95, 0.9))
sep.rb.bin(0.1, c(5, 1), c(0.1, 0.9), c(10, 5), c(0.9, 0.9))
sep.rb.bin(0.01, c(5, 1), c(0.1, 0.9), c(90, 50), c(0.9, 0.9))
```

sep.rb.bin.varse

sep.rb.bin.varse

calculate population sensitivity for varying unit sensitivity - binomial

Description

calculate population sensitivity for varying unit sensitivity - binomial

Usage

```
sep.rb.bin.varse(pstar, rr, ppr, df)
```

Arguments

pstar	design prevalence (scalar)
rr	relative risk values (vector)
ppr	population proportions corresponding to rr values (vector of equal length)
df	dataframe of values for each sensitivity level col $1 = risk$ group index, col $2 = unit Se$, col $3 = n$

Value

list of 3 elements, a scalar of population-level sensitivity a vector of EPI values and a vector of corresponding Adjusted risks

```
# examples for sep.rb.bin.varse
rg<- c(1, 1, 2, 2)
se<- c(0.92, 0.85, 0.92, 0.85)
n<- c(80, 30, 20, 30)
df<- data.frame(rg, se, n)</pre>
sep.rb.bin.varse(0.01, c(5, 1), c(0.1, 0.9), df)
rg<- c(1, 1, 2, 2)
se<- c(0.95, 0.8, 0.95, 0.8)
n<- c(20, 10, 10, 5)
df<- data.frame(rg, se, n)</pre>
sep.rb.bin.varse(0.05, c(3, 1), c(0.2, 0.8), df)
rg < -c(rep(1, 30), rep(2, 15))
se<-c(rep(0.95, 20), rep(0.8, 10), rep(0.95, 10), rep(0.8, 5))
n < - rep(1, 45)
df<- data.frame(rg, se, n)</pre>
sep.rb.bin.varse(0.02, c(3, 1), c(0.2, 0.8), df)
rg<- c(1, 2, 3, 1, 2, 3)
se<- c(0.95, 0.95, 0.95, 0.8, 0.8, 0.8)
n<- c(20, 10, 10, 30, 5, 5)
df<- data.frame(rg, se, n)</pre>
sep.rb.bin.varse(0.01, c(5, 3, 1), c(0.1, 0.3, 0.6), df)
```

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sep.	rb.	hypergeo

calculate population sensitivity for single risk factor - hypergeometric

Description

calculate population sensitivity for single risk factor - hypergeometric

Usage

```
sep.rb.hypergeo(pstar, rr, N, n, se)
```

Arguments

pstar	design prevalence (scalar)
rr	relative risk values (vector)
N	Population size per risk category (vector same length as rr and ppr)
n	sample size per risk category (vector same length as rr and ppr)
se	unit sensitivity (fixed or vector same length as rr, ppr, n)

Value

list of 3 elements, a scalar of population-level sensitivity a vector of EPI values and a vector of corresponding Adjusted risks

Examples

```
# examples for sep.rb.bin sep.rb.hypergeo(0.1, c(5, 3, 1), c(10, 10, 80), c(5, 5, 5), 0.9) sep.rb.hypergeo(0.1, c(5, 1), c(15, 140), c(10, 5), c(0.95, 0.9)) sep.rb.hypergeo(0.1, c(5, 1), c(23, 180), c(10, 5), c(0.9, 0.9)) sep.rb.hypergeo(0.01, c(5, 1), c(100, 900), c(90, 50), c(0.9, 0.9))
```

sep.rb.hypergeo.varse calculate population sensitivity for varying unit sensitivity - hypergeometric

Description

calculate population sensitivity for varying unit sensitivity - hypergeometric

Usage

```
sep.rb.hypergeo.varse(pstar, rr, N, df)
```

Arguments

pstar	design prevalence (scalar)
rr	relative risk values (vector)
N	vector of population size corresponding to rr values (vector of equal length)
df	dataframe of values for each sensitivity level col $1 = risk$ group index, col $2 = unit Se$, col $3 = n$

sep.rb2.binom 31

Value

list of 5 elements, a scalar of population-level sensitivity a vector of EPI values, a vector of corresponding Adjusted risks a vector of sample sizes (n) per risk group and a vector of mean unit sensitivities per risk group

Examples

```
# examples for sep.rb.hypergeo.varse
rg<- c(1, 1, 2, 2)
se<- c(0.92, 0.85, 0.92, 0.85)
n<- c(80, 30, 20, 30)
df<- data.frame(rg, se, n)</pre>
sep.rb.hypergeo.varse(0.01, c(5, 1), c(200, 1800), df)
rg<- c(1, 1, 2, 2)
se<- c(0.95, 0.8, 0.95, 0.8)
n<- c(20, 10, 10, 5)
df<- data.frame(rg, se, n)</pre>
sep.rb.hypergeo.varse(0.05, c(3, 1), c(100, 400), df)
rg<- c(rep(1, 30), rep(2, 15))
se<- c(rep(0.95, 20), rep(0.8, 10), rep(0.95, 10), rep(0.8, 5))
n < - rep(1, 45)
df<- data.frame(rg, se, n)</pre>
sep.rb.hypergeo.varse(0.02, c(3, 1), c(100, 400), df)
rg<- c(1, 2, 3, 1, 2, 3)
se<- c(0.95, 0.95, 0.95, 0.8, 0.8, 0.8)
n<- c(20, 10, 10, 30, 5, 5)
df<- data.frame(rg, se, n)</pre>
sep.rb.hypergeo.varse(0.01, c(5, 3, 1), c(100, 300, 600), df)
```

sep.rb2.binom

1-stage risk-based sampling with 2 risk factors - binomial

Description

1-stage risk-based sampling with 2 risk factors - binomial

Usage

```
sep.rb2.binom(pstar, rr1, ppr1, rr2, ppr2, n, se)
```

Arguments

pstar	design prevalence
rr1	relative risks for first level risk factor
ppr1	population proportions for first level risk factor
rr2	relative risks for second level risk factor, matrix, rows = levels of rr1, cols = levels of rr2
ppr2	population proportions for second level risk factor, matrix, rows = levels of rr1, cols = levels of rr2

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n	matrix of number tested for each risk group (stratified
se	test unit sensitivity

Value

list of 4 elements, a scalar of population-level sensitivity a matrix of EPI values, a vector of corresponding Adjusted risks for the first risk factor and a matrix of adjusted risks for the second risk factor

Examples

```
# examples for sep.rb2.binom
pstar<- 0.01
rr1<- c(3, 1)
ppr1<- c(0.2, 0.8)
rr2<- rbind(c(4,1), c(4,1))
ppr2<- rbind(c(0.1, 0.9), c(0.3, 0.7))
se<- 0.8
n<- rbind(c(50, 20), c(20, 10))
sep.rb2.binom(pstar, rr1, ppr1, rr2, ppr2, n, se)</pre>
```

sep.rb2.hypergeo

1-stage risk-based sampling with 2 risk factors - hypergeometric

Description

1-stage risk-based sampling with 2 risk factors - hypergeometric

Usage

```
sep.rb2.hypergeo(pstar, rr1, rr2, N, n, se)
```

Arguments

pstar	design prevalence
rr1	relative risks for first level risk factor
rr2	relative risks for second level risk factor
N	matrix of population size for each risk group (stratified
n	matrix of number tested (sample size) for each risk group (stratified
se	test unit sensitivity

Value

list of 6 elements, a scalar of population-level sensitivity a matrix of EPI values, a vector of corresponding Adjusted risks for the first risk factor and a matrix of adjusted risks for the second risk factor, a vector of population proportions for the first risk factor and a matrix of population proportions for the second risk factor

sep.sys 33

Examples

```
# examples for sep.rb2.hypergeo
pstar<- 0.01
rr1<- c(3, 1)
rr2<- rbind(c(4,1), c(4,1))
N<- rbind(c(100, 500), c(300, 1000))
n<- rbind(c(50, 20), c(20, 10))
se<- 0.8
sep.rb2.hypergeo(pstar, rr1, rr2, N, n, se)</pre>
```

sep.sys

population sensitivity from sampling of individual clusters

Description

population sensitivity from sampling of individual clusters

Usage

```
sep.sys(H = NA, N = NA, n, pstar.c, pstar.u, se = 1)
```

Arguments

Н	population size = number of clusters or NA if not known	
N	population size within clusters NA if not provided, otherwise a vector of same length as \boldsymbol{n}	
n	sample size (tested)	
pstar.c	cluster (herd) level design prevalence single value either proportion or integer	
pstar.u	unit (animal) level design prevalence single value either proportion or integer	
se	unit sensitivity of test (proportion)	

Value

vector of population-level sensitivities

Note

if pstar.c is not a proportion N must be entered (and N>=n)

```
# examples for sep.sys - checked
H<- 500
N<- rep(1000, 150)
N[5]<- NA
n<- rep(30, 150)
pstar.u<- 0.1
pstar.c<- 0.01
se<- 0.98
sep.sys(H, N, n, pstar.c, pstar.u, se)
sep.sys(NA, N, n, 0.02, 0.05, 0.95)</pre>
```

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```
N<- round(runif(105)*900+100)
n<- round(runif(105)*30+10)
sse<- sep.sys(1000, N, n, 0.02, 0.05, 0.9)
data.frame(N, n, sse[[2]])</pre>
```

sep.var.se

population sensitivity for varying unit sensitivity

Description

population sensitivity for varying unit sensitivity

Usage

```
sep.var.se(N = NA, se, pstar)
```

Arguments

N population size (number of units/clusters. $N \ge length(se)$) or NA if unknown

se vector of unit sensitivity values pstar specified design prevalence

Value

value vector of population-level sensitivities

Examples

```
# examples of sep.var.se - checked
sens<- c(rep(0.9, 50), rep(0.95, 100))
sep.var.se(NA, sens, 0.01)
sep.var.se(se=sens, pstar=0.01)
sep.var.se(N=500, sens, 0.01)
sep.var.se(NA, runif(150, 0.95, 0.99), 0.02)
sep.var.se(500, runif(150, 0.95, 0.99), 0.02)</pre>
```

sph

Population specificity

Description

Population specificity

Usage

```
sph(n, sp)
```

Arguments

n integer scalar or vector of number tested sp unit specificity of test (proportion) sph.binom 35

Value

value vector of population-level specificities

Examples

```
# examples for sph - checked
sph(10, 0.9)
sph(c(10, 20, 50, 100), 0.99)
sph(100, c(0.999, 0.99, 0.98, 0.95, 0.9))
```

sph.binom

Binomial population specificity for imperfect test

Description

Calculates population specificity for a large or unknown population, using the Binomial distribution and adjusting for cut-point number of reactors

Usage

```
sph.binom(n, c = 1, sp)
```

Arguments

```
n sample size
c The cut-point number of reactors to classify a herd/flock as positive, default=1, if reactors < c result is negative, >= c is positive
sp test unit specificity
```

Value

a vector of population-level specificities

```
# examples for sph.imperfect.sp
sph.binom(30, 2, 0.98)
sph.binom(30, 1, 0.98)
sph.binom(1:5*10, 2, 0.98)
sph.binom(100, 1:5, 0.98)
sph.binom(100, 3, 95:100/100)
sph.binom(c(5, 10, 15, 20, 30, 50, 100, 200), 2, 0.98)
```

36 sse.combined

Description

Calculates population specificity for a finite population and imperfect test, using Hypergeometric distribution

Usage

```
sph.hp(N, n, c = 1, sp)
```

Arguments

N	population size
n	sample size
С	The cut-point number of reactors to classify a herd/flock as positive, default=1, if reactors < c result is negative, >= c is positive
sp	test unit specificity

Value

a vector of population-level specificities

Examples

```
# examples of sph.hp
sph.hp(150, 30, 2, 0.98)
sph.hp(150, 30, 1, 0.98)
sph.hp(150, 1:5*10, 2, 0.98)
sph.hp(500, 30, 2, 95:100/100)
```

sse.combined

update between components to account for lack of independence

Description

update between components to account for lack of independence

Usage

```
sse.combined(C = NA, pstar.c, rr, ppr, sep)
```

Arguments

С	NA or vector of population sizes (number of clusters) for each risk group	
pstar.c	cluster level design prevalence	
rr	cluster level relative risks	
ppr	cluster level population proportions (not required if C is specified)	
sep	list of sep values for clusters in each component and corresponding risk group. Each element is a dataframe, first column= clusterid, 2nd =rg.c, 3rd col = sep	

sse.rb.2stage 37

Value

list of 2 elements, a matrix (or vector if C not specified) of population-level (surveillance system) sensitivities (binomial and hypergeometric and adjusted vs unadjusted) and a matrix of adjusted and unadjusted component sensitivities for each component

Examples

```
# example for sse.combined (checked in excel combined components.xlsx)
C<- c(300, 1200)
pstar<- 0.01
rr<- c(3,1)
ppr<- c(0.2, 0.8)
comp1<- data.frame(id=1:100, rg=c(rep(1,50), rep(2,50)), cse=rep(0.5,100))
comp2<- data.frame(id=seq(2, 120, by=2), rg=c(rep(1,25), rep(2,35)), cse=runif(60, 0.5, 0.8))
comp3<- data.frame(id=seq(5, 120, by=5), rg=c(rep(1,10), rep(2,14)), cse=runif(24, 0.7, 1))
sep<- list(comp1, comp2, comp3)
sse.combined(C, pstar, rr, sep = sep)
sse.combined(C=NA, pstar, rr, ppr, sep = sep)</pre>
```

sse.rb.2stage

calculate system sensitivity for 2 stage risk-based sampling

Description

calculate system sensitivity for 2 stage risk-based sampling

Usage

```
sse.rb.2stage(C = NA, pstar.c, pstar.u, rr.c, ppr.c, rr.u, ppr.u, N = NA, n,
    rg, se)
```

Arguments

С	Population size (number of clusters), NA = unknown
pstar.c	cluster level design prevalence
pstar.u	unit level design prevalence
rr.c	cluster level relative risks (vector), NA if no cluster level risk factor if risk factor does not apply at either level use $rr = c(1,1)$
ppr.c	cluster level population proportions for risk categories (vector), NA if no cluster level risk factor
rr.u	unit level relative risks (vector), NA if no unit level risk factor if risk factor does not apply at either level use $rr = c(1,1)$
ppr.u	matrix, 1 row for each cluster, columns = unit level risk groups
N	cluster sizes NA or matrix of N for each risk group for each cluster, N=NA means cluster sizes not provided)
n	matrix, 1 row for each cluster, columns = unit level risk groups
rg	vector of cluster level risk group for each cluster
se	unit sensitivity for each cluster, scalar or vector of values for each cluster

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Value

list of 2 elements, a scalar of population-level (surveillance system) sensitivity and a vector of cluster-level sensitivities

Examples

```
# examples for sse.rb.2stage
pstar.c<- 0.02
pstar.u<- 0.1
rr.c<- c(5, 1)
ppr.c<- c(0.1, 0.9)
rr.u<- c(3, 1)
se<- 0.9
n < - cbind(rep(10, 50), rep(5, 50))
rg<- c(rep(1, 30), rep(2, 20))
ppr.u<- cbind(rep(0.2, 50), rep(0.8, 50))
N<- cbind(rep(30, 50), rep(120, 50))
C<- 500
sse.rb.2stage(C=NA, pstar.c, pstar.u, rr.c, ppr.c, rr.u, ppr.u, N=NA, n, rg, se)
sse.rb.2stage(C, pstar.c, pstar.u, rr.c, ppr.c, rr.u, ppr.u, N=NA, n, rg, se)
sse.rb.2stage(C=NA, pstar.c, pstar.u, rr.c, ppr.c, rr.u, ppr.u, N, n, rg, se)
sse.rb.2stage(C, pstar.c, pstar.u, rr.c, ppr.c, rr.u, ppr.u, N, n, rg, se)
```

tp

True prevalence

Description

Estimates true prevalence and confidence limits for given sample size and result, according to specified method

Usage

```
tp(x, n, se, sp, type = "blaker", conf = 0.95)
```

Arguments

X	number of positive units (scalar)
n	sample size (no. units sampled) (scalar)
se	test sensitivity (scalar)
sp	test specificity (scalar)
type	method for estimating CI, one of c("normal", "c-p", "sterne", "blaker", "wilson", "all")
conf	desired level of confidence for CI, default = 0.95 (scalar)

Value

list with 2 elements, a matrix of apparent prevalence and lower and upper confidence limits and a matrix of true prevalence and lower and upper confidence limits using the chosen method(s)

tp.normal 39

Examples

```
# examples for tp
x<- 20
n<- 120
se<- 0.9
sp<- 0.99
conf<- 0.95
tp(x, n, se, sp, "all")
tp(x, n, se, sp, "c-p")
tp(x, n, 0.95, 0.9, "c-p")</pre>
```

tp.normal

Normal approximation confidence limits for true prevalence

Description

Estimates true prevalence and confidence limits for estimates based on normal approximation

Usage

```
tp.normal(x, n, se, sp, conf = 0.95)
```

Arguments

Х	number of positive results in sample	
n	sample size	
se	test unit sensitivity	
sp	test unit specificity	
conf	desired level of confidence for CI, default = 0.95	

Value

list with 2 elements, a matrix of apparent prevalence and wilson lower and upper confidence limits and a matrix of true prevalence and normal approximation lower and upper confidence limits

```
# examples for tp.normal
tp.normal(25, 120, 0.9, 0.99)
tp.normal(seq(5, 25, by=5), 120, 0.9, 0.99)
```

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