

Rgof and R2sample: Testing and Benchmarking for the Univariate Goodness-of-Fit and Two-Sample Problems

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Abstract In a goodness-of-fit problem one attempts to see whether a data set might have been generated by some theoretical probability distribution, possibly with unknown parameters. In a (non-parametric) two-sample problem one wants to check whether two data sets could have come from the same unspecified distribution. Both are among the oldest and best studied problems in Statistics. While they seem quite different, there are a number of similarities between these problems, not the least of which is that many methods exist that have versions for both. The two packages discussed in this article bring together a large number of methods and many different scenarios, most of which do not yet have existing implementations. They also include routines that allow a developer of a new method to quickly compare its performance (aka power) to those included in the package for a large number of cases.

1 Introduction

Both the goodness-of-fit and the nonparametric two-sample problem have histories going back a century, with many contributions by some of the most eminent statisticians. In the goodness-of-fit problem we have a sample (x_1, \dots, x_n) drawn from a random variable X . We also have a probability distribution F , possibly with unknown parameters, and we wish to test $H_0 : X \sim F$. In the two-sample problem we also have a second sample (y_1, \dots, y_m) drawn from some distribution G , and here we want to test $H_0 : F = G$, that is we want to test whether the two data sets were generated by the same (unspecified) distribution.

The literature on both of these problems is vast and steadily growing. Detailed discussions can be found in [D'Agostini and Stephens \(1986\)](#), [Thas \(2010\)](#), [Raynor et al. \(2012\)](#). For an introduction to Statistics and hypothesis testing in general see [Casella and Berger \(2002\)](#) or [Bickel and Doksum \(2015\)](#).

Some tests such as the Kolmogorov-Smirnov test are already implemented for both problems in base R [R Core Team \(2021\)](#). Many others can be run through various existing packages, for example the Anderson-Darling goodness-of-fit test is available in the R package *ADGofTest* [Bellosta \(2011\)](#). There are a number of packages with tests that focus on a specific distribution, for example the *nortest* [Gross and Ligges \(2015\)](#) package has five tests for composite normality. There are also packages that allow the user to run several tests, for example the *twosamples* [Dowd \(2022\)](#), *dgof* [Taylor and Emerson \(2009\)](#), *EnvStats* [Millard and Kowarik \(2017\)](#) and *goftest* [Faraway et al. \(2007\)](#) packages.

However, there are no packages that bring together as many tests as *R2sample* and *Rgof*. Moreover, some methods do not currently have any R implementations, for example Zhang's test, Lehmann-Rosenblatt, the Wasserstein p1 test and almost all tests for discrete data.

Both packages have the following features:

- many methods are implemented for both continuous and discrete data.
- the methods are implemented using both *Rcpp* [Eddelbuettel et al. \(2024\)](#) and parallel programming.
- the packages include routines to run several test and then find a corrected p value for the combination of tests.

- some of the methods allow for data with weights.
- the routines allow for a random sample size, assumed to come from a Poisson distribution.
- in the two-sample problem some methods make use of large-sample formulas, therefore allowing for very large data sets.
- the routines can also use any other user-defined tests.
- the packages include routines to easily carry out power studies and draw power graphs.
- the packages include routines to easily compare the power of a new test to those included in the packages.
- in the two-sample package the user can provide a routine that generates new data from a model. This can be used as an alternative to the permutation method to find p values.

There are several reasons for including tests for discrete data. In the context of a computer program this means a finite (and usually fairly small) number of different values which then repeat many times.

- Tests for discrete data such as from Binomial or Poisson distributions are of interest in their own right.
- There are currently almost no implementations of either goodness-of-fit or two-sample methods for discrete data in *R*.
- It also makes it possible to apply the tests to very large continuous data sets via discretization. While a goodness-of-fit test for a continuous data set with (say) 100,000 observations can be done in a matter of a few minutes, for larger data sets the calculations will be quite time consuming. Data sets with many millions of observations are not uncommon today. Binning the data and then running the corresponding discrete tests however is quite fast.
- There are also situations where the underlying distribution is continuous but the data is collected in binned form. This is for example often the case for data from high energy physics experiments and from astronomy because of finite detector resolution. In some fields this is referred to as histogram data. For the purpose of either the goodness-of-fit or two-sample problems standard discrete data and histogram data can be treated the same, with the midpoints of the bins used as observations where such are needed.

For the two-sample problem p values are found via the permutation method. If the data sets are large for some of the tests the p values can be found via large sample approximations. In the goodness-of-fit case p values are always found via simulation. While large sample approximations are known for some methods such as Kolmogorov-Smirnov and Anderson-Darling, there are no known large sample theories for most of the other tests. Moreover, in the more common situation where the distribution under the null hypothesis depends on parameters, which have to be estimated from the data, even those tests no longer have known large sample theories and one is forced to use simulation to find p values.

The packages *Rgof* [Rolke \(2023b\)](#) and *R2sample* [Rolke \(2023a\)](#) are available from CRAN.

2 Goodness-of-fit / two-sample hybrid problem

As was mentioned in the abstract, while the goodness-of-fit problem and the two-sample problem are quite different, they also share certain features such as methods that exist for both. On a deeper level they are both hypothesis tests in the Fisherian sense, in that they are

tests without an alternative hypothesis. These tests are usually done for confirmation, that is in the goodness-of-fit case the researcher wants to make sure that his assumed probability model is reasonably good, without any consideration of how it might be wrong.

There is yet another connection between these types of problems. Sometimes one wants to carry out a goodness-of-fit test. However, the model under the null hypothesis is quite complex with a large number of nuisance parameters. Therefore calculating values from the distribution function requires integration in high dimensions and is at present not feasible. It is however possible to sample from the distribution. So the problem now changes from a goodness-of-fit to a two-sample problem.

If the null hypothesis in the goodness-of-fit problem does not fully specify the distribution but just its functional form one can then estimate the parameters from the data. However, in this situation the permutation method for estimating the p value fails, it is extremely conservative. Instead the user can provide a routine to generate new data, essentially using a parametric bootstrap approach.

3 The types of problems

The problems that can be analyzed with these packages are as follows:

- **Goodness-of-Fit Problem - Continuous Data:** We have a sample x of size of n drawn from some random variable X . F is a continuous probability distribution, which may depend on unknown parameters. We want to test $X \sim F$.
- **Goodness-of-Fit Problem - Discrete Data:** We have a set of values $vals$ and a vector of counts x . F is a discrete probability distribution, which may depend on unknown parameters. We want to test $X \sim F$.
- **Two-sample Problem - Continuous Data:** We have a sample x of size of n , drawn from some unknown continuous probability distribution F , and a sample y of size m , drawn from some unknown continuous probability distribution G . We want to test $F = G$.
- **Two-sample Problem - Discrete Data:** We have a set of values $vals$ and vectors of counts x and y , drawn from some unknown discrete probability distributions F and G . We want to test $F = G$.

In all of the above problems, the sample size can either be fixed or follow a Poisson distribution with a known rate. In all cases the data can be weighted. In all cases the user can provide his/her own testing method.

4 The methods

In the following we list the methods included in the packages. Most are well known and have been in use for a long time. For their details see the references. They are:

| Method | <i>Rgof</i> | | <i>R2sample</i> | |
|--------------------|-------------|----------|-----------------|----------|
| | Continuous | Discrete | Continuous | Discrete |
| Chi-Square Tests | Yes | Yes | Yes | Yes |
| Kolomarov-Smirnov | Yes | Yes | Yes | Yes |
| Kuiper | Yes | Yes | Yes | Yes |
| Cramer-von Mises | Yes | Yes | Yes | Yes |
| Anderson-Darling | Yes | Yes | Yes | Yes |
| Zhang's tests | Yes | No | Yes | No |
| Wasserstein | Yes | Yes | Yes | Yes |
| Watson's test | Yes | Yes | No | No |
| Lehmann-Rosenblatt | No | No | Yes | Yes |

There are of course many other tests that could have also been implemented in the routines. All the tests included share the following features. They are true omnibus tests, that is not designed with any specific alternative in mind. For this reason we did not include the class of Neyman's smooth tests, for example. Moreover they are all tests that do not depend on some tuning parameters. The exception here are the chi-square tests, which depend on the choice of the number and shape of the bins. The chi-square tests are included because they are so well known and widely used, even so their power often leaves much to be desired.

We denote the cumulative distribution function (cdf) by F , its empirical distribution function (edf) by \hat{F} . In the case of the two-sample problem we also have the edf of the second data set \hat{G} and the edf of the combined data set \hat{H} .

1. Chi-Square Tests

In the case of continuous data the routines include eight chi-square tests, with either equal size (ES) or equal probability (EP) bins, either a large ($nbins[1]=50$) or a small ($nbins[2]=10$) number of bins and with either the Pearson (P) or the log-likelihood (L) formula. Here and in what follows $nbins$ and similar items are arguments to the routines that the user can change. So the combination of a large number of equal size bins and Pearson's chi-square formula is denoted by *ES-l-P*, etc.

In the case of discrete data the type and the number of classes is already given, and for a second test these are combined for a total of $nbins[2]=10$. Again both chi-square formulas are used. So here the case of a large number of bins and Pearson's formula is denoted by *l-P*.

In all cases neighboring bins with low counts are joined until all bins have a count of at least $minexpcount=5$. In all cases the p values are found using the usual chi-square approximation.

If parameters have to be estimated, this is done via the user-provided routine *phat*. As long as the method of estimation used is consistent and efficient and the expected counts are large enough the chi-square statistic will have a chi-square distribution, as shown by Fisher (1922) and Fisher (1924).

Alternatively we can use the argument *ChiUsePhat=FALSE*. In that case the value provided by *phat* is used as a starting point but the parameters are estimated via the method of minimum chi-square. This method has the desirable feature that if the null hypothesis is rejected for this set of values, it will always be rejected for any other as well. For a discussion of this estimation method see Berkson (1980).

2. Kolmogorov-Smirnov (KS)

This test is based on the largest absolute distance between F and \hat{F} in the goodness-of-fit problem and between \hat{F} and \hat{G} in the two-sample problem. The tests were first proposed in Kolmogorov (1933), Smirnov (1939) and are among the most widely used tests today. There is a known large sample distribution of the test statistic in the two-sample problem, which

is used either if both sample sizes exceed 10000 or if the argument *UseLargeSample=TRUE* is set. In the goodness-of-fit case the large sample theory is known only in the case of a fully specified distribution under the null hypothesis. Because this is rarely of interest the large sample approximation is not used.

3. Kuiper (K)

This test is closely related to Kolmogorov-Smirnov, but it uses the sum of the largest positive and negative differences as a test statistic. It was first proposed in [Kuiper \(1960\)](#).

4. Cramer-vonMises (CvM)

This test is based on the integrated squared differences:

- Goodness-of-Fit: $\int_{-\infty}^{\infty} (F(x) - \hat{F}(x))^2 dF(x)$
- Two-Sample: $\int_{-\infty}^{\infty} (\hat{F}(x) - \hat{G}(x))^2 d\hat{H}(x)$

The goodness-of-fit version is discussed in [Cramer \(1928\)](#) and [von Mises \(1928\)](#). The two-sample version was proposed in [Anderson \(1962\)](#).

5. Anderson-Darling (AD)

This test is similar to the Cramer-vonMises test but with an integrand that emphasizes the tails:

- Goodness-of-Fit: $\int_{-\infty}^{\infty} \frac{(F(x) - \hat{F}(x))^2}{F(x)(1-F(x))} dF(x)$
- Two-Sample: $\int_{-\infty}^{\infty} \frac{(\hat{F}(x) - \hat{G}(x))^2}{\hat{H}(x)(1-\hat{H}(x))} d\hat{H}(x)$

It was first proposed in [Anderson and Darling \(1952\)](#). The two-sample version is discussed in [Pettitt \(1976\)](#).

6. Zhang's tests (ZA, ZK and ZC)

These tests were proposed in [Zhang \(2002\)](#) and [Zhang \(2006\)](#). They are variations of test statistics based on the likelihood ratio and different weight functions. Note that these tests do not work for discrete data, that is, they never achieve the correct type I error rate. They are therefore not run for discrete data.

7. Wasserstein p=1 (Wassp1)

A test using the Wasserstein p=1 metric. It is based on a comparison of quantiles. In the goodness-of-fit case these are the quantiles of the data set and the quantiles of the cdf, and in the two-sample problem they are the quantiles of the individual data sets and the quantiles of the combined data set. If $n = m$ the test statistic in the continuous case takes a very simple form: $\frac{1}{n} \sum_{i=1}^n |x_i - y_i|$. In the goodness-of-fit problem for continuous data the user has to supply a function that calculates the inverse of the cdf under the null hypothesis. For a discussion of the Wasserstein distance see [Vaserstein \(1969\)](#).

There are also a number of tests which are only implemented for either the goodness-of-fit or the two-sample problem:

8. Watson's Test (W), Goodness-of-Fit Problem

This test is closely related to the Cramer-vonMises test. It adjust that tests statistic via a squared difference of the mean of $\hat{F}(x_i)$ and 0.5. It was proposed in [Watson \(1961\)](#).

9. Lehmann-Rosenblatt (LR), Two-sample Problem

Let r_i and s_i be the ranks of x and y in the combined sample, then the test statistic is given by

$$\frac{1}{nm(n+m)} \left[n \sum_{i=1}^n (r_i - 1)^2 + m \sum_{i=1}^m (s_i - 1)^2 \right]$$

For details see [Lehmann \(1951\)](#) and [Rosenblatt \(1952\)](#).

5 Simultaneous inference

As no single test can be relied upon to consistently have good power, it is reasonable to employ several of them. We would then reject the null hypothesis if any of the tests does so, that is, if the smallest p value is less than the desired type I error probability α .

This procedure clearly suffers from the problem of simultaneous inference, and the true type I error probability will be much larger than α . It is however possible to adjust the p value so it does achieve the nominal type I error. A sketch of the algorithm is as follows:

- generate a new data set under the null hypothesis, run the desired tests and record the smallest p value.
- repeat $B(=1000)$ times.
- use the empirical distribution function \hat{F}_p of the B smallest p values to estimate their distribution function.
- apply \hat{F}_p to the smallest p value of the data set. This is essentially the probability integral transform.

Here is an example: say the null hypothesis specifies a uniform $[0, 1]$ distribution and a sample size of 250. Next we find the smallest p value in each run for two selections of four methods. One includes the methods by Wilson, Anderson-Darling, Zhang's ZC and a chi square test with a small number of bins and using Pearson's formula. This selection has good power against a large number of alternatives. As a second selection we use the methods by Kolmogorov-Smirnov, Kuiper, Anderson-Darling and Cramer-vonMises. For the case where the null hypothesis specifies a Uniform $[0, 1]$ distribution these tests turn out to be highly correlated.

Next we find the empirical distribution function for the two sets of p values and draw their graphs. We also add the curve for the cases of four identical tests and the case of four independent tests, which of course is the Bonferroni correction. These are shown in figure 1.

As one would expect, the two curves for the p values fall between the extreme cases of total dependence and independence. Moreover, the curve of our good selection is closer to the curve of independence than the selection of correlated methods.

Finally we simply have to apply this function to the smallest p value found for the actual data.

`Rgof::gof_test_adjusted_pvalues` and `R2sample::twosample_test_adjusted_pvalues` find these adjusted p values. Their arguments are the same as those of `Rgof::gof_test` and `R2sample::twosample_test`, see the section Usage 2.7. For an example that uses this adjustment method in the context of simultaneous confidence bands see [Aldor-Noima et al. \(2013\)](#).

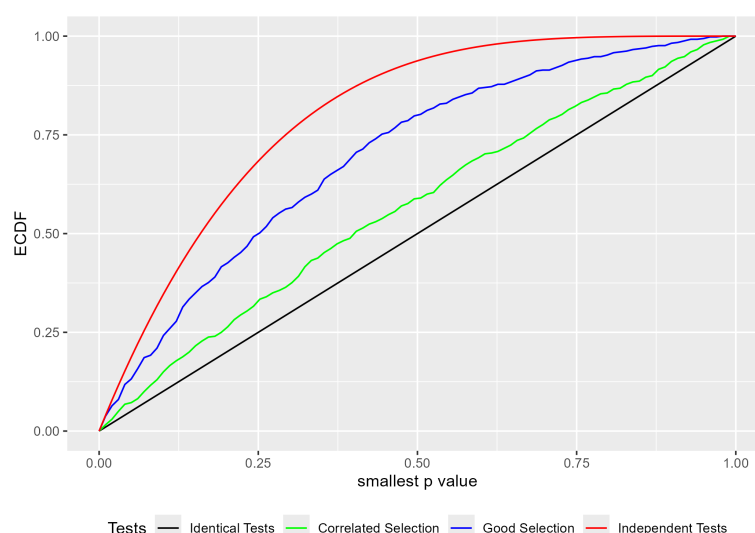


Figure 1: Distribution functions of the smallest p value for four dependence cases.

6 Special circumstances

6.1 Random sample size

In some cases the sample size is not determined at the beginning of the experiment but is a consequence of other factors. As an example, in high energy physics experiments the sample size is determined by the luminosity (aka energy) at which the accelerator is run, and by how long. In such a situation the distributions of the test statistics are different from the fixed sample size case, and there are no known null distributions. In the case of the chi-square tests, for example, the bin counts are now independent according to the theory of marked Poisson processes, and therefore the degrees of freedom need to be adjusted accordingly. Often though it is reasonable to assume that the sample size has a Poisson distribution. If so the routines in *Rgof* and *R2sample* have an argument *rate* to indicate a random sample size from a Poisson distribution with rate λ .

6.2 Weighted samples

Another variation is as follows. Say we have a continuous random variable X and a weight function w . There also exists a random variable Y such that $E[f(Y)] = E[f(X)w(x)]$ for (almost) any function f . In other words, these are weights as one encounters in importance sampling. Say we wish to test whether the distribution of Y is F but using observations from X with their weights. This is done very easily with the routines by supplying the weights as arguments. These weights can then be used to find for example the empirical distribution function, and with it run tests such as Kolmogorov-Smirnov or Anderson-Darling.

One field where this situation is common is high energy physics. There we have the Standard Model, the current best model for explaining the physics of collision experiments in particle accelerators such as the Large Hadron Collider at CERN. Say we wish to test some specific part of this theory, that is we want to do a goodness-of-fit test. However, the Standard Model depends on dozens of parameters. The calculations of the probabilities needed for a goodness-of-fit test are at present not feasible. Among other issues they would require integration in very high dimensions. However, it is possible to generate a Monte Carlo sample from the Standard Model, so instead we can run a two-sample test, comparing the data to the Monte Carlo sample. There is however another problem. The Monte Carlo sampling of the Standard Model is very computationally expensive. There exist a number of such samples, each for a specific set of the parameters of the Standard Model. Now if the test we wish to do requires a sample with a slightly different set of parameters we can use

an existing sample and importance sampling. The routines in the packages discussed here make this very easy.

7 Usage

7.1 Goodness-of-fit problem - testing

The routine to carry out hypothesis tests is *Rgof::gof_test*. It's arguments are

- *x*: a data set, either the continuous outcomes or the counts in the discrete case.
- *vals=NA*: all possible values of the discrete random variable, or NA if data is continuous.
- *pnull*: a function to calculate values for the cdf.
- *rnull*: a function to generate new data under the null hypothesis.
- *w=function(x) -99*: a weight function if weights are present, or -99 if not.
- *phat=function(x) -99*: a function to estimate parameters, or -99 if null hypothesis is simple and no parameters are estimated.
- *TS*: routine to calculate test statistics other than those included.
- *TSextra*: a list passed to *TS*.
- *nbins=c(50, 10)*: number of bins to use in chi-square tests.
- *rate=0*: rate of Poisson if sample size is random, 0 if sample size is fixed.
- *Range=c(-Inf, Inf)*: range of continuous random variable.
- *B=5000*: number of simulation runs.
- *minexpcount=5*: required minimal expected counts for chi-square tests.
- *maxProcessors=1*: number of cores to use for parallel processing, 1 means no parallel programming.
- *doMethod="all"*: vector with names of methods, if not all are to be included.

The format of the routines *pnull*, *rnull* and *w* has to be as follows. In the continuous case we will use as an example the normal distribution and in the discrete case the Binomial distribution with 10 tries.

- Continuous data, no parameter estimation: a function of one variable, the data. For example, *pnull=function(x) pnorm(x)*
- Continuous data, with parameter estimation: a function of two variables, the data and a vector of parameter estimates. For example, *pnull=function(x,p) pnorm(x,p[1],p[2])*
- Discrete data, no parameter estimation: a function without arguments. For example, *pnull=function() pbinom(0:10, 10, 0.5)*
- Discrete data, with parameter estimation: a function of one variable, a vector of parameter estimates. For example, *pnull=function(p) pbinom(0:10, 10, p)*

Continuous data

As an example we generate $N = 1000$ observations from a standard normal distribution. Then we test to see whether the data comes from a normal distribution with the mean estimated from the data, so in this case the null hypothesis is true:

```
pnull = function(x, mu=0) pnorm(x, mu) # cdf under null hypothesis
rnull = function(mu=0) rnorm(1000, mu) # generate data under null hypothesis
phat = function(x) mean(x) # estimate parameter
x = rnull() # data from distribution under the null hypothesis
Rgof::gof_test(x, NA, pnull, rnull, phat=phat)

#> $statistics
#>      KS      K      AD      CvM      W      ZA      ZK      ZC
#>  0.0208  0.0407  0.6500  0.0926  0.0911  3.2940  1.4780 12.4900
#>  ES-l-P  ES-s-P  EP-l-P  EP-s-P  ES-l-L  ES-s-L  EP-l-L  EP-s-L
#> 37.4800  9.6060 51.7300 11.9200 38.5700  9.4030 50.2600 12.4700
#>
#> $p.values
#>      KS      K      AD      CvM      W      ZA      ZK      ZC
#> 0.4712 0.4712  0.2328 0.2546  0.2354  0.6292  0.7692  0.5828
#> ES-l-P  ES-s-P  EP-l-P  EP-s-P  ES-l-L  ES-s-L  EP-l-L  EP-s-L
#> 0.4011  0.2120  0.3302  0.1550  0.3542  0.2250  0.3840  0.1315
```

If we wish to find an adjusted p value for a combination of tests we can run

```
Rgof::gof_test_adjusted_pvalue(x, NA, pnull, rnull, phat=phat)

#> p values of individual tests:
#> W : 0.128
#> ZC : 0.232
#> AD : 0.08
#> ES-s-P : 0.212
#> adjusted p value of combined tests: 0.2267
```

Next we generate a data set from a t distribution with 5 degrees of freedom, so now the null hypothesis is false. Here and in the examples that follow we only show the p values of the tests:

```
y = rt(1000, 5) # data where null hypothesis is false
Rgof::gof_test(y, NA, pnull, rnull, phat=phat)[["p.values"]]

#>      KS      K      AD      CvM      W      ZA      ZK      ZC
#>  0.0106  0.0106  0.0000  0.0048  0.0034  0.0000  0.0000  0.0000
#>  ES-l-P  ES-s-P  EP-l-P  EP-s-P  ES-l-L  ES-s-L  EP-l-L  EP-s-L
#>  0.0000  0.0000  0.0000  0.0000  0.0000  0.0000  0.0000  0.0408
```

If the quantile function (aka inverse) of the cdf is known it can be included. It is then used in some of the chi-square tests and the Wasserstein test. It can be passed to the routine that finds the test statistic(s) via the list TSextra:

```
TSextra = list(qnull=function(x, mu) qnorm(x, mu))
Rgof::gof_test(x, NA, pnull, rnull, phat=phat, TSextra=TSextra)[["p.values"]]

#>      KS      K      AD      CvM      W      ZA      ZK      ZC  Wassp1
#>  0.4710  0.4710  0.2288  0.2444  0.2238  0.6198  0.7664  0.5762  0.1760
#>  ES-l-P  ES-s-P  EP-l-P  EP-s-P  ES-l-L  ES-s-L  EP-l-L  EP-s-L
#>  0.4011  0.2120  0.5634  0.4413  0.3542  0.2250  0.5182  0.4210
```

A user can also use this routine to run their own test. For example, let's say we wish to include the Neyman's smooth test from the *DDST* (P Niecek 2016) package:

```
NeymanSmoothTest = function(x, pnull, param) {
  ts=as.numeric(unlist(ddst::ddst.norm.test(x))[1])
  names(ts) = "DDST"
  ts
}
Rgof::gof_test(x, NA, pnull, rnull, phat=phat, TS=NeymanSmoothTest)[["p.values"]]

#> DDST
#> 0.5482

Rgof::gof_test(y, NA, pnull, rnull, phat=phat, TS=NeymanSmoothTest)[["p.values"]]

#> DDST
#> 0
```

The routine has to have the following form:

```
newTS(x, pnull, param, TSextra)
```

x is the data set and $pnull$ the distribution function under the null hypothesis, as described above. $param$ is the estimated parameters in the case of a composite null hypothesis and is ignored in the case without parameter estimation. The argument $TSextra$, a list of items also needed for calculating the test statistic, is optional.

Next we assume that the sample size was random and drawn from a Poisson distribution with rate 950. One of the consequences of this is that now the degrees of freedom of the chi-square tests is the number of bins - number of estimated parameters rather than number of bins - 1 - number of estimated parameters.

```
Rgof::gof_test(x, NA, pnull, rnull, phat=phat, TSextra=TSextra, rate=950)[["p.values"]]

#> KS K AD CvM W ZA ZK ZC Wassp1
#> 0.4736 0.4736 0.2406 0.2580 0.2372 0.6206 0.7642 0.5782 0.1904
#> ES-l-P ES-s-P EP-l-P EP-s-P ES-l-L ES-s-L EP-l-L EP-s-L
#> 0.2602 0.1210 0.4009 0.2779 0.2936 0.1517 0.4539 0.2957
```

As an example for the use of importance sampling weights we generate the data from a mixture of two normal random variables but as above test for a simple normal distribution with unknown mean.

```
rnull = function(mu=0) c(rnorm(500, -1), rnorm(500, 1))
x = rnull()
w=function(x, mu=0) dnorm(x, mu)/(dnorm(x, -1)/2+dnorm(x, 1)/2)
Rgof::gof_test(x, NA, pnull, rnull, w=w, phat=phat)[["p.values"]]

#> KS K CvM AD
#> 0.7990 0.7718 0.6330 0.4516

Rgof::gof_test(y, NA, pnull, rnull, w=w, phat=phat)[["p.values"]]

#> KS K CvM AD
#> 0 0 0 0
```

Discrete data

Here we will consider the following example. The null hypothesis specifies a binomial distribution with $n = 100$ trials and a success probability p , estimated from the data. As an example where the null hypothesis is false we generate data that is a mixture of a binomial distribution with $p = 0.5$ and a discrete uniform distribution on the integers from 30 to 70.

```
set.seed(1234)
vals = 0:100 # all possible values
pnull = function(p=0.5) pbinom(0:100, 100, p)
rnull = function(p=0.5) table(c(0:100,rbinom(1000, 100, p)))-1
phat = function(x) mean(0:100*x)/1000
x = rnull()
Rgof::gof_test(x, vals, pnull, rnull, phat=phat)$p.values

#>      KS      K      AD      CvM      W      Wassp1      l-P      s-P      l-L      s-L
#> 0.2320 0.2752 0.4858 0.3132 0.2460 0.6257 0.1599 0.5592 0.1456 0.1456

y = table(c(0:100, rbinom(900, 100, 0.5), sample(30:70, size=100, replace=TRUE)))-1
Rgof::gof_test(y, vals, pnull, rnull, phat=phat)$p.values

#>      KS      K      AD      CvM      W      l-P      s-P      l-L      s-L
#> 0.3846 0.3290 0.0000 0.0174 0.0114 0.0000 0.0000 0.0000 0.0000
```

Here we have an example where most tests correctly reject the null hypothesis but some do not.

Note that the routine *rnull* has to insure that all values of *vals* are present, even if many have counts of zero.

Again the user can provide his/her own test statistic. The routine has to be as follows:

```
newTS(x, pnull, param, vals, TSextra)
```

Here *x* is the counts and *pnull* the distribution function under the null hypothesis as described above. *param* is the estimated parameters in the case of a composite null hypothesis and is ignored in the case without parameter estimation. *vals* is the set of values where $P(X = vals) > 0$. The argument *TSextra*, a list of items also needed for calculating the test statistic, is optional.

7.2 Goodness-of-fit problem - power estimation

To estimate the power of the various tests we can use the function *gof_power*. It's arguments are the same as *gof_test*, as well as

- *ralt*: a function that generates data under the alternative hypothesis.
- *param_alt*: a vector of values to be passed to *ralt*.
- *With.p.value*=FALSE: set to TRUE if the new user supplied routine calculates p values.
- *alpha*=0.05: type I error probability to be used for test.
- *B*=1000: number simulation runs.

As an example say we wish to estimate the power of the tests when the null hypothesis specifies a normal distribution, but the data comes from a t distribution. We have 500 observations, and both the mean and standard deviation are estimated from the data. The package also includes the routine *plot_power*, which has as its argument the output from the *gof_power* command and draws the power curve. It is shown in figure 2.

```

pnull = function(x, p=c(0,1)) pnorm(x, p[1], p[2]) # cdf under null hypothesis
rnull = function(p=c(0,1)) rnorm(500, p[1], p[2]) # generate data under null hypothesis
phat = function(x) c(mean(x), sd(x)) # estimate parameters
TSextra = list(qnull = function(x, p=c(0,1)) qnorm(x, p[1], p[2])) # quantile function
ralt = function(df) rt(500, df) # generate data under alternative
tmp=Rgof::gof_power(pnull, NA, rnull, ralt, param_alt=4*1:10, phat=phat, TSextra = TSextra)
Rgof::plot_power(tmp, "df", "Standard Normal vs t Distributions")

```

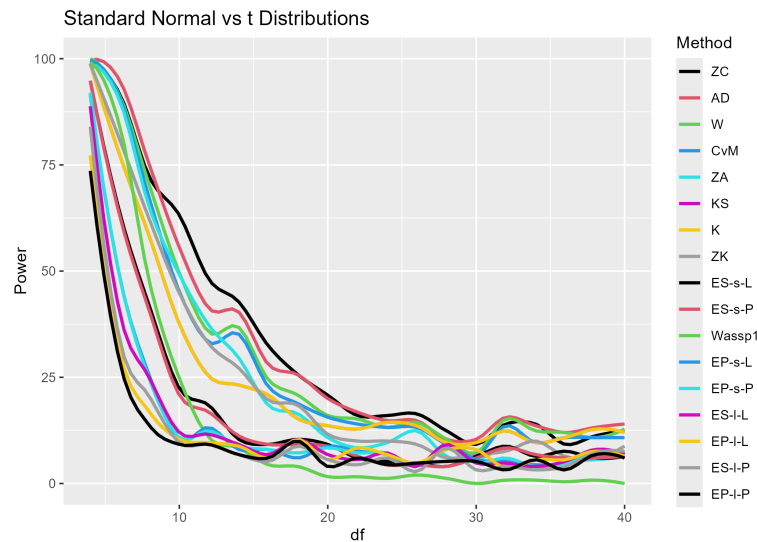


Figure 2: Power graph for goodness-of-fit-tests of normal vs t distributions, with mean and standard deviation estimated from the data.

plot_power has the arguments

- *pwr*: a matrix, usually the output of the *gof_power* command.
- *xname*: name of parameter of *ralt*.
- *title=""*: title of graph
- *Smooth=TRUE*: should curves be smoothed?
- *span=0.25*: parameter for smoothing routine.

Power estimation for discrete data works the same. As an example consider the following. One data set comes from a Poisson distribution, restricted to the set of integers from 70 to 140. The second data set is a 50-50 mixture of Poisson random variables with a rate of 100 and a rate of $100 + \lambda$. The power graph is in figure 3.

```

vals=70:140
pnull=function(lambda) (ppois(70:140, lambda)-ppois(69,lambda))/(ppois(140, lambda)-ppois(69,lambda))
rnull=function(lambda) {
  vals=70:140
  x=rpois(1000, lambda)
  x[x<70]=70
  x[x>140]=140
  x=table(c(70:140, x))-1
}
phat=function(x) sum(70:140*x)/1000
ralt=function(lambda) {
  vals=70:140

```

```

x=c(rpois(500, 100), rpois(500, 100+lambda))
x[x<70]=70
x[x>140]=140
x=table(c(70:140, x))-1
}
tmp=Rgof::gof_power(pnull, vals, rnull, ralt, param_alt=2*0:8, phat=phat)
Rgof::plot_power(tmp, "lambda", "Poisson vs Mixture of Poisson Distributions")

```

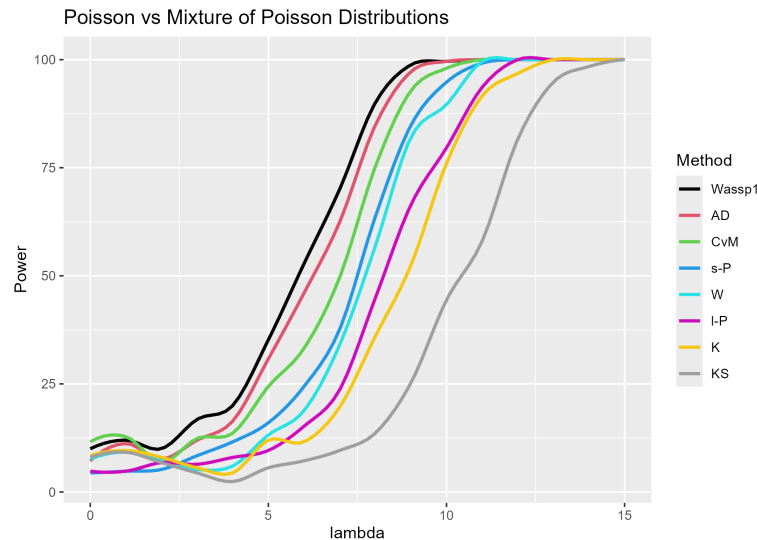


Figure 3: Power graph for goodness-of-fit-tests of Poisson vs a mixture of Poisson distributions, with the rate estimated from the data.

7.3 Two-sample problem - testing

p values are generally found using the permutation method. The idea of a permutation test is simple. Say we have data sets x_1, \dots, x_n and y_1, \dots, y_m . They are combined into one large data set $x_1, \dots, x_n, y_1, \dots, y_m$, permuted and split again in sets of size n and m. Under the null hypothesis these new data sets come from the same distribution as the actual data. Therefore calculating the tests statistics for them and repeating many times one can build up the distributions of the test statistics and find p values from them.

In the case of continuous data the routines also allow for the use of large sample formulas for some of the tests. In the discrete case none of the methods (outside of the chi-square tests) has a known large sample theory.

In the discrete case the permutation method is somewhat more complicated. Say we have data sets with values v_1, \dots, v_k and counts $x_1, \dots, x_k, y_1, \dots, y_k$. One can then simply expand these to yield a large data set with $x_1 + y_1$ v_1 's, $x_2 + y_2$ v_2 's and so on. Then this vector is permuted and split as described above. The drawback of this sampling method is that its calculation speed increases with the sample size and would be impossible for data sets with very large counts.

Alternatively *R2sample* provides the following option. One can show that the distribution of the permuted data sets is as follows: let $n = \sum x_i$ and $m = \sum y_i$, then

$$P(\mathbf{X} = \mathbf{a} | \mathbf{x}, \mathbf{y}) = \left[\prod_{j=1}^k \binom{x_j + y_j}{a_j} \right] / \binom{n+m}{n}$$

for any \mathbf{a} such that $0 \leq a_i \leq x_i; i = 1, \dots, k$ and $\sum a_i = n$.

It is possible to sample from this distribution as follows: Let N and M be the sample sizes of the two data sets, respectively. Let $p = N/(N + M)$ be the proportion of events

in the first data set. Say that in a data set we have x_1 observations of the smallest discrete value v_1 in the first data set and y_1 in the second. We can then generate a random sample by drawing an observation from a Binomial with parameters $x_1 + y_1$ and p . We repeat this for all values of the discrete random variable. We also need to insure, though, that the total number of observations in the first simulated data set is again N and in the second data set M . If this is not so, we randomly choose two values, giving a higher priority for those with high counts, and flipping one observation between the two data sets. This is repeated until the simulated data set has N events total.

The routine is `twosample_test`. It's arguments are

- `x, y`: the two data sets, either the observations in the continuous case or the counts in the discrete case. `x` can also be a list with elements `x` and `y`, and then `y` is ignored.
- `vals=NA`: the possible values of the discrete random variables, or `NA` for continuous data.
- `TS`: routine to calculate test statistics other than those included
- `TSextra`: a list passed to `TS`
- `wx = rep(1, length(x))`: weights for `x` data
- `wy = rep(1, length(y))`: weights for `y` data
- `B=5000`: number of simulation runs
- `nbins=c(50,10)`: number of bins to use in chi-square tests
- `minexpcount=5`: required minimal expected counts for chi-square tests
- `maxProcessors`: number of cores to use for parallel processing, default is 1 less than are detected on computer
- `UseLargeSample`: should large sample approximations be used instead of permutation?
- `samplingmethod="Binomial"`: sampling method for discrete data.
- `rnull`: a function that generates data from a model, possibly with parameter estimation. This is needed in the goodness-of-fit/two-sample hybrid problem.
- `doMethod="all"`: vector with names of methods, if not all are to be included.

The arguments match those of `gof_test`, where this makes sense.

Continuous data

The `x` and `y1` data sets come from a standard normal distribution, and `y2` from a normal distribution with mean 1.

```
x = rnorm(100)
y1 = rnorm(150)
y2 = rnorm(150, 1)
R2sample::twosample_test(x, y1)[["p.values"]]

#>      KS   Kuiper    CvM     AD      LR      ZA      ZK      ZC
#> 0.9514 0.8730 0.9242 0.9126 0.8992 0.7480 0.4502 0.7888
#> Wassp1 ES large ES small EP large EP small
#> 0.8580 0.6309 0.2454 0.4334 0.9496

R2sample::twosample_test(x, y2)[["p.values"]]
```

```
#>      KS   Kuiper      CvM      AD      LR      ZA      ZK      ZC
#>      0       0       0       0       0       0       0       0
#>  Wassp1 ES large ES small EP large EP small
#>      0       0       0       0       0
```

Again, the user can provide their own test statistic:

```
DiffStandardizedMeans = function(x, y) {
  TS = abs(mean(x)/sd(x)-mean(y)/sd(y))
  names(TS) = "DSM"
  TS
}
R2sample::twosample_test(x, y1, TS=DiffStandardizedMeans)[["p.values"]]

#>      DSM
#>      0.585

R2sample::twosample_test(x, y2, TS=DiffStandardizedMeans)[["p.values"]]

#> DSM
#>      0
```

The user supplied routine has to be a function of the two data sets *x* and *y* and optionally a list *TSextra*.

As an example for weighted data, let's say the *x* data set actually came from a *t* distribution with 5 degrees of freedom:

```
x = rt(100, 5)
wx = dnorm(x)/dt(x, 5)
R2sample::twosample_test(x, y1, wx=wx)[["p.values"]]

#>      KS   Kuiper      CvM      AD   ES large ES small EP large EP small
#>      0.2400 0.2354 0.2522 0.2560 0.7627 0.8632 0.3814 0.8636

R2sample::twosample_test(x, y2, wx=wx)[["p.values"]]

#>      KS   Kuiper      CvM      AD ES large ES small EP large EP small
#>      0.0004 0.0004 0.0000 0.0000 0.0001 0.0000 0.0270 0.0000
```

If the data sets are very large using permutation to derive the null distribution of the test statistics can be very slow. In this case one can use the argument *UseLargeSample=TRUE*. This will be done automatically if both sample sizes are at least 10000.

```
x = rnorm(1e5)
y1 = rnorm(1e5)
y2 = rnorm(1e5, 0.02)
R2sample::twosample_test(x, y1)[["p.values"]]

#>      KS   Kuiper      CvM      AD
#>      0.098 0.1214 0.1602 0.1330
#>      ES large ES small EP large EP small
#>      0.2683 0.3347 0.0974 0.1988

R2sample::twosample_test(x, y2)[["p.values"]]

#>      KS   Kuiper      CvM      AD
#>      0.0112 0.0542 0.0046 0.0013
#>      ES large ES small EP large EP small
#>      0.0935 0.0074 0.6170 0.0717
```

Discrete data

As an example for the case of discrete data we will use two data sets from geometric random variables with slightly different rates. x and y have to have the same length as $vals$ and $x+y$ has to be positive for all values in $vals$.

```
x = table(rgeom(1000, 0.7))
y = table(rgeom(1000, 0.8))
vals = unique(c(names(x), names(y))) # all values from either x or y
x1 = rep(0, length(vals))
names(x1)=vals
y1 = x1
x1[names(x)]=x
y1[names(y)]=y
vals = as.numeric(vals)
R2sample::twosample_test(x1, y1, vals)[["p.values"]]

#>      KS Kuiper      CvM      AD      LR      ZA Wassp1  large  small
#>      0        0        0        0        0        0        0        0
```

Again the user can supply their own test. The routine has to be a function of the two data sets x and y and a vector $vals$ of possible values of the discrete random variable. A list $TSextra$ is optional.

7.4 Two-sample problem - power estimation

The package *R2sample* includes the routine *twosample_power*. The arguments TS , $TSextra$, B , $nbins$, $minexpcount$, $UseLargeSample$ and $maxProcessor$ are the same as in *twosample_test*. In addition we have

- f : a function that generates a list with two vectors called x and y and (in the case of discrete data) a vector $vals$. The function can have zero, one or two arguments.
- ... arguments passed to f .
- $With.p.value=FALSE$: set to $TRUE$ is user supplied routine calculates p values.
- $alpha=0.05$: type I error probability for the tests.

As an example for continuous data we again consider the case of normal vs t distributions. One data set with 10001 observations comes from a standard normal distribution, the other with 10002 observations from a t distribution with df degrees of freedom. Because of the large sample sizes the large sample approximations are used.

```
f=function(df) {
  x=rnorm(10001)
  y=rt(10002, df)
  list(x=x, y=y)
}
tmp=R2sample::twosample_power(f, df=seq(5, 100, 5))
R2sample::plot_power(tmp, "df", "Data from Standard Normal vs t Distributions")
```

For discrete data we again consider the case of a Poisson random variable vs a 50-50 mixture of Poisson random variables:

```
f=function(a) {
  vals=70:140
```

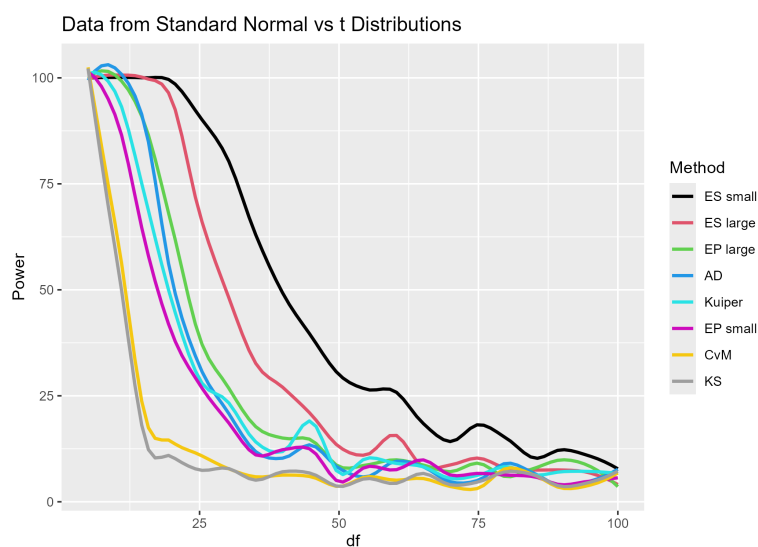


Figure 4: Power graph for two-sample tests where one data set comes from a standard normal distribution and the other from a t distribution.

```
x=rpois(1000, 100)
x[x<70]=70
x[x>140]=140
x=table(c(70:140, x))-1
y=c(rpois(500, 100), rpois(500, 100+a))
y[y<70]=70
y[y>140]=140
y=table(c(70:140, y))-1
I=seq_along(vals)[x+y>0]
list(x=x[I], y=y[I], vals=vals[I])
}
tmp=R2sample::twosample_power(f, a=seq(0,5,0.25))
R2sample::plot_power(tmp, "lambda", "Data from Poisson vs Mixture of Poisson Distributions")
```

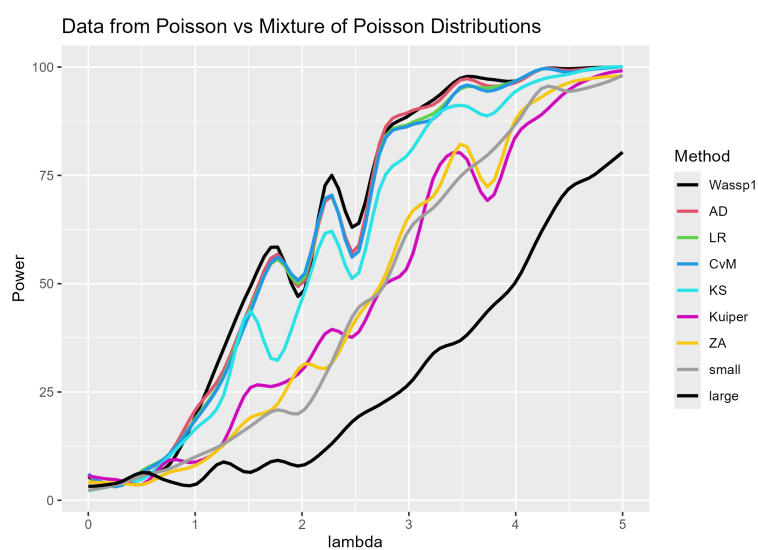


Figure 5: Power graph for two-sample tests where one data set comes from Poisson distribution and the other from a mixture of two Poisson distributions.

8 Benchmarking

Say a researcher has developed a new method for the univariate goodness-of-fit problem and wants to see how it stacks up in comparison to the standard methods such as the chi-square tests or the Kolmogorov-Smirnov test. Both *Rgof* and *R2sample* packages include the routine *run_studies*, which makes this very easy.

As a specific example say we wish to see whether the Kolmogorov-Smirnov test or the Anderson-Darling test has better power when the null hypothesis specifies a normal distribution with mean and standard deviation unspecified, but in reality the data comes from a *t* distribution. Say we run the *Rgof::gof_power* command for a sample size of 500, a *t* distribution with 8 degrees of freedom and a true type I error of 5%, and we find that the Kolmogorov-Smirnov test has a power of 43% whereas the Anderson-Darling test has a power of 70%. It is then true that the Anderson-Darling test will also have a higher power for any other combination of sample size, degrees of freedom and true type I error. In order to assess the ranking of the methods it therefore suffices to run each case study with just one combination of *n*, α and the parameter under the alternative.

8.1 Goodness-of-fit problem

Here the user needs to create a function that finds either the test statistic or, if possible, the *p* value of the new test. Its arguments should be as described previously. Say the routine is called *myTS* and is designed for continuous data and calculates the test statistic. Then the user can run

```
run_studies(myTS)
```

This will run 20 different case studies and provide some information on the relative power of the new method when compared to those included in *Rgof*. For a list of the case studies see the appendix.

If the routine actually calculates a *p* value, run instead

```
run_studies(myTS, With.p.value=TRUE)
```

This will of course be much faster as it does not require simulation.

The arguments of *run_studies* are

- *TS*: the name of the new test routine
- *study*: the name of the study to run, or all studies if missing
- *TSextra*: a list of additional info passed to *TS*, if such are needed
- *With.p.value*=FALSE: TRUE if routine finds *p* values
- *BasicComparison*=TRUE: if TRUE the values for sample size, type I error etc from included studies are used.
- *nsample* = 500: desired sample size
- *alpha* = 0.05: desired type I error probability
- *param_alt*: (list of) parameters for alternative distributions
- *maxProcessor*: number of cores to use for parallel programming, number of cores - 1 if missing
- *B* = 1000: number of simulation runs

As an example we will use the R built-in *ks.test* to do the Kolmogorov-Smirnov test:

```

myTS=function(x, pnull, param) {
  if(length(formals(pnull))==1) # case studies without parameter estimation
    mypnull=function(x) pnull(x)
  else mypnull=function(x) pnull(x, param) # case studies with parameter estimation
  z=ks.test(x, mypnull)[["p.values"]]
  names(z)="RKS"
  z
}
pwr=run.studies(myTS, With.p.value = TRUE)

#> Average number of studies a method is close to the best::
#> EP-1-P ES-1-P EP-1-L RKS ES-1-L EP-s-P EP-s-L KS K ES-s-P
#> 5.450 5.9 6.200 6.900 6.950 8.875 9.175 9.400 9.400 10.175
#> W CvM ES-s-L ZA ZK Wassp1 ZC AD
#> 10.525 10.550 10.850 11.625 11.850 12.050 12.300 13.050

```

Note that the performance of RKS is much lower than that of the Kolmogorov-Smirnov test included in the *Rgof* package. This is due to the fact the *R* routine `ks.test` does not actually allow for parameter estimation.

As another example say the user routine calculates the test statistic and the user wants to find the power of the methods in the case where the null hypothesis specifies a normal distribution with mean and standard deviation estimated from the data, but the true distribution is a *t* distributions with *df* degrees of freedom. He also wants a sample size of 1000 and a true type I error of 0.1 Then he can run

```
run_studies(myTS, "normal.t.est", nsample=1000, alpha=0.1, param_alt=5*1:10)
```

8.2 Two-sample problem

This works in the exact same way as in the goodness-of-fit problem.

8.3 Real data examples

R sunspots data

This data set has the monthly mean relative sunspot numbers from 1749 to 1983, collected at Swiss Federal Observatory, Zurich until 1960, then Tokyo Astronomical Observatory. It is part of the base *R* program. It is of course a time series, but we will treat it here as if it were independent data.

A histogram of the data suggests that an exponential model might be a good fit. However, all the tests in *Rgof* reject that null hypothesis:

```

library(ggplot2)
dta=data.frame(Sunspots=c(sunspots))
ggplot(dta, aes(x=Sunspots)) +
  geom_histogram(color="black", fill="white") +
  ggtitle("Sunspots Data")

pnull = function(x, p=1) pexp(x, p)
TSextra=list(qnull= function(x, p=1) qexp(x, p))
rnull = function(p) rexp(N, p)
phat = function(x) 1/mean(x)
Rgof::gof_test(c(sunspots), NA, pnull, rnull, phat=phat, TSextra = TSextra, Range=c(0, Inf))

```

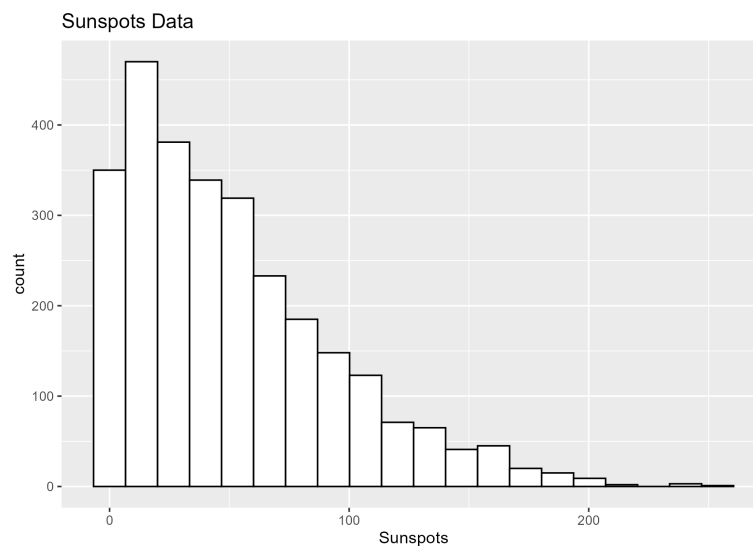


Figure 6: Histogram of sunspots data set.

```
#> $statistics
#>      KS      K      AD      CvM      W      ZA      ZK      ZC
#> 0.0631 0.1260 0.0000 2.8873 1.8894 0.0000 0.0000 0.0000
#> Wassp1 ES-l-P ES-s-P EP-l-P EP-s-P ES-l-L ES-s-L EP-l-L
#> 5.2559 154.7900 100.0300 252.8700 84.6450 170.2800 122.7200 255.0200
#> EP-s-L
#> 84.1610
#>
#> $p.values
#>      KS      K      AD      CvM      W      ZA      ZK      ZC Wassp1 ES-l-P ES-s-P
#>      0      0      0      0      0      0      0      0      0      0      0
#> EP-l-P EP-s-P ES-l-L ES-s-L EP-l-L EP-s-L
#>      0      0      0      0      0      0
```

Death by horsekicks data

This is the famous data set first discussed in [Bortkewitsch \(1898\)](#) and analyzed in many statistics text books and articles, see for example [Preece et al. \(1988\)](#). It is the number of soldiers in the Prussian army killed by being kicked by a horse between 1875 to 1894. The data is

| Number of Deaths | Frequencies |
|------------------|-------------|
| 0 | 109 |
| 1 | 65 |
| 2 | 22 |
| 3 | 3 |
| 4 | 1 |
| 5+ | 0 |

We will run the tests to see whether a Poisson model might fit this data set:

```
vals=0:5
x=c(109, 65, 22, 3, 1, 0)
pnul1 = function(lambda=1) c(ppois(0:4,lambda), 1)
rnul1 = function(lambda=1) {
  x = rpois(200, lambda)
```

```

x[x>5]=5
table(c(0:5, x))-1
}
phat = function(x) sum(0:5*x)/200
Rgof::gof_test(x, vals, pnull, rnull, phat=phat)

#> $statistics
#>      KS      K      AD      CvM      W Wassp1      l-P      s-P      l-L      s-L
#> 0.5434 0.5481 0.0323 0.0026 4.5350 0.0124 0.0628 0.0628 0.0625 0.0625
#>
#> $p.values
#>      KS      K      AD      CvM      W Wassp1      l-P      s-P      l-L      s-L
#> 0.4830 0.6372 0.9434 0.9412 0.5288 0.9386 0.8021 0.8021 0.8026 0.8026

```

All methods fail to reject the null hypothesis, so the Poisson distribution is a good model for the horsekick data.

9 Conclusion

The R packages *Rgof* and *R2sample* bring together a large number of methods for the univariate goodness-of-fit problem and the univariate two-sample problem. The routines make it easy to run these tests simultaneously. They are implemented for both continuous and discrete data and can handle a number of different situations such as random sample sizes and weighted data. The packages also include routines for power estimation as well as routines for benchmarking new tests.

10 Appendix

10.1 Case Studies for Goodness-of-fit Problem

Without parameter estimation

- | | |
|---|--|
| 1. uniform.linear | U[0,1] vs a linear model on [0,1] with slope s. |
| 2. uniform.quadratic some curvature a. | U[0,1] vs a quadratic model with vertex at 0.5 and |
| 3. uniform.bump | U[0,1] vs U[0,1]+N(0.5,0.05). |
| 4. uniform.sine | U[0,1] vs U[0,1]+Sine wave |
| 5. beta22.betaaa | Beta(2,2) vs Beta(a,a) |
| 6. beta22.beta2a | Beta(2,2) vs Beta(2,a) |
| 7. normal.shift | N(0,1) vs N(μ ,1) |
| 8. normal.stretch | N(0,1) vs N(0, σ) |
| 9. normal.t | N(0,1) vs t(df) |
| 10. normal.outlier1 | N(0,1) vs N(0,1)+U[2,3] |
| 11. normal.outlier2 | N(0,1) vs N(0,1)+U[-3,-2]+U[2,3] |
| 12. exponential.gamma | Exp(1) vs Gamma(1,b) |
| 13. exponential.weibull | Exp(1) vs Weibull(1,b) |
| 14. exponential.bump | Exp(1) vs Exp(1)+N(0.5,0.05) |

15. trunc.exponential.linear Exp(1) vs Linear, on [0,1]

With parameter estimation

16. normal.t.est $N(\mu, \sigma)$ vs $t(df)$
 17. exponential.weibull.est Exp(λ) vs Weibull(1,b)
 18. trunc.exponential.linear.est Exp(λ) vs Linear, on [0,1]
 19. exponential.gamma.est Exp(λ) vs Gamma(1,b)
 20. normal.cauchy.est $N(\mu, \sigma)$ vs Cauchy (Breit-Wigner)

10.2 Case Studies for two-sample problem

The first 14 case studies are the same as those of the goodness-of-fit problem. The others are

15. gamma.normal Gamma(μ) vs $N(\bar{x}, sd(x))$, here the mean of the normal distribution are the sample mean and sample standard deviation of the x data set.
 16. normal.normalmixture $N(0,1)$ vs $N(-\mu,1)+N(\mu,1)$
 17. uniform.uniformmixture $U[0,1]$ vs. $\alpha U[0,1/2]+(1-\alpha)U[1/2,1]$
 18. uniform.betamixture $U[0,1]$ vs. $\alpha U[0,1/2]+(1-\alpha)\text{Beta}(2,2)$
 19. chisquare.noncentral $\chi^2(5)$ vs. $\chi^2(5, \tau)$
 20. uniform.triangular $U[0,1]$ vs. triangular

References

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