

Sample selection models for count data in R

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Abstract We provide a detailed hands-on tutorial for the R package **SemiParSampleSel** (version 1.5). The package implements selection models for count responses fitted by penalized maximum likelihood estimation. The approach can deal with non-random sample selection, flexible covariate effects, heterogeneous selection mechanisms and varying distributional parameters. We provide an overview of the theoretical background and then demonstrate how **SemiParSampleSel** can be used to fit interpretable models of different complexity. We use data from the German Socio-Economic Panel survey (SOEP v28, 2012. doi:[10.5684/soep.v28](https://doi.org/10.5684/soep.v28)) throughout the tutorial.

Keywords Copula · Non-random sample selection · Penalized regression spline · Selection bias · Count response · Tutorial

1 Introduction

The sample selection model was introduced by [Gronau \(1974\)](#), [Lewis \(1974\)](#) and [Heckman \(1976\)](#) to deal with the situation in which the observations available for statistical analysis are not from a random sample of the population, and discussed by [Heckman \(1990\)](#) among others. This issue occurs when individuals select themselves into (or out of) the sample based on a combination of observed and unobserved characteristics. Estimates based on models that ignore such a non-random selection

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may be biased and inconsistent. This situation can be rectified using sample selection models, which typically consist of a two-equation system: a binary selection equation determining whether a particular statistical unit will be available in the outcome equation.

Let us consider a case study which uses data from the German Socio-Economic Panel survey (SOEP v28 2012) which will be analyzed in more detail in Sect. 3. The aim of the study is to estimate the determinants of labor mobility as well as the average number of job changes. Non-random selection arises if the sample consisting of individuals who are in the job market differ in important characteristics from the sample of individuals who are not part of the market. If the link between the decision to be part of the market and number of direct job changes (i.e., changes without an intervening spell of unemployment) is through observables, then selection bias can be avoided by accounting for these variables. However, if the link is also through unobservables then inconsistent parameter estimates are produced when using classic univariate modeling approaches. There are several other aspects that may complicate modeling labor mobility. Variables such as employment and length of education in years may have non-linear impacts on decision of being part of the job market and on the number of direct job changes, possibly due to productivity and life-cycle effects; imposing a priori a linear relationship (or non-linear by simply using quadratic polynomials, for example) could mean failing to capture interesting complex relationships. In addition, the assumption of bivariate normality employed in many sample selection models between, in this case, decision to be part of the job market and number of job changes may be too restrictive for applied work and it is typically made for mathematical convenience. Finally, the outcome of interest is a count variable.

The literature on sample selection models is vast and many variants of such models have been proposed; without claim of exhaustiveness here we mention some works. Chib et al. (2009) and Wiesenfarth and Kneib (2010) introduced two estimation methods to deal with non-linear covariate effects. Specifically, the approach of the former authors is based on Markov chain Monte Carlo simulation techniques and uses a simultaneous equation system that incorporates Bayesian versions of penalized smoothing splines. The latter further extended this approach by introducing a Bayesian algorithm based on low rank penalized B-splines for non-linear and varying-coefficient effects and Markov random-field priors for spatial effects. Marra and Radice (2013) proposed a frequentist counterpart which is computationally fast. Greene (1997), Terza (1998), Winkelmann (1998), Miranda (2004), Miranda and Rabe-Hesketh (2006) and Bratti and Miranda (2011) discuss fully parametric methods for estimating count data models allowing for overdispersion. These are based on the Poisson distribution and normally distributed unobserved heterogeneity. The common and limiting factor of these approaches is the assumption of bivariate normality.

Various methods that relax the assumption of normality have been proposed over the years; these include semiparametric (e.g., Gallant and Nychka 1987; Lee 1994; Powell 1994; Newey 2009) and nonparametric methods (e.g., Das et al. 2003; Lee 2008; Chen and Zhou 2010). Another way to relax the normality assumption is to use non-Gaussian parametric distributions. Recently, Marchenko and Genton (2012) and Ding (2014) extended the sample selection model to deal with heavy-tailedness by using the bivariate Student-t distribution. Another example of non-Gaussian paramet-

ric approach is copula modeling which allows for a great deal of flexibility in specifying the joint and marginal distributions of the selection and outcome equations (e.g., [Smith 2003](#); [Prieger 2002](#); [Hasebe and Vijverberg 2012](#); [Schwiebert 2013](#)). In the context of count responses, [Marra and Wyszynski \(2016\)](#) propose a copula-based approach, where covariates can be modeled flexibly using splines. There are advantages and disadvantages to both approaches (semi/non-parametric and parametric). The strongest point of the semi/non-parametric approach is the property of maintaining consistency of such estimators even disposing, in part or altogether, of distributional assumptions. However, semi/non-parametric methods are usually restricted when it comes to including a large set of covariates in the model and the resulting estimates are inefficient relatively to fully parametrized models (e.g., [Bhat and Eluru 2009](#)). To date, packages implementing semi/non-parametric procedures are CPU-intensive and the set of options provided is often quite limited. As for the parametric approach, many scholars agree upon its greater computational feasibility as compared to semi/non-parametric approaches, which allows for the use of familiar tools such as maximum likelihood without requiring simulation methods or numerical integration. As pointed out by [Smith \(2003\)](#), maximum likelihood techniques allow for the simultaneous estimation of all model parameters, and such methods, if the usual regularity conditions hold and the model is correctly specified, ensure consistent, efficient and asymptotically normal estimators. While a fully parametric copula approach is less flexible than semi/non-parametric approaches, it still allows the user to assess the sensitivity of results to different modeling assumptions. Specifically, the wide selection of potential copulae allows the modeler to perform sensitivity analysis to assess changes in results.

Some of the methods described above are implemented in popular software packages like SAS ([SAS Institute Inc 2013](#)), Stata ([StataCorp 2011](#)), LIMDEP ([Greene 2007](#)), EViews ([IHS Global Inc. 2015](#)) and R ([R Development Core Team 2016](#)). For example, the conventional Heckman sample selection model can be fitted in SAS using the **proc qlim** procedure and in Stata using **heckman** statement. The non-parametric method by [Lee \(2008\)](#) can be employed using the Stata package **leebounds** and the bivariate Student-t distribution Heckman model using **heckt**. The Poisson count data model by [Miranda and Rabe-Hesketh \(2006\)](#) can be employed in **Stata** using **ssm**. In R the sample selection packages are **sampleSelection** ([Toomet and Henningsen 2008](#)), **bayesSampleSelection** ([Wiesenfarth and Kneib 2010](#)), **ssmrob** ([Zhelonkin et al. 2013](#)) and **SemiParBIVProbit** ([Marra and Radice 2015](#)). **sampleSelection** and **bayesSampleSelection** make the assumption of bivariate normality between the model equations. **sampleSelection** and **ssmrob** assume a priori that continuous regressors have linear or pre-specified non-linear relationships to the responses, whereas **ssmrob** relaxes the assumption of bivariate normality by providing a robust two-stage estimator of Heckman's approach. **sampleSelection** and **SemiParBIVProbit** support binary responses for the outcome equation, with the latter allowing for non-linear covariate effects and non-Gaussian bivariate distributions. The R package **SemiParSampleSel** ([Marra et al. 2017b](#); [Wojtyś et al. 2016](#)) deals simultaneously with non-random sample selection, non-linear covariate effects and non-normal bivariate distribution between the model equations. Covariate-response relationships are flexibly modeled using a spline approach, whereas non-normal distributions are dealt with by using copula

functions. Note that `copulaSampleSel()` in **GJRM** (Marra and Radice 2017) works as `SemiParSampleSel()` in **SemiParSampleSel**.

In this paper, we further extend the **SemiParSampleSel** package (v. 1.5) by allowing the outcome to be modeled as a discrete random variable, and by allowing the mean, the higher order moments and the copula dependence parameter to be heterogeneous by specifying flexible linear predictor equations for each of them. Our approach of allowing multiple parameters to vary by observation follows the same rationale provided by Rigby and Stasinopoulos (2005), who extended generalized additive models to the context of more complex response distributions. As suggested by Marra and Wyszynski (2016), we expand the number of available outcome distributions; these include, for instance, beta binomial and zero inflated margins (see “Appendix 1”).

The rest of the paper is organized as follows: in Sect. 2, we provide a brief theoretical overview of the sample selection modeling approach for count data and its properties (this is based on: Marra and Wyszynski 2016). Section 3 presents **SemiParSampleSel** by describing its main infrastructure and how to use the package to obtain interpretable statistical models. Using **SemiParSampleSel**, we provide a step-by-step illustration on how to use sample selection models for count data in R to build a prediction model for SOEP data. A summary of the paper is given in Sect. 4.

2 An overview of sample selection models for count data

2.1 Model definition

In the sample selection problem, our aim is to fit a regression model when some observations of the outcome variable, Y_{2i} for $i = 1, \dots, n$, are missing not at random. We will use a latent continuous variable Y_{1i}^* such that $Y_{1i} = \mathbf{1}(Y_{1i}^* > 0)$, where $\mathbf{1}$ is the indicator function and Y_{1i} governs whether or not an observation on the variable of primary interest is generated. We assume normality for Y_{1i}^* and a discrete distribution, \mathcal{F} , (see “Appendix 1” for all possible choices and Fig. 1 for illustration) for Y_{2i} . That is, $Y_{1i}^* \sim \mathcal{N}(\mu_{1i}, 1)$ (which yields a probit model for Y_{1i}) and $Y_{2i} \sim \mathcal{F}(\mu_{2i}, \sigma_i, \nu_i)$, where μ_{1i}, μ_{2i} are location parameters; σ_i is the scale parameter and ν_i is the shape parameter. Note that we are considering parametrization for the most generic case of \mathcal{F} , although other parametrizations such as $\mathcal{F}(\mu_{2i})$ and $\mathcal{F}(\mu_{2i}, \sigma_i)$ are possible.

We can represent the random sample using a pair of variables (Y_{1i}, Y_{2i}) . Let F denote the joint cumulative distribution function (cdf) of (Y_{1i}, Y_{2i}) and let F_1 and F_2 be the marginal cdfs pertaining to Y_{1i} and Y_{2i} , respectively. The model is then defined by using the copula representation (Sklar 1959)

$$F(y_{1i}, y_{2i}) = C(F_1(y_{1i}), F_2(y_{2i}); \theta_i),$$

for some two-place function C , where θ_i is an association parameter measuring the dependence between the two marginal cdfs. For details on binding continuous and discrete margins see Marra and Wyszynski (2016). In the presented package, the families currently implemented are normal, Clayton, Joe, Frank, Gumbel, Farlie–Gumbel–Morgenstern (FGM), and Ali–Mikhail–Haq (AMH; for examples of copulae,

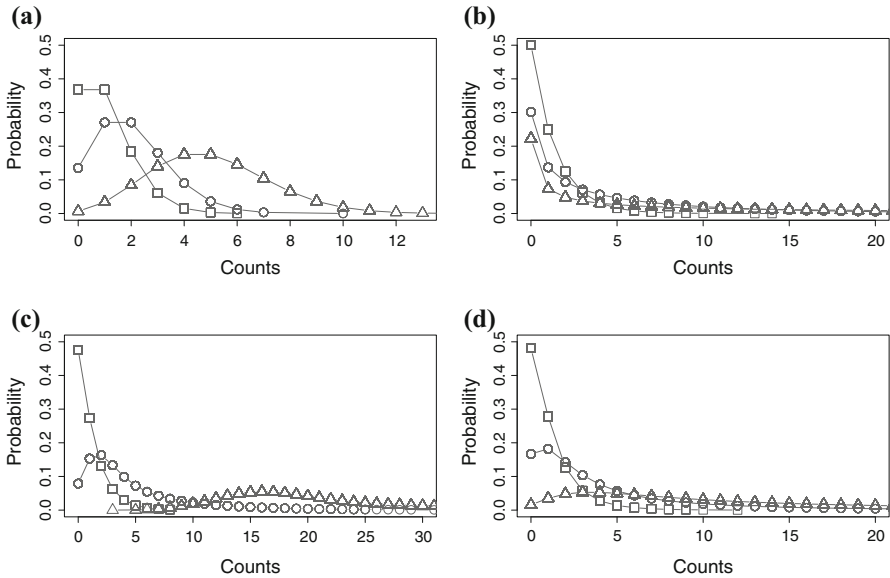


Fig. 1 Probability mass functions of the Poisson (a), negative binomial (b), Delaporte (c) and Poisson inverse Gaussian (d) distribution. The parameter values have been chosen arbitrarily to show different shapes of the distributions. For Poisson, μ is 1, 2 and 5, represented as rectangular, circular and triangular lines respectively. Similarly, for negative binomial and Poisson inverse Gaussian, μ and σ are (1, 1), (5, 2) and (30, 3). For Delaporte, μ , σ and ν are (1, 0.1), (5, 2, 0.3) and (30, 3, 0.5). Note that Delaporte can have thinner or thicker tails depending on the choice of parameters. At the same time, the tails of Poisson are thinner than those of Delaporte (see: [Marra and Wyszynski 2016](#))

see Table 6 in “Appendix 1”). Rotations by 90, 180 and 270 degrees for Clayton, Joe and Gumbel can be obtained (see also Fig. 2 [Brechmann and Schepsmeier 2013](#)). Despite the fact that θ cannot be interpreted directly, it can be transformed into Kendall’s τ ranging on the interval $[-1, 1]$ yielding a general interpretation for all copulae (for a discussion on Kendall’s τ for continuous and discrete margins see: [Genest and Neslehova 2007](#); [Marra and Wyszynski 2016](#)).

The log-likelihood function for the sample selection model can be generically expressed as a product over two disjoint subsets of the sample: one for the observations with a missing value of the response of interest and the other for the remaining observations ([Smith 2003](#)). In the first case, the likelihood takes the simple form of $\Pr(Y_1 = 0)$, which is equivalent to $F_1(0)$. In the second case, the joint likelihood can be expressed, using the multiplication rule, as $P(Y_2 = y_2, Y_1 = 1)$. We dropped the observation index i to avoid clutter:

$$\begin{aligned}
 L &= \prod_0 \Pr(Y_1 = 0) \prod_1 P(Y_2 = y_2, Y_1 = 1) \\
 &= \prod_0 \Pr(Y_1^* \leq 0) \prod_1 f_{2|1}(y_2|y_1^* > 0) \Pr(Y_1^* > 0).
 \end{aligned}$$

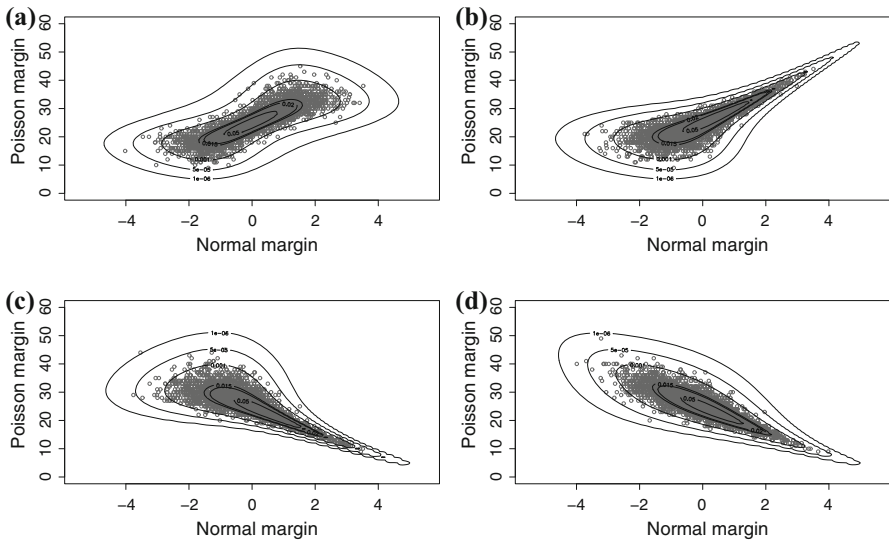


Fig. 2 Contour plots of Frank (a), Joe (b), Clayton 90 (c) and Gumbel 270 (d) copulae. 5000 deviates were generated for each copula. The first margin is Poisson, whereas the second is standard normal. Kendall’s τ was set to 0.7. The plots have different shapes depending on the copula. For instance, Joe copula shows greater tail dependence in *upper right corner*, whereas Clayton 90 shows greater tail dependence in *lower right corner* (see: Marra and Wyszynski 2016)

Note that for the continuous response

$$\begin{aligned}
 f_{2|1}(y_2|y_1^* > 0) &= \frac{\partial}{\partial y_2} \frac{F_2(y_2) - F(0, y_2)}{F_1(1)} \\
 &= \frac{1}{1 - F_1(0)} \frac{\partial}{\partial y_2} (F_2(y_2) - F(0, y_2)) \\
 &= \frac{1}{1 - F_1(0)} (f_2(y_2) - \frac{\partial}{\partial y_2} F(0, y_2)), \tag{1}
 \end{aligned}$$

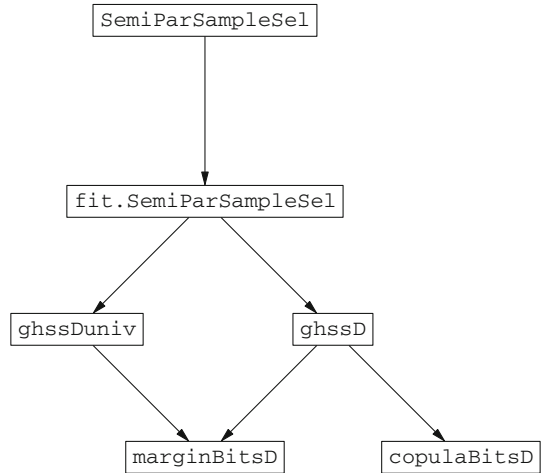
and the log-likelihood will be

$$\ell = \sum_0 \log F_1(0) + \sum_1 \log (f_2(y_2) - \frac{\partial}{\partial y_2} F(0, y_2)).$$

In the third line of (1), for the discrete y_2 , the derivative cannot be computed, since $F_2(y_2)$ is discontinuous on the integers in its domain. We will calculate the derivative with respect to y_2 using finite differences (Nikoloulopoulos and Karlis 2009)

$$\begin{aligned}
 f_{2|1}(y_2|y_1^* > 0) &= \frac{1}{1 - F_1(0)} \{F_2(y_2) - F_2(y_2 - 1)\} - \frac{1}{1 - F_1(0)} \{F(0, y_2) - F(0, y_2 - 1)\} \\
 &= \frac{1}{1 - F_1(0)} [\{F_2(y_2) - F_2(y_2 - 1)\} - \{F(0, y_2) - F(0, y_2 - 1)\}]
 \end{aligned}$$

Fig. 3 Modular relations between functions in **SemiParSampleSel** R package. The *arrows* indicate the direction in which the functions are called. For instance, `SemiParSampleSel` calls `fit.SemiParSampleSel`



$$\begin{aligned}
 &= \frac{1}{1 - F_1(0)} \{F_2(y_2) - F_2(y_2 - 1) - F(0, y_2) + F(0, y_2 - 1)\} \\
 &= \frac{1}{1 - F_1(0)} \{f_2(y_2) - F(0, y_2) + F(0, y_2 - 1)\}.
 \end{aligned}$$

The model log-likelihood will be given by (Marra and Wyszynski 2016)

$$\ell = \sum_0 \log F_1(0) + \sum_1 \log(f_2(y_2) - F(0, y_2) + F(0, y_2 - 1)). \tag{2}$$

Figure 3 illustrates how the main **SemiParSampleSel** function decomposes the likelihood in (2) in a modular fashion. The probability mass function, $f_2(y_2)$, and cumulative distribution functions, $F_1(0)$, $F_2(y_1)$ and $F_2(y_1 - 1)$, and their derivatives are obtained by `marginBitsD`. The copulae, $F(0, y_2)$ and $F(0, y_2 - 1)$, and their derivatives with respect to cumulative distribution functions are computed with `copulaBitsD`. The `ghssD` function makes use of `marginBitsD` and `copulaBitsD` and constructs an object encompassing the likelihood of the sample selection model and its first and second derivatives. The univariate counterpart of the sample selection model, `ghssDuniv`, likewise utilizes information provided by `marginBitsD`. The parameter estimation is conducted by `fit.SemiParSampleSel` and both inputs and outputs are processed by `SemiParSampleSel`. Hence, one can implement potentially any margin or copula with great ease without changing the general structure of the likelihood. For instance, let us denote the parameters associated with the outcome linear predictor as δ_2 . The general expression of the first derivative of (2) with respect to δ_2 is

$$\frac{\partial \ell}{\partial \delta_2} = \sum_1 \frac{1}{f_2(y_2) - F(0, y_2) + F(0, y_2 - 1)}$$

$$\times \left(\frac{\partial f_2(y_2)}{\partial \delta_2} - \frac{\partial F(0, y_2)}{\partial F_2(y_2)} \frac{\partial F_2(y_2)}{\partial \delta_2} + \frac{\partial F(0, y_2 - 1)}{\partial F_2(y_2 - 1)} \frac{\partial F_2(y_2 - 1)}{\partial \delta_2} \right), \quad (3)$$

In this case, `copulaBitsD` will evaluate two derivatives and three derivatives will be processed by `marginBitsD`. The entire expression will be obtained via `ghssD`. Hence, implementing an additional copula or margin requires solely changes in either `copulaBitsD` or `marginBitsD`.

2.2 Linear predictor specification

We assume that the mean, scale, shape and copula parameters, μ_{1i} , μ_{2i} , σ_i , ν_i and θ_i , are linked with the predictors η_{vi} , $v = 1, \dots, 5$, i.e., $\mu_{1i} = \eta_{1i}$, $g_\mu(\mu_{2i}) = \eta_{2i}$, $g_\sigma(\sigma_i) = \eta_{3i}$, $g_\nu(\nu_i) = \eta_{4i}$ and $g_\theta(\theta_i) = \eta_{5i}$, where the link functions g depend on the distributions of y_{2i} and on the copula functions (see ‘‘Appendix 1’’). For simplicity, and without loss of generality, we suppress the v subscript and define the generic linear predictor as

$$\eta_i = \mathbf{u}_i^\top \boldsymbol{\alpha} + \sum_{k=1}^K s_k(z_{ki}), \quad i = 1, \dots, n,$$

where vector $\mathbf{u}_i^\top = (1, u_{2i}, \dots, u_{pi})$ is the i^{th} row of $\mathbf{U} = (\mathbf{u}_1, \dots, \mathbf{u}_n)^\top$, the $n \times P$ model matrix containing P parametric model components (e.g., intercept, dummy and categorical variables), $\boldsymbol{\alpha}$ is a parameter vector, and the s_k are unknown smooth functions of the K continuous covariates z_{ki} . The smooth functions are subject to the centering (identifiability) constraint $\sum_i s_k(z_{ki}) = 0$, $k = 1, \dots, K$ (Wood 2017).

The smooth functions are represented using regression splines, where, in the one-dimensional case, a generic $s_k(z_{ki})$ is approximated by a linear combination of known spline basis functions, $b_{kj}(z_{ki})$, and regression parameters, β_{kj} , i.e., $s_k(z_{ki}) = \sum_{j=1}^{J_k} \beta_{kj} b_{kj}(z_{ki}) = \boldsymbol{\beta}_k^\top \mathbf{B}_k(z_{ki})$, where J_k is the number of spline bases used to represent s_k , $\mathbf{B}_k(z_{ki})$ is the i^{th} vector of dimension J_k containing the basis functions evaluated at the observation z_{ki} , i.e., $\mathbf{B}_k(z_{ki}) = \{b_{k1}(z_{ki}), b_{k2}(z_{ki}), \dots, b_{kJ_k}(z_{ki})\}^\top$, and $\boldsymbol{\beta}_k$ is the corresponding parameter vector. Calculating $\mathbf{B}_k(z_{ki})$ for each i yields J_k curves (encompassing different degrees of complexity) which multiplied by some real valued parameter vector $\boldsymbol{\beta}_k$ and then summed will give a (linear or non-linear) estimate for $s_k(z_k)$ (see, for instance, Marra and Radice (2010) for a more detailed overview). Basis functions should be chosen to have convenient mathematical and numerical properties. B-splines, cubic regression and low rank thin plate regression splines are supported in our implementation (see Wood (2017) for full details on these spline bases). Our implementation also supports varying coefficients’ models, obtained by multiplying one or more smooth terms by some predictor(s), smooth functions of two or more (e.g., spatial) covariates, random effect and Markov random field smooth functions, to name but a few (Wood 2017). These cases follow a similar construction as described above. See, for instance, Wood (2017); Marra et al. (2017a).

In principle, the parameters of the sample selection model are identified even if the same regressors appear in both linear predictors (e.g., Wiesenfarth and Kneib 2010).

However, better estimation results are generally obtained when the set of regressors in the selection equation contains at least one or more regressors (usually known as exclusion restrictions) that are not included in the outcome equation (e.g., [Marra and Radice 2013](#)).

2.3 Estimation approach

Unpenalized estimation can result in smooth term estimates that are too rough and overfitting (e.g., [Wood 2017](#)). This issue is dealt with by using a roughness penalty term ([Ruppert et al. 2003](#); [Wood 2017](#)). Denote the log-likelihood function as $\ell(\delta)$, where $\delta^T = (\delta_1^T, \dots, \delta_5^T)$. Note that this parameter vector's definition is the most generic given the marginal distributions considered in this paper (this is consistent with zero inflated negative binomial, for instance). For outcome margins parametrized only in terms of μ (e.g., Poisson) we would have $\delta^T = (\delta_1^T, \dots, \delta_3^T)$, whereas for distributions parametrized in terms of both μ and σ (e.g. beta binomial) we have $\delta^T = (\delta_1^T, \dots, \delta_4^T)$. For each smooth $s_{v_{k_v}}(z_{v_{k_v}})$ we have a penalty such that $\beta^T_{v_{k_v}} \mathbf{S}_{v_{k_v}} \beta_{v_{k_v}}$, where $\mathbf{S}_{v_{k_v}}$ is a positive semi-definite penalty matrix with known coefficients. The quadratic expression is used, since it measures the second-order roughness of the smooth terms in the model. The form of the penalty $\mathbf{S}_{v_{k_v}}$ will depend on the selected spline basis ([Wood 2017](#)). The function to maximize is

$$\ell_p(\delta) = \ell(\delta) - \frac{1}{2} \delta^T \mathbf{S}_\lambda \delta. \tag{4}$$

where $\mathbf{S}_\lambda = \text{diag}(\mathbf{0}^T_{P_1}, \lambda_{1K_1} \mathbf{S}_{1K_1}, \dots, \lambda_{1K_1} \mathbf{S}_{1K_1}, \mathbf{0}^T_{P_2}, \lambda_{2K_2} \mathbf{S}_{2K_2}, \dots, \lambda_{2K_2} \mathbf{S}_{2K_2}, \mathbf{0}^T_{P_3}, \lambda_{3K_3} \mathbf{S}_{3K_3}, \dots, \lambda_{3K_3} \mathbf{S}_{3K_3}, \mathbf{0}^T_{P_4}, \lambda_{4K_4} \mathbf{S}_{4K_4}, \dots, \lambda_{4K_4} \mathbf{S}_{4K_4}, \mathbf{0}^T_{P_5}, \lambda_{5K_5} \mathbf{S}_{5K_5}, \dots, \lambda_{5K_5} \mathbf{S}_{5K_5})$ in the most generic case; $\lambda_{v_{k_v}}$ represents smoothing parameters which control for the trade-off between fit and smoothness.

The estimation algorithm is structured in two main steps which are iterated until convergence:

Step.1 For a given vector $\delta^{[a]}$, and maintaining smoothing parameter fixed at $\lambda^{[a]}$, find the new iterate for δ using trust region approach ([Chapter4 Nocedal and Wright 2006](#)):

$$\min_{\mathbf{p}} \check{\ell}_p(\delta^{[a]}) \stackrel{\text{def}}{=} - \left\{ \ell_p(\delta^{[a]}) + \mathbf{p}^T (\mathbf{g}^{[a]} - \mathbf{S}_\lambda \hat{\delta}^{[a]}) + \frac{1}{2} \mathbf{p}^T (\mathcal{H}^{[a]} - \mathbf{S}_\lambda) \mathbf{p} \right\} \text{ so that } \|\mathbf{p}\| \leq r^{[a]},$$

$$\delta^{[a+1]} = \arg \min_{\mathbf{p}} \check{\ell}_p(\delta^{[a]}) + \delta^{[a]},$$

where $\|\cdot\|$ denotes the Euclidean norm and $r^{[a]}$ represents the radius of the trust region. After dropping the iteration index, the score vector \mathbf{g} is defined by five subvectors $\mathbf{g}_o = \partial \ell(\delta) / \partial \delta_o$ for $o = 1, \dots, 5$, while the Hessian matrix has a 5×5 matrix block structure with $(r, h)^{th}$ element $\mathcal{H}_{r,h} = \partial^2 \ell(\delta) / \partial \delta_r \partial \delta_h^T$, $r, h = 1, \dots, 5$. At each iteration of the algorithm, $\check{\ell}_p(\delta^{[a]})$ is minimized subject to the constraint that the solution falls within a trust region with radius

$r^{[a]}$. The proposed solution is then accepted or rejected and the trust region expanded or shrunken based on the ratio between the improvement in the objective function when going from $\delta^{[a]}$ to $\delta^{[a+1]}$ and that predicted by the quadratic approximation. Note that, near the solution, the trust region Newton algorithm typically behaves as a Newton algorithm (Chapter 4 Nocedal and Wright 2006).

Step 2 For a given smoothing parameter vector value $\lambda^{[a]}$, and maintaining $\delta^{[a+1]}$ fixed, find an estimate of λ :

$$\text{minimize } \frac{1}{n^*} \|\sqrt{\mathbf{W}}(\mathbf{z} - \mathbf{X}\tilde{\delta})\|^2 - 1 + \frac{2}{n^*} \text{tr}(\mathbf{A}_\lambda) \quad \text{w.r.t. } \lambda, \quad (5)$$

where $\sqrt{\mathbf{W}}$ is a weight non-diagonal matrix square root, \mathbf{z} is the pseudo-data vector consisting of 5-dimensional subvectors \mathbf{z}_i defined as $\mathbf{z}_i = \mathbf{X}_i \delta^{[a]} + \mathbf{W}_i^{-1} \mathbf{d}_i$, $\mathbf{d}_i = \{\partial \ell(\delta)_i / \partial \eta_{1i}, \dots, \partial \ell(\delta)_i / \partial \eta_{5i}\}^\top$, \mathbf{W}_i is a 5×5 matrix with $(r, h)^{th}$ element $(\mathbf{W}_i)_{rh} = -\partial^2 \ell(\delta)_i / \partial \eta_{ri} \partial \eta_{hi}$, $r, h = 1, \dots, 5$, $\mathbf{X}_i = \text{diag} \left\{ \left(\mathbf{u}_{1i}^\top, \mathbf{B}_{1i}^\top \right), \dots, \left(\mathbf{u}_{5i}^\top, \mathbf{B}_{5i}^\top \right) \right\}$, $\tilde{\delta} = (\mathbf{X}^\top \mathbf{W} \mathbf{X} + \mathbf{S}_\lambda)^{-1} \mathbf{X}^\top \mathbf{W} \mathbf{z}$, $n^* = 5n$, $\sqrt{\mathbf{W}} \mathbf{A} = \sqrt{\mathbf{W}} \mathbf{X} (\mathbf{X}^\top \mathbf{W} \mathbf{X} + \mathbf{S}_\lambda)^{-1} \mathbf{X}^\top \mathbf{W}$ is the hat matrix, and $\text{tr}(\mathbf{A}_\lambda)$ the estimated degrees of freedom (*edf*) of the penalized model (e.g., Marra and Wyszynski 2016). The iteration index has been dropped to avoid clutter. We will use the approach by Wood (2004) to perform minimization of (5), which turns out to be computationally efficient and stable.

2.4 Confidence intervals, variable and model selection

Inferential theory for penalized estimators is complicated by the presence of smoothing penalties which undermines the usefulness of classic frequentist results for practical modeling. As shown in Marra and Radice (2013), reliable pointwise confidence intervals for the terms of a regression spline sample selection model can be constructed using

$$\delta | \mathbf{z} \rightsquigarrow \mathcal{N}(\hat{\delta}, \mathbf{V}_\delta), \quad (6)$$

where $\hat{\delta}$ is an estimate of δ and $\mathbf{V}_\delta = (-\mathcal{H} + \mathbf{S}_\lambda)^{-1}$. The structure of \mathbf{V}_δ is such that it includes both a bias and variance component in a frequentist sense, which is why such intervals exhibit close to nominal coverage probabilities (Marra and Wood 2012). Given (6), confidence intervals for linear and non-linear functions of the model parameters can be easily obtained. For instance, for a generic $\hat{s}_k(z_{ki})$ these can be obtained using

$$\hat{s}_k(z_{ki}) \rightsquigarrow \mathcal{N}(s_k(z_{ki}), \mathbf{B}_k(z_{ki})^\top \mathbf{V}_{\delta_k} \mathbf{B}_k(z_{ki})),$$

where \mathbf{V}_{δ_k} is the submatrix of \mathbf{V}_δ corresponding to the regression spline parameters associated with k^{th} function. Intervals for non-linear functions of the estimated model

coefficients can be conveniently obtained by simulation from the posterior distribution of δ . As for the parametric model components, using (6) is equivalent to using classic likelihood results because such terms are not penalized. Intervals for average values of σ , ν and θ can be obtained by simulating from the posterior distribution of δ as follows:

1. Draw n_{sim} random vectors from (6).
2. Calculate n_{sim} simulated realizations of the function of interest. For example, since $g_{\sigma}(\sigma_i) = \eta_{3i}$, $\sigma_i^{sim} = (\sigma_{1i}^{sim}, \sigma_{2i}^{sim}, \dots, \sigma_{n_{sim}i}^{sim})$ where $\sigma_{oi}^{sim} = g_{\sigma_i}^{-1}(\eta_{3oi}^{sim})$, $o = 1, 2, \dots, n_{sim}$. For each o , obtain the mean value of σ_{oi}^{sim} across all observations.
3. Using σ^{sim} , calculate the lower, $\xi/2$, and upper, $1 - \xi/2$, quantiles. For 95 per cent Bayesian confidence intervals, ξ is set to 0.05.

Copula models with a single dependence parameter can be thought of as non-nested models. As suggested by Zimmer and Trivedi (2006) among others, one approach for choosing between copula models is to use either the Akaike or (Schwarz) Bayesian information criterion (*AIC* and *BIC*, respectively). In our case, $AIC = -2\ell(\hat{\delta}) + 2edf$ and $BIC = -2\ell(\hat{\delta}) + \log(n)edf$, where the log-likelihood is evaluated at the penalized parameter estimates and $edf = \text{tr}(\hat{A}_{\hat{\delta}})$. Other model selection methods include Vuong and Clarke test (Vuong 1989; Clarke 2007), which will be illustrated in the next section.

3 The package SemiParSampleSel

As pointed out above, the R package **SemiParSampleSel** implements model-based penalized maximum likelihood estimation that results in interpretable models. As mentioned in Sect. 2.1, the **SemiParSampleSel** package offers a modular nature that allows to specify a wide range of non-random selection models. The non-random sample selection model is specified as the combination of a distributional assumption and structural assumptions. Unlike in the case of Greene (1997), the distributional assumption specifies the conditional distribution of the outcome. The structural assumption specifies the types of effects that are to be used in the model, i.e., it represents the deterministic structure of the model. Usually, it specifies how the predictors are related to the conditional mean, variance and shape of the outcome variable and how the selection mechanism and the outcome are related.

The function `SemiParSampleSel()` provides an interface to fit sample selection models for continuous and count data. Before we show how one can use the function to fit this model to estimate the determinants of labor mobility, we give a short overview on the function:

```
SemiParSampleSel(formula, data = list(), BivD = 'N',
                  margins = c('probit', 'NB'), \ldots)
```

The model is specified using a `formula` which is a list of five formulas, one for the mean of the selection equation, three for the mean, variance and shape of the outcome equation and one for the copula parameter. These are `glm()` like formulas except that smooth terms can be included in the equations as for `gam()` in **mgcv**. For instance, the formulas for the mean of the selection equation may look like:

```
y.sel ~ as.factor(x1) + s(x2, bs = ``cr``, k = 10, m = 2) + ...
```

and that for the scale parameter of the outcome equation:

```
~ as.factor(x1) + s(x3, x4) + \cdots
```

where `y.sel` represents the binary selection variable, `x1` is a categorical predictor, and the `s()` terms are used to specify smooth functions of the continuous predictors `x2`, `x3` and `x4`. Argument `bs` specifies the spline basis; possible choices include `cr` (cubic regression spline), `cs` (shrinkage version of `cr`), `tp` (thin plate regression spline) and `ts` (shrinkage version of `tp`). Bivariate smoothing, e.g., `s(x3, x4)`, is achieved using `bs = "tp"`. `k` is the basis dimension (default is 10) and `m` the order of the penalty (default is 2). More details and options on smooth term specification can be found in the documentation of **mgcv**. **SemiParSampleSel** does not currently support the use of tensor product smooths. The data set is provided as a `data.frame` via the `data` argument. The type of bivariate copula linking the selection and outcome equations can be specified through `BivD`. Possible choices are "N", "C0", "J0", "FGM", "F", "AMH" and "G0" which stand for bivariate normal, Clayton, Joe, Farlie-Gumbel-Morgenstern, Frank, Ali-Mikhail-Haq and Gumbel. Rotated versions (90, 180 and 270 degrees) of the asymmetric copulae are also implemented ("C90", "C180", "C270", "J90", ...). For more details on available copulae see "Appendix 1". The argument `margins` specifies the marginal distributions of the selection and outcome equations, given in the form of a two-dimensional vector which is equal to `c("probit", "NB")` for normal and negative binomial margins. The first margin currently admits only the normal distribution ("probit"). The second margin can also be "N", "GA", "NB", "D", "PIG" or "S" which stand for normal, gamma, negative binomial, Delaporte, Poisson inverse Gaussian and Sichel (see "Appendix "). Details on all the other arguments, including starting value and control options, and the fitted-object list that the function returns can be found in [Marra et al. \(2017b\)](#). See also data simulation function in Appendix 2, for example.

Apart from `SemiParSampleSel()`, the package also encompasses several other functions, which offer some numerical and graphical interpretations. These will be presented in the next section:

- `resp.check(y, margin = "P")`. This function preliminarily checks whether the (non-missing) response follows one of the discrete distributions. This is done by generating a kernel density and normalized and randomized QQ-plot (for details see: [Stasinopoulos and Rigby 2007](#)). Distributions that require a binomial denominator (e.g., binomial) need to specify `bd` as an argument.
- `conv.check()` provides some information about the convergence of the algorithm.
- `summary()`. The summary function of `SemiParSampleSel()`; analogical to the one in `glm` and `gam`.
- `AIC()` and `BIC()` return Akaike and Bayesian information criteria.
- `VuongClarke()` performs Vuong and Clarke test for comparing two competing models. For example, these can be models that differ on the choice of copula.
- `aver()` provides the average predicted values for the entire data set. When `univariate = TRUE`, the average prediction for the univariate model ignoring non-random sample selection is returned.

Table 1 Variables of SOEP data

Name	Description
selection	Binary selection variable; 1 if the individual is part of the job market; 0 otherwise.
EXPFT	Full-time employment in years.
Single	Binary variable; 1 if the individual is single; 0 otherwise.
WhiteCollar	Binary variable; 1 if the individual is a white collar worker; 0 otherwise.
LEduc	Length of education in years.
SPDSup	Binary variable; 1 if the individual is a strong or a very strong supporter of the Socialdemocratic Party of Germany; 0 otherwise.

- `plot()` function illustrates smooth terms when they are specified for a given equation. For instance, `eq = 1` will yield a smooth function for the selection equation of the model.
- `post.check()` takes the response vector and produces a QQ-plot based on the estimates obtained from the sample selection model. Similarly to `resp.check()`, the residuals are normalized and randomized. Distributions that require a binomial denominator need to specify `bd` as an argument.

3.1 Case study: an application to labor mobility

The data is from the German Socio-Economic Panel survey of 1984 (SOEP v28 2012) and the aim of the study is to estimate the determinants of labor mobility in the presence of non-random sample selection. As mentioned in the introduction, non-random selection arises if the sample consisting of individuals who are on the job market differ in important characteristics from the sample of individuals who are not part of the job market. To rectify this situation, we will employ sample selection models implemented in **SemiParSampleSel**.

We first load the package and the data set.

```
R> library(SemiParSampleSel)
R> data(SOEP)
```

The data set contains 2,651 observations and the outcome variable is the number of direct job changes (DJC, i.e. changes without an intervening spell of unemployment). Other available variables are presented in Table 1.

Following the work by Winkelmann (1998), the linear predictors for the selection and outcome equations, respectively, can be specified as follows

$$\eta_1 = \alpha_{10} + \alpha_{11}\text{Single} + \alpha_{12}\text{WhiteCollar} + s_{11}(\text{EXPFT})$$

$$\eta_2 = \alpha_{20} + \alpha_{21}\text{SPDSup} + \alpha_{22}\text{WhiteCollar} + s_{21}(\text{LEduc}),$$

where the non-linear specification of `LEduc` and `EXPFT` arises from the fact that these covariates embody productivity and life-cycle effects that are likely to affect the probability to be part of the job market non-linearly. These will be modeled using thin plate regression splines with 10 bases and penalties based on second order derivatives. We will employ linear predictors for the scale and association parameters, since `WhiteCollar` may have an impact on variability of the response. Thus,

$$\begin{aligned}\eta_3 &= \alpha_{30} + \alpha_{31}\text{WhiteCollar} \\ \eta_4 &= \alpha_{40} + \alpha_{41}\text{WhiteCollar},\end{aligned}$$

In R, the specification above implies determining formulas as

```
R> sel.eq <- selection ~ Single + WhiteCollar + s(EXPFT)
R> out.eq <- DJC ~ SPDSup + WhiteCollar + s(LEduc)
R> sigma.eq <- ~ WhiteCollar
R> theta.eq <- ~ WhiteCollar
```

Prior to fitting a count data sample selection model, a distribution for the outcome variable may be chosen by looking at the histogram of the response along with the estimated density from the assumed distribution, and at the randomized and normalized responses (Rigby and Stasinopoulos 2005). The latter will provide an approximate guide to the adequacy of the chosen distribution. These should behave approximately as normally distributed variables (even though the original observations are not). Note that this preliminary check has to be treated with caution, since the distribution of the outcome is assumed to be unconditional on any covariate effects.

```
R> resp.check(SOEP$DJC, margin = ``P``)
R> resp.check(SOEP$DJC, margin = ``NB``)
R> resp.check(SOEP$DJC, margin = ``D``)
R> resp.check(SOEP$DJC, margin = ``PIG``)
R> resp.check(SOEP$DJC, margin = ``S``)
```

These plots (reported in Fig. 4; for the remainder see Appendix 3) show that the density and QQ-plots are the most supported with “PIG” being the best. Based on this, we fit the count sample selection model with Poisson inverse Gaussian margin for the outcome variable using the function `SemiParSampleSel()` to the SOEP data. The normal copula is used to link the selection and the outcome equations. The reason for choosing normal copula is to preliminarily check the direction and magnitude of association between both equations.

```
R> fit1 <- SemiParSampleSel(list(sel.eq, out.eq, sigma.eq, theta.eq),
+                           data = SOEP, BivD = ``N``,
+                           margins = c(``N``, ``PIG``),
+                           iterlimsp = 50)
```

Before viewing the output, it is necessary to check the convergence of the algorithm. This can be done by submitting `conv.check()` command.

```
R> conv.check(fit1)
```

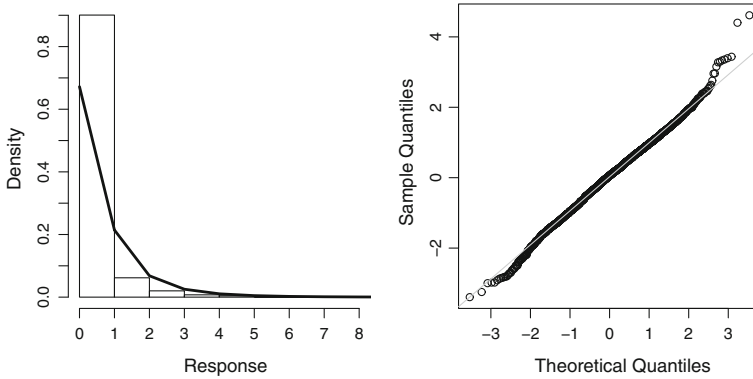


Fig. 4 Poisson inverse Gaussian kernel density and QQ-plot

Largest absolute gradient value: 8.984057e-05
 Observed information matrix is positive definite
 Eigenvalue range: [0.4848655,5477.451]

Trust region iterations before smoothing parameter estimation: 9
 Loops for smoothing parameter estimation: 5
 Trust region iterations within smoothing loops: 12

The algorithm converged successfully, since the gradient is equal to zero and the Hessian is positive definite. Note that the condition number is equal to $\frac{5477.451}{0.4848655} \approx 10^4$. The summary() function yields

```
R> summary(fit1)
```

ERRORS' DISTRIBUTION: Bivariate Normal

SELECTION EQ.

Family: Bernoulli

Link function: probit

Formula: sel ~ s(EXPFT) + Single + WhiteCollar

	Estimate	Std. Error	z value	Pr(> z)
(Intercept)	1.32760	0.04038	32.878	< 2e-16 ***
Single	0.05431	0.17878	0.304	0.761
WhiteCollar	0.58262	0.11849	4.917	8.79e-07 ***

 Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

Smooth components' approximate significance:

	edf	Ref.df	Chi.sq	p-value
s(EXPFT)	5.383	6.508	34.56	1.01e-05 ***

Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

OUTCOME EQ.

Family: Poisson inverse Gaussian

Link function: log

Formula: $l \sim \text{WhiteCollar} + s(\text{LEduc}) + \text{SPDSup}$

	Estimate	Std. Error	z value	Pr(> z)	
(Intercept)	-0.24748	0.10721	-2.308	0.02098	*
WhiteCollar	-0.39729	0.15376	-2.584	0.00977	**
SPDSup	0.10818	0.09063	1.194	0.23260	

Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

Smooth components' approximate significance:

	edf	Ref.df	Chi.sq	p-value	
s(LEduc)	2.955	3.604	21.01	0.000219	***

Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

EQUATION 3

Link function: log(sigma)

Formula: $\sigma \sim \text{WhiteCollar}$

	Estimate	Std. Error	z value	Pr(> z)	
(Intercept)	0.9093	0.2273	4.001	6.31e-05	***
WhiteCollar	-0.2048	0.3829	-0.535	0.593	

Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

EQUATION 4

Link function: atanh(theta)

Formula: $\theta \sim \text{WhiteCollar}$

	Estimate	Std. Error	z value	Pr(> z)	
(Intercept)	-1.0343	0.1968	-5.255	1.48e-07	***
WhiteCollar	0.1458	0.3879	0.376	0.707	

Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

n = 2651 n.sel = 2445 sigma = 2.378(1.358,3.97)
theta = -0.761(-0.898,-0.43) total edf = 18.338

The output shows all estimates pertaining to the selection and outcome equations, and the equations linked with the scale and association parameters. In particular, the white-collar effect on the selection is statistically significant and implies that

Table 2 Akaike and Bayesian information criteria for the Poisson inverse Gaussian models for SOEP data

Model	AIC	BIC	Model	AIC	BIC
Normal	6139.7	6247.5	Gumbel 90	6145.8	6252.6
Frank	6117.3	6223.4	Gumbel 270	6129.7	6236.1
Joe 90	6149.1	6256.1	Clayton 90	6116.9	6223.2
Joe 270	6116.9	6223.1	Clayton 270	6151.6	6258.9

white-collar workers are more likely to be active on the labour market than non-white-collar workers. In addition, the white-collar effect on the outcome is also statistically significant and suggests that being a white-collar worker decreases chances of changing jobs and hence labour mobility. At the bottom of the output one can see the mean values across all observations for $\hat{\sigma}$ and $\hat{\theta}$ together with the corresponding confidence intervals.

Let us estimate a Poisson inverse Gaussian Frank (`fit2`) and Joe 270 (`fit3`) models with the same specification as the normal Poisson inverse Gaussian. The Akaike and Bayesian information criteria to compare the competing models are obtained using `AIC()` and `BIC()` functions.

```
R> AIC(fit1, fit2, fit3)
      df      AIC
fit1 18.33783 6139.654
fit2 18.03288 6117.332
fit3 18.06156 6116.889
R> BIC(fit1, fit2, fit3)
      df      BIC
fit1 18.33783 6247.530
fit2 18.03288 6223.414
fit3 18.06156 6223.140
```

The left-hand side column indicates the estimated degrees of freedom, whereas the right-hand side denotes the criterion. Table 2 shows the AIC and BIC scores for the normal Poisson inverse Gaussian model and other Poisson inverse Gaussian models with different copulae. These are Frank and the 90 and 270 degrees rotated versions of Joe, Gumbel and Clayton. The reason for choosing these rotations is that the normal model initially indicated negative dependence of $\hat{\theta} = -0.761$ and these rotations envisage negative dependence. The differences in AIC and BIC are small. Nevertheless, the best-performing model in accordance with both criteria turns out to be Frank, Joe 270 and Clayton 90. Hence, these models may be considered if the modeler is interested in making predictions of the number of job changes.

Alternatively, one can use the Vuong test to compare between two competing models. The `VuongClarke()` function returns

```
R> VuongClarke(fit2, fit3)
```

```
Vuong's test: it is not possible to discriminate between the two models.
Clarke's test: Model 1 is preferred over Model 2.
```

By default, the significance level was set to 0.05. The Vuong test does not indicate any preferred model. On the other hand, Clarke test prefers Frank over Joe 270 model.

We will carry out the remainder of the analysis using Frank model. The average predictions for all observations are obtained by

```
R> aver(fit2, univariate = TRUE)
```

Estimated average with 95% confidence interval:

```
0.526 (0.487,0.565)
```

```
R> aver(fit2, univariate = FALSE)
```

Estimated average with 95% confidence interval:

```
0.747 (0.649,0.845)
```

The former statement returns an average prediction for the univariate model ignoring non-random sample selection; the latter returns an average prediction for the sample selection model. To extract the association parameters with their corresponding confidence bounds one needs to submit

```
R> out.sum <- summary(fit2)
```

```
R> out.sum$theta
```

```
[1] -16.69619
```

```
R> out.sum$CIth
```

```
[1] -39.270136 -5.193194
```

```
R> out.sum$tau
```

```
[1] -0.7840277
```

```
R> out.sum$CIkt
```

```
[1] -0.9024080 -0.4686178
```

Table 3 summarizes the corresponding average predictions and association parameters of the univariate, normal, Frank, and the rotated versions of Clayton, Joe and Gumbel models. The values in brackets indicate 95% confidence interval bounds. For all models, the confidence intervals of $\hat{\theta}$ do not reach their bound, which implies that non-random sample selection is likely to be present. Note that Kendall's $\hat{\tau}$ indicates a strong negative correlation between the selection and the outcome. The average predictions do not differ substantially from another. In fact, the confidence intervals of most of the models overlap. Thus, the copula assumption does not seem to have a major impact on the predictions for sample selection models.

The post-estimation QQ-plots can be obtained by submitting the following command line

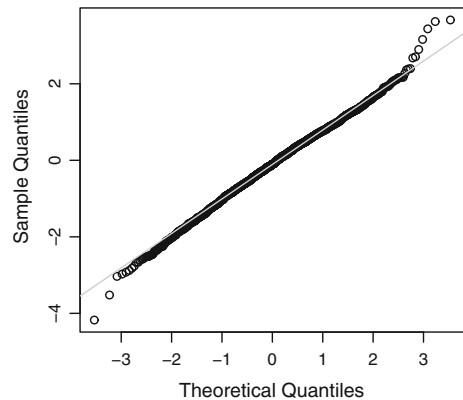
```
R> post.check(fit2)
```

Figure 5 shows post-estimation QQ-plots based on estimates from the Poisson inverse Gaussian sample selection model with Frank copula. Note that the plot is similar to the one in Fig. 4 and hence the model provides a good fit to the data.

Table 3 Average predictions of job changes and estimates of association parameters for Delaporte model and Frank, and 90 and 270 degrees rotated Clayton, Joe and Gumbel

	\bar{y}	$\hat{\theta}$	$\hat{\tau}$
Univariate	0.53 (0.49,0.57)	– –	– –
Normal	0.75 (0.61,0.88)	– 0.76 (– 0.90, – 0.42)	– 0.55 (– 0.71, – 0.28)
Frank	0.75 (0.65,0.85)	– 16.70 (– 39.27, – 5.19)	– 0.78 (– 0.90, – 0.47)
Joe 90	0.74 (0.58,0.91)	– 1.67 (– 3.12, – 1.07)	– 0.27 (– 0.53, – 0.04)
Joe 270	0.75 (0.65,0.84)	– 15.24 (– 47.31, – 8.13)	– 0.88 (– 0.96, – 0.79)
Gumbel 90	0.76 (0.61,0.91)	– 1.67 (– 2.75, – 1.09)	– 0.40 (– 0.64, – 0.09)
Gumbel 270	0.76 (0.64,0.87)	– 3.21 (– 5.14, – 1.95)	– 0.69 (– 0.81, – 0.49)
Clayton 90	0.75 (0.65,0.84)	– 14.36 (– 47.14, – 7.24)	– 0.88 (– 0.96, – 0.78)
Clayton 270	0.75 (0.56,0.94)	– 0.86 (– 2.68, – 0.09)	– 0.30 (– 0.57, – 0.04)

The values in brackets indicate 95% confidence interval bounds

Fig. 5 QQ-plot for Poisson inverse Gaussian and Frank copula

The smooth plots for both equations can be obtained by submitting the following command lines

```
R> par(mfrow = c(1,2))
R> plot(fit2, eq = 1)
R> plot(fit2, eq = 2)
```

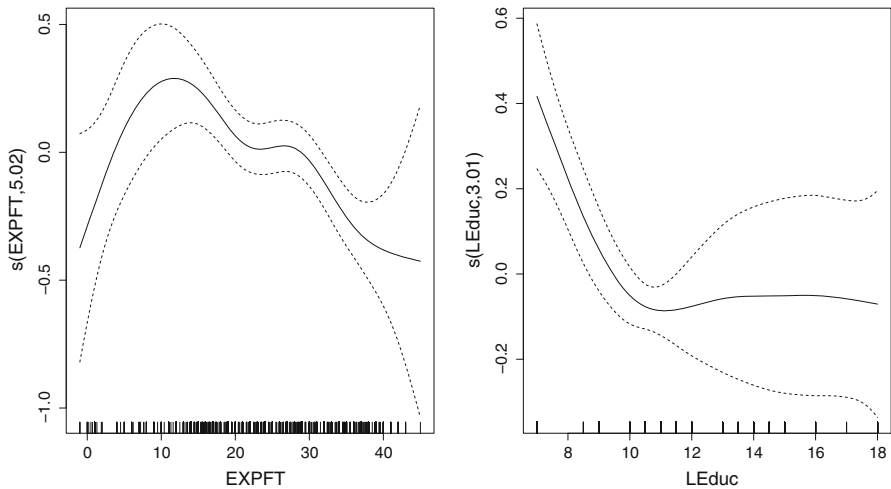


Fig. 6 Selection (*left*) and outcome (*right*) equation smooth for Poisson inverse Gaussian outcome and Frank copula

Figure 6 shows the smooths for the continuous covariates EXPFT (left) and LEduc (right). The selection equation smooth can be interpreted such that the likelihood of being active on the labor market as a dependent worker increases when gaining full time work experience. Then, the likelihood of being active gradually decreases as individuals start to receive pensions or enter into self-employment. The outcome smooth suggests that the longer the education of an individual lasts the less likely he is to change his job. This is due to increasing specialization in his professional area as he becomes more qualified. Therefore, it is harder for him to find an alternative occupation or there are no financial incentives for changing jobs (e.g. medical doctors).

To obtain each σ_i and τ_i (and the corresponding 95 per cent confidence intervals) one needs to submit `fit2$sigma` and `fit2$tau` (`summary(fit2)$CIsig.1` and `summary(fit2)$CIkt.1`). The confidence intervals for white-collars and non-white-collars given each parameter are visualized in Fig. 7. For instance, since the confidence interval for white-collars in τ overlaps with zero, for the group of white-collar workers non-random sample selection is likely to be absent.

4 Summary

In this paper, we introduced the new features of `SemiParSampleSel` R package. The functions included in the package can be used to estimate sample selection models for a wide selection of discrete responses and incorporate flexible covariate effects. The modeling approach allows for a specification of the bivariate distribution using copulae. By doing so, the modeler can check the assumption of bivariate normality.

The function `resp.check()` allows the modeler to perform exploratory analysis on the non-missing response permitting the distributional assumption of the outcome to be checked. `SemiParSampleSel()` can be used to estimate the model under the

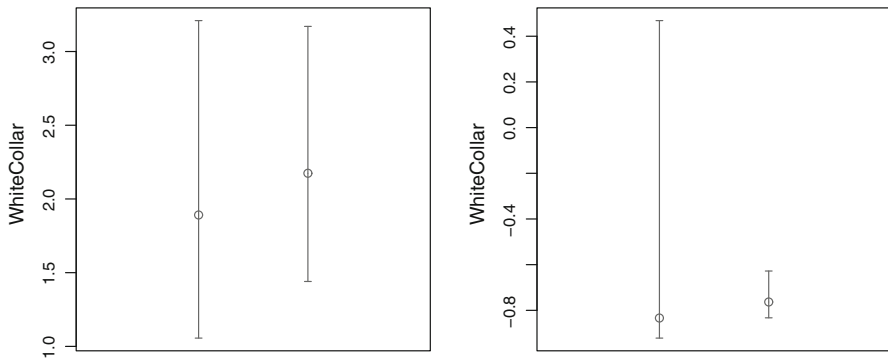


Fig. 7 Point estimate plots with the corresponding confidence intervals for white-collar workers (`WhiteCollar = 1`; *left*) and non-white-collar workers (`WhiteCollar = 0`; *right*). The *left-hand panel* depicts the plot for σ and the *right-hand plot* for τ

desired specification, whereas the `summary()` function returns the output. `aver()` function calculates the average prediction of the univariate and sample selection model for all observations. `AIC()`, `BIC()` and `VuongClarke` can be used for choosing between two competing models; `post.check()` creates post-estimation QQ-plots and `plot()` returns smooth plots for equations, where continuous covariates were specified in terms of splines.

The approach can be extended to trivariate system models. These account for the endogeneity of a treatment variable and for non-random sample selection in the outcome. Also, copulae with two parameters can be introduced - these would lead to a better control of tail-dependence, despite the risk of association parameters losing their interpretation (e.g., [Brechmann and Schepsmeier 2013](#)).

In the context of selection margin, one could employ skew probit links as derived from the standard skew-normal distribution by [Azzalini \(1985\)](#). Introducing a parameter which regulates the distribution's skewness may have very attractive properties from the probability point of view ([Azzalini and Arellano-Valle 2013](#)).

Future research may also involve applying the new `SemiParSampleSel` functions to problems in other fields. For instance, [Greene \(1998\)](#) estimates a sample selection model for predicting the number of non-payments by individuals who own a credit card. In this case, individuals select themselves into the sample of credit card owners based on observed characteristics (such as gender and age) and unobserved ones (e.g. unreported income). In particular, it would be of interest to explore how average predictions change if non-random sample selection is accounted for.

Another application involves patients who are newly-referred to a specialist doctor by a general practitioner and often have to be assessed using ultrasound imaging devices ([Ciurtin et al. 2016](#)). In many circumstances, public health systems cannot afford an assessment of every patient. Non-random sample selection may be present in the assessment scores, since patients are referred based on characteristics that cannot be accounted for e.g. fatigue of the consultant or patient's personality. To rectify this situation, sample selection models can be applied.

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Appendices

1 Margins and copulae implemented in `SemiParSampleSel`

See Tables 4, 5 and 6.

Table 4 Discrete distributions implemented in `SemiParSampleSel` R package. The parameter ranges corresponding to log, logit and identity links are $(0, \infty)$, $[0, 1]$ and $(-\infty, \infty)$. For binomial-type distributions, the binomial denominator is determined by the modeler. Note that the support of the logarithmic distribution does not include zero

Distribution	μ link (μ range)	σ link (σ range)	ν link (ν range)	Binomial denominator	Argument call
Poisson	log	-	-	-	"P"
Delaporte	log	log	logit	-	"D"
Poisson inverse	log	log	-	-	"PIG"
Gaussian					
Sichel	log	log	identity	-	"S"
Beta binomial	logit	log	-	q	"BB"
Binomial	logit	-	-	q	"BI"
Geometric	log	-	-	-	"GEOM"
Logarithmic	logit	-	-	-	"LG"
Negative binomial type I	log	log	-	-	"NBI I"
Negative binomial type II	log	log	-	-	"WARING"
Waring	log	log	-	-	"YULE"
Yule	log	-	-	-	

Table 5 Zero-inflated and zero-altered discrete distributions implemented in **SemiParSampleSel** R package

Distribution	μ link (μ range)	σ link (σ range)	ν link (ν range)	Binomial denominator	Argument call
Zero inflated beta binomial	logit	log	logit	q	"ZIBB"
Zero altered beta binomial	logit	log	logit	q	"ZABB"
Zero inflated binomial	logit	logit	–	q	"ZIBI"
Zero altered binomial	logit	logit	–	q	"ZABI"
Zero adjusted logarithmic	logit	logit	–	–	"ZALG"
Zero inflated negative binomial	log	log	logit	–	"ZINBI"
Zero altered negative binomial	log	log	logit	–	"ZANBI"
Zero altered Poisson	log	logit	–	–	"ZAP"
Zero inflated Poisson	log	logit	–	–	"ZIP"
Zero inflated Poisson type II	log	logit	–	–	"ZIP2"
Zero inflated Poisson inverse Gaussian	log	log	logit	–	"ZIPIG"

The parameter ranges corresponding to log and logit links are $(0, \infty)$ and $[0, 1]$. For binomial-type distributions, the binomial denominator is determined by the modeler. Note that the support of the logarithmic distribution does not include zero

Table 6 Examples of families of bivariate copulae

Name	Copula $C_\theta(u, v)$	Parameter space of θ	Parameter space of Kendall's τ	Kendall's τ in terms of θ	θ^*
FGM	$uv(1 + \theta(1-u)(1-v))$	$[-1, 1]$	$[-2/9, 2/9]$	$\frac{2}{9}\theta$	$\tanh^{-1}(\theta)$
Normal	$\Phi_\theta(\Phi^{-1}(u), \Phi^{-1}(v))$	$[-1, 1]$	$[-1, 1]$	$\frac{2}{\pi} \arcsin(\theta)$	$\tanh^{-1}(\theta)$
AMH	$uv/(1 - \theta(1-u)(1-v))$	$[-1, 1]$	$[-0.1817, \frac{1}{3}]$	$1 - \frac{2}{3\theta^2}(\theta + (1-\theta)^2)$	$\tanh^{-1}(\theta)$
Clayton	$(u^{-\theta} + v^{-\theta} - 1)^{-1/\theta}$	$(0, \infty)$	$(0, 1)$	$\log(1 - \theta)$	$\log(\theta - \epsilon)$
Frank	$-\theta^{-1} \log(1 + (e^{-\theta u} - 1)(e^{-\theta v} - 1))$	$(-\infty, \infty) \setminus \{0\}$	$(-1, 1)$	$\frac{\theta}{\theta+2}$	$\theta - \epsilon$
Gumbel	$\exp(-((-\log u)^\theta + (-\log v)^\theta)^{1/\theta})$	$[1, \infty)$	$[0, 1)$	$1 - \frac{4}{\theta}[1 - D_1(\theta)]$	$\log(\theta - 1)$
Joe	$1 - ((1-u)^\theta + (1-v)^\theta - (1-u)^\theta(1-v)^\theta)^{1/\theta}$	$(1, \infty)$	$(0, 1)$	$1 - \frac{4}{\theta^2} D_2(\theta)$	$\log(\theta - 1 - \epsilon)$

$\Phi_\theta(\cdot, \cdot)$ denotes bivariate standard normal cumulative distribution function with correlation coefficient θ , whereas u and v represent the margins. $D_1(\theta) = \frac{1}{\theta} \int_0^\theta \frac{t}{\exp(t)-1} dt$ is the Debye function and $D_2(\theta) = \int_0^1 t \log(t)(1-t)^{\frac{2(1-\theta)}{\theta}} dt$. ϵ denotes a quantity set to 10^{-8} to ensure that the dependence parameters lie in their respective ranges

2 Data simulation function

For reader's interest, we also provide a code for simulating data suffering from non-random sample selection for Poisson, negative binomial, Delaporte, Poisson inverse Gaussian and Sichel distributions. The potential copulae include the ones mentioned in Table 6. The list of possible outcomes and copulae is not exhaustive and can be extended. In the case below, two covariates are simulated which enter both the selection and outcome equation. n stands for the number of observations to be generated, s . τ is correlation between the outcome and the selection equation defined in terms of Kendall's tau, rhC is the Pearson correlation coefficient for the covariates; `outcome.margin` and `copula` are the outcome margin and copula defined by the user.

```
R> bcds <- function(n, s.tau=0.2, s.sigma=1, s.nu=0.5, rhC=0.2,
+               outcome.margin='PO', copula='FGM') {
+   # Generating covariates
+   SigmaC <- matrix( c(1,rhC,rhC,1), 2 , 2)
+   covariates <- rmvnorm(n,rep(0,2),SigmaC, method='svd')
+   covariates <- pnorm(covariates)
+   x1 <- covariates[,2]; x2 <- round(covariates[,1])
+   # Establishing copula object
+   Cop <- switch(copula,
+               FGM = fgmCopula(dim = 2, param = iTau(fgmCopula(), s.tau)),
+               BN = ellipCopula(family = 'normal', dim = 2,
+                               param = iTau(normalCopula(), s.tau)),
+               AMH = archmCopula(family = 'amh', dim = 2,
+                                param = iTau(amhCopula(), s.tau)),
+               Clayton = archmCopula(family = 'clayton', dim = 2,
+                                     param = iTau(claytonCopula(), s.tau)),
+               Frank = archmCopula(family = 'frank', dim = 2,
+                                   param = iTau(frankCopula(), s.tau)),
+               Gumbel = archmCopula(family = 'gumbel', dim = 2,
+                                    param = iTau(gumbelCopula(), s.tau)),
+               Joe = archmCopula(family = 'joe', dim = 2,
+                                 param = iTau(joeCopula(), s.tau)) )
+   # Creating equations
+   f1 <- function(x) 0.4*(-4 - (5.5*x-2.9) + 3*(4.5*x-2.3)^2 - (4.5*x-2.3)^3)
+   f2 <- function(x) x*sin(8*x)
+   mu_s <- 1.0 + f1(x1) - 2.0*x2
+   mu_o <- exp(1.1 + f2(x1) - 1.9*x2)
+   # Creating margin-dependent object
+   speclist <- switch(outcome.margin,
+                   PO = list(mu = mu_o),
+                   NBI = list(mu = mu_o, sigma = s.sigma),
+                   DEL = list(mu = mu_o, sigma = s.sigma, nu = s.nu),
+                   PIG = list(mu = mu_o, sigma = s.sigma),
+                   SICHEL = list(mu = mu_o, sigma = s.sigma, nu = s.nu) )
+   spec <- mvdc(copula = Cop, c('norm', outcome.margin),
+               list(list(mean = mu_s, sd=1), speclist))
+   # Simulating
+   simGen <- rMvdc(n, spec)
+   y <- ifelse(simGen[,1]>0, simGen[,2], -99)
+   Y
+   # Data frame
+   dataSim <- data.frame(y,x1,x2)
+   dataSim
+ }
```

For instance, to simulate data and estimate a sample selection model with Poisson margin and Frank copula one needs to submit the following code

```
R> library(SemiParSampleSel)
R> set.seed(1)
+ dataSim <- bcds(2000, s.tau=0.5, rhC=0.5,
+               outcome.margin='PO', copula='Frank')
R> # Generating selection variable
R> dataSim$y.probit<-ifelse(dataSim$y>=0, 1, 0)
R> out <- SemiParSampleSel(list(y.probit ~ s(x1) + x2, y ~ s(x1) + x2),
+                            data=dataSim, BivD = 'F', margins = c('N', 'P'))
```

3 Preliminary analysis plots

See Fig. 8

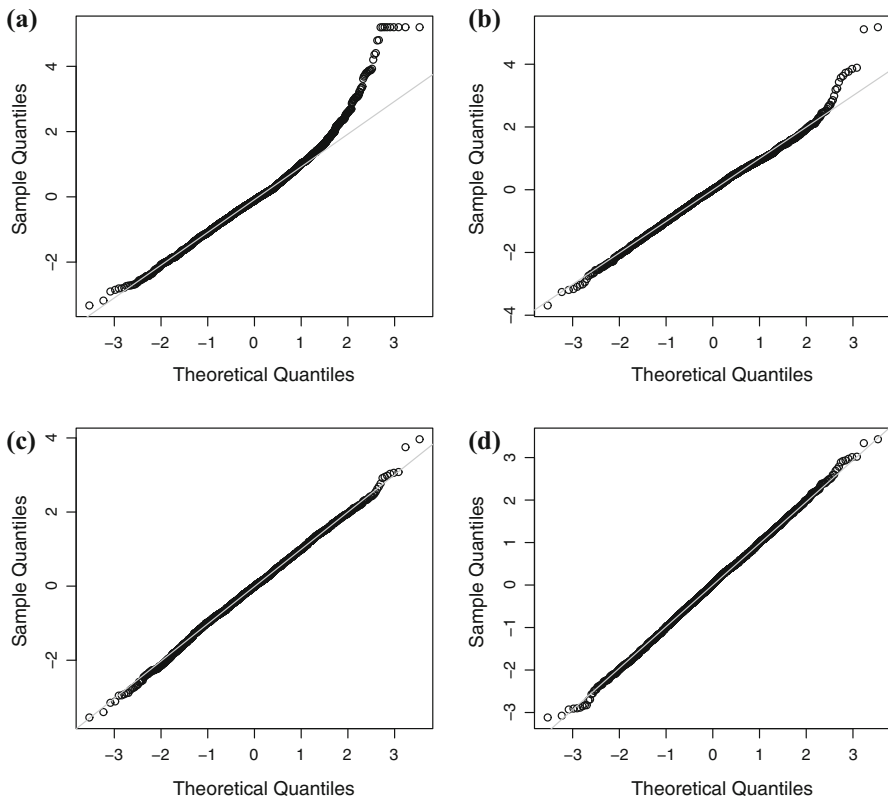


Fig. 8 Poisson (a), negative binomial (b), Delaporte (c) and Sichel (d) QQ-plots

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