



Robust mixture regression model fitting by Laplace distribution



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ABSTRACT

A robust estimation procedure for mixture linear regression models is proposed by assuming that the error terms follow a Laplace distribution. Using the fact that the Laplace distribution can be written as a scale mixture of a normal and a latent distribution, this procedure is implemented by an EM algorithm which incorporates two types of missing information from the mixture class membership and the latent variable. Finite sample performance of the proposed algorithm is evaluated by simulations. The proposed method is compared with other procedures, and a sensitivity study is also conducted based on a real data set.

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1. Introduction

Least absolute deviation (LAD) regression has been widely used in practice if robust estimation is desired. The research on its computation and theoretical properties is abundant in the literature. A detailed survey on this topic can be found in Dielman (1984, 2005). It is known that the outliers impact more heavily on mixture linear regression models than on the usual linear regression models, since the outliers not only affect the estimation of the regression parameters, but also possibly totally blur the mixture structure. In this paper, LAD will be applied to a class of mixture linear regression models. Simulation studies show that the proposed estimators of the regression coefficients are robust.

To be specific, let X be a p -dimensional vector of explanatory variables and Y be a scalar response variable. The relationship between Y and X is often investigated through a linear regression model. In the mixture linear regression setup, we assume that with probability π_i , $i = 1, 2, \dots, g$, (X', Y) comes from one of the following $g \geq 2$ linear regression models

$$Y = X'\beta_i + \sigma_i\varepsilon_i, \quad i = 1, 2, \dots, g, \quad (1)$$

where $\sum_{i=1}^g \pi_i = 1$, the β_i 's are unknown p -dimensional vectors of regression coefficients, and the σ_i 's are unknown positive scalars. The random errors ε_i 's are assumed to be independent of the X_i 's. It is commonly assumed that the density functions of ε_i 's are members in a location-scale family with means 0 and variances 1. In the following discussion, the design variable X is assumed to be random, but the proposed estimation procedure also works for the fixed design.

If $g = 1$, the LAD estimator of β is the minimizer of the target function $Q(\beta) = \sum_{j=1}^n |Y_j - X_j'\beta|$, where $(X_j', Y_j)_{j=1}^n$ is a sample from model (1). Many algorithms have been developed in the literature to tackle the minimization problem $\hat{\beta} = \operatorname{argmin}_{\beta} Q(\beta)$, such as linear programming, least angle regression, modified maximum likelihood method by Li and

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Arce (2004), among others. An often adopted but ad-hoc scheme for finding the MLE of β is to obtain the root of the derivative of $Q(\beta)$. Here σ^2 is treated as a nuisance parameter. By doing this, we obtain

$$\frac{\partial Q(\beta)}{\partial \beta} = - \sum_{j=1}^n X_j \operatorname{sgn}(Y_j - X'_j \beta) = 0, \tag{2}$$

where $\operatorname{sgn}(\cdot)$ is the sign function which takes $-1, 0, 1$ if the argument is negative, 0, and positive, respectively. Let $w_j = 1/|Y_j - X'_j \beta|$, and rewrite the Eq. (2) as $\sum_{j=1}^n w_j X_j (Y_j - X'_j \beta) = 0$. Thus by supplying an initial value β_0 for β , the updated value β can be found by the weighted least square solution

$$\beta_1 = \left(\sum_{j=1}^n w_j X_j X'_j \right)^{-1} \sum_{j=1}^n w_j X_j Y_j, \tag{3}$$

where $w_j = 1/|Y_j - X'_j \beta_0|$. By iterating the procedure, one can eventually find an approximate solution to $\operatorname{argmin}_{\beta} Q(\beta)$.

A very interesting connection between the iterated weighted least square procedure stated above and an EM algorithm in conjunction with the Laplace distribution is found in Phillips (2002). For the sake of completeness, we briefly describe the procedure proposed in Phillips (2002).

Andrews and Mallows (1974) showed that a Laplace distribution can be expressed as a mixture of a normal distribution and another distribution related to the exponential distribution. To be specific, suppose Z and V are two random variables, V has a distribution with density function $v^{-3} \exp(-(2v^2)^{-1})$, $v > 0$, and given $V = v$, the conditional distribution of Z is normal with mean 0 and variance $\sigma^2/(2v^2)$. Then Z marginally has a Laplace distribution with density function $h_{\epsilon}(z) = \exp(-\sqrt{2}|z|/\sigma)/(\sqrt{2}\sigma)$. Based on this, Phillips (2002) developed an EM algorithm to search for the minimizer of $Q(\beta)$.

If V could be observed, then the complete log-likelihood function of $\theta = (\beta, \sigma^2)$, based on the sample $\mathbf{P} = (X_j, Y_j, V_j)_{j=1}^n$, is

$$L(\theta; \mathbf{P}) = -\frac{n}{2} \log(\pi \sigma^2) - \frac{1}{\sigma^2} \sum_{j=1}^n V_j^2 (Y_j - X'_j \beta)^2 - \sum_{j=1}^n \log V_j^2 - \frac{1}{2} \sum_{j=1}^n \frac{1}{V_j^2}.$$

Following the two steps in the EM algorithm procedure, and assuming that $\theta^{(k)} = (\beta^{(k)}, \sigma^{2(k)})$ is the value for the k th iteration, then in the $(k + 1)$ th iteration, we have to first calculate the conditional expectation of the complete log likelihood function $L(\theta; \mathbf{P})$, given the observed data set $(Y_j, X_j)_{j=1}^n$ and $\theta = \theta^{(k)}$, which has the following form

$$E[L(\theta; \mathbf{P})|\mathbf{S}] = -\frac{n}{2} \log(\pi \sigma^2) - \frac{\sum_{j=1}^n E[V_j^2 | \theta^{(k)}, (X_j, Y_j)_{j=1}^n] (Y_j - X'_j \beta)^2}{\sigma^2} - \sum_{j=1}^n E[\log V_j^2 | \theta^{(k)}, (X_j, Y_j)_{j=1}^n] - \frac{1}{2} \sum_{j=1}^n E \left[\frac{1}{V_j^2} \middle| \theta^{(k)}, (X_j, Y_j)_{j=1}^n \right].$$

In the second step, the conditional expectation is maximized over θ . Denote $w_j = E[V_j^2 | \theta^{(k)}, (X_j, Y_j)_{j=1}^n]$, and notice that the third and fourth terms on the right hand side do not involve the unknown regression parameters. Therefore, to maximize the above conditional expectation is equivalent to maximize the following terms with respect to θ ,

$$-\frac{n}{2} \log \sigma^2 - \frac{\sum_{j=1}^n w_j (Y_j - X'_j \beta)^2}{\sigma^2}.$$

Interestingly, Phillips (2002) showed $w_j = E[V_j^2 | \theta^{(k)}, (X_j, Y_j)_{j=1}^n] = \sigma^{(k)}/(\sqrt{2}|Y_j - X'_j \beta^{(k)}|)$. This implies that the solution $\beta^{(k+1)}$ is the same as the one based on (3) and the iteratively reweighted least squares procedure is an application of the EM algorithm. It is also easy to see that $\sigma^{2(k+1)}$ can be estimated by $2 \sum_{j=1}^n w_j (Y_j - X'_j \beta^{(k+1)})^2/n$.

The robustness property of the LAD procedure, and the natural connection between LAD estimation and maximum likelihood estimation for the regression coefficients given Laplace distributed random error when $g = 1$, motivate us to consider the possible extension of the algorithm to the mixture model setup. When $g \geq 2$, we assume that for each i , $i = 1, 2, \dots, g$, ϵ_i follows a Laplace distribution with location 0 and scale parameter $1/\sqrt{2}$, which results in the variance of ϵ_i being 1. Then it is easily seen that for a sample $\mathbf{S} = \{(X'_j, Y_j), j = 1, 2, \dots, n\}$ from the model (1), the log-likelihood function of $\theta = (\beta_1, \sigma_1^2, \pi_1, \beta_2, \sigma_2^2, \pi_2, \dots, \beta_g, \sigma_g^2, \pi_g)$ can be written as

$$L(\theta; \mathbf{S}) = \sum_{j=1}^n \log \left[\sum_{i=1}^g \frac{\pi_i}{\sqrt{2}\sigma_i} \exp \left(-\frac{\sqrt{2}|Y_j - X'_j \beta_i|}{\sigma_i} \right) \right]. \tag{4}$$

Usually no explicit MLE is available. In the following, another missing component will be incorporated into the log-likelihood function (4), so that the solution can be obtained by a standard use of the EM algorithm.

Wei (2012) proposed a robust estimation procedure for the mixture linear regression models based on the t distribution by extending McLachlan and Peel (2000)’s work. The research conducted in this paper deals with the same questions as in Wei (2012), but with the LAD technique, or the Laplace distribution, instead of the less commonly used t -distribution, used for achieving robustness. The natural connection between the LAD procedure and the MLE procedure based on Laplace error makes the proposed procedure more appealing.

The paper is organized as follows. The EM algorithm is developed in Section 2, together with some discussion on how to control the outliers in the x -direction. Section 3 conducts some numerical simulations to evaluate the finite sample performance of the proposed method, and compares it with some other existing methods. Finally, a sensitivity study for various estimation procedures is conducted in Section 3 based on a real data set.

2. EM algorithm for robust mixture regression

In model (1), assume that ε_i 's follow a Laplace distribution with mean 0 and scale parameter $1/\sqrt{2}$. For $i = 1, 2, \dots, g$, $j = 1, 2, \dots, n$, denote G_{ij} as latent Bernoulli variables such that

$$G_{ij} = \begin{cases} 1, & \text{if } j\text{th observation } (X_j, Y_j) \text{ is from } i\text{th component;} \\ 0, & \text{otherwise.} \end{cases}$$

If the complete data set $\mathbf{T} = \{(X_j, Y_j, G_{ij})\}_{i=1,2,\dots,g; j=1,2,\dots,n}$ is observable, the complete log likelihood function of $\theta = (\beta_1, \sigma_1^2, \pi_1, \beta_2, \sigma_2^2, \pi_2, \dots, \beta_g, \sigma_g^2, \pi_g)$ can be written as

$$L(\theta; \mathbf{T}) = \sum_{j=1}^n \sum_{i=1}^g G_{ij} \log \left[\frac{\pi_i}{\sqrt{2}\sigma_i} \exp \left(-\frac{\sqrt{2}|Y_j - X'_j\beta_i|}{\sigma_i} \right) \right]. \tag{5}$$

From Andrews and Mallows (1974), we know that a Laplace distributed random variable is a scale mixture of a normal random variable and another variable related to the exponential distribution. Also see Section 1 for detail. Denote V_j , coupled with (X_j, Y_j) , as the latent scale variable, $j = 1, 2, \dots, n$. Then the complete log-likelihood function of θ , based on $\mathbf{D} = \{X_j, Y_j, V_j, G_{ij}\}_{i=1,2,\dots,g; j=1,2,\dots,n}$, has the form

$$\begin{aligned} L(\theta; \mathbf{D}) &= \sum_{j=1}^n \sum_{i=1}^g G_{ij} \log \left[\pi_i \frac{V_j}{\sqrt{\pi}\sigma_i} \exp \left(-\frac{V_j^2(Y_j - X'_j\beta_i)^2}{\sigma_i^2} \right) \frac{1}{V_j^3} \exp \left(-\frac{1}{2V_j^2} \right) \right] \\ &= \sum_{j=1}^n \sum_{i=1}^g G_{ij} \log \pi_i - \frac{1}{2} \sum_{j=1}^n \sum_{i=1}^g G_{ij} \log(\pi\sigma_i^2) - \sum_{j=1}^n \sum_{i=1}^g \frac{G_{ij}V_j^2(Y_j - X'_j\beta_i)^2}{\sigma_i^2} \\ &\quad - \sum_{j=1}^n \sum_{i=1}^g G_{ij} \log V_j^2 - \frac{1}{2} \sum_{j=1}^n \sum_{i=1}^g \frac{G_{ij}}{V_j^2}. \end{aligned} \tag{6}$$

Based on the EM algorithm principle, in the E-step, we have to calculate the condition expectation $E[L(\theta; \mathbf{D})|\mathbf{S}, \theta^{(0)}]$, where $\mathbf{S} = \{(X_j, Y_j)\}_{j=1}^n$, and $\theta^{(0)} = (\beta_1^{(0)}, \sigma_1^{2(0)}, \pi_1^{(0)}, \dots, \beta_g^{(0)}, \sigma_g^{2(0)}, \pi_g^{(0)})$ is a proper initial value for θ . Since the last two terms in (6) do not involve the unknown regression parameters, we can simply drop them from the analysis. Thus, to find $E[L(\theta; \mathbf{D})|\mathbf{S}, \theta^{(0)}]$, we only have to calculate the following two terms

$$\tau_{ij} = E[G_{ij}|\mathbf{S}, \theta^{(0)}], \quad \delta_{ij} = E[V_j^2|\mathbf{S}, \theta^{(0)}, G_{ij} = 1].$$

One can show that

$$\tau_{ij} = \frac{\pi_i^{(0)} \sigma_i^{-1(0)} \exp(-\sqrt{2}|Y_j - X'_j\beta_i^{(0)}|/\sigma_i^{(0)})}{\sum_{m=1}^g \pi_m^{(0)} \sigma_m^{-1(0)} \exp(-\sqrt{2}|Y_j - X'_j\beta_m^{(0)}|/\sigma_m^{(0)})}, \quad \delta_{ij} = \frac{\sigma_i^{(0)}}{\sqrt{2}|Y_j - X'_j\beta_i^{(0)}|}. \tag{7}$$

The calculation for δ_{ij} follows the same argument as in Phillips (2002). In the M-step, the following expression will be maximized with respect to π_i 's, β_i 's and σ_i^2 's,

$$\sum_{j=1}^n \sum_{i=1}^g \tau_{ij} \log \pi_i - \frac{1}{2} \sum_{j=1}^n \sum_{i=1}^g \tau_{ij} \log \sigma_i^2 - \sum_{j=1}^n \sum_{i=1}^g \frac{\tau_{ij}\delta_{ij}(Y_j - X'_j\beta_i)^2}{\sigma_i^2}, \tag{8}$$

and the maximizer will be used for the next iteration.

In summary, we propose the following EM algorithm to maximize (4).

EM algorithm

- (1) Choose an initial value for $\theta = (\beta_1, \sigma_1^2, \pi_1, \dots, \beta_g, \sigma_g^2, \pi_g)$.
- (2) E-Step: at the $(k + 1)$ -th iteration, calculate $\tau_{ij}^{(k+1)}$ and $\delta_{ij}^{(k+1)}$ from Eq. (7) with (0) replaced by (k) .
- (3) M-Step: at the $(k + 1)$ -th iteration, use the following formulas to calculate the maximizer of (8):

$$\pi_i^{(k+1)} = \frac{1}{n} \sum_{j=1}^n \tau_{ij}^{(k)}, \quad \beta_i^{(k+1)} = \left(\sum_{j=1}^n \tau_{ij}^{(k+1)} \delta_{ij}^{(k+1)} X_j X_j' \right)^{-1} \left(\sum_{j=1}^n \tau_{ij}^{(k+1)} \delta_{ij}^{(k+1)} X_j Y_j \right),$$

and

$$\sigma_i^{2(k+1)} = \frac{2 \sum_{j=1}^n \tau_{ij}^{(k+1)} \delta_{ij}^{(k+1)} (Y_j - X_j' \beta_i^{(k+1)})^2}{\sum_{j=1}^n \tau_{ij}^{(k+1)}}.$$

- (4) Repeat steps (2), (3) until convergence is obtained.

If we further assume that all σ_i^2 's are equal, then in the above EM algorithm, a common initial value for σ_i^2 should be used, and σ^2 can be updated in the M-step by

$$\sigma^{2(k+1)} = \frac{2 \sum_{j=1}^n \sum_{i=1}^g \tau_{ij}^{(k+1)} \delta_{ij}^{(k+1)} (Y_j - X_j' \beta_i^{(k+1)})^2}{n}.$$

The robustness of the EM procedure above follows from the adoption of LAD regression. It is also obvious from the formulas of the updated β_i 's in each iteration. Note that the factor δ_{ij}^{k+1} is reciprocally related to the term $|Y_j - X_j' \beta_i^{(k)}|$, meaning that larger residuals give smaller values of δ_{ij}^{k+1} , hence downweight the corresponding observations when calculating the estimates.

The EM algorithm proposed above for calculating the estimate of β is an iterated reweighted least square (IRLS) procedure, as the one proposed in Schlossmacher (1973) for the one population case and the weights are given by $\tau_{ij}^{(k+1)} \delta_{ij}^{(k+1)}$ for the $k + 1$ -th iteration. Extra attention should be paid when programming the proposed EM algorithm. In the case of $g = 1$, Schlossmacher (1973) warned that if a perfect LAD fit occurs, i.e., $Y_j - X_j' \hat{\beta}_i = 0$ for some i, j and $\hat{\beta}_i$, then the algorithm will eventually give $Y_j - X_j' \beta_i^k \approx 0$ when iteration proceeds. As a result, δ_{ij}^{k+1} , which is reciprocally related to $|Y_j - X_j' \beta_i^k|$, will be very large, and numerical instability would follow. Although Phillips (2002) noticed that this problem rarely arises in the case of $g = 1$, this does occur often in our case, which is not surprising, simply because more than one regression model provides more chance for a perfect LAD fitting. But simply adopting Schlossmacher (1973)'s weight scheme by setting $\delta_{ij}^{k+1} = 0$ whenever $|Y_j - X_j' \beta_i^k| < e$ for a pre-assigned $e > 0$ is not quite reasonable. It makes sense to allocate large weights for small residuals and small weights for large residuals and a cogent argument on this issue is provided in Phillips (2002). In our simulation study, we simply adopt a hard threshold rule to control the extremely small LAD residuals in each iteration step. Under this rule, $\delta_{ij}^{(k+1)}$ will be assigned a value of 10^6 for any perfect LAD fit. We also tried other threshold values, such as $10^8, 10^{10}$ in the simulation, all these choices generate almost identical results. For the sake of brevity, we only report the simulation results by using 10^6 as the threshold value. A smooth alternative to the above hard threshold, as one referee recommended, is to replace $|Y_j - X_j' \beta_i^k|$ in the denominator of δ_{ij}^{k+1} by $|Y_j - X_j' \beta_i^k| + c$, where $c = 10^{-6}$ or some other small constants.

It is well known that in the IRLS procedure, numerical instability could occur if the weights are very small. A common way to deal with this issue is to impose a hard threshold on τ_{ij}^{k+1} obtained in the $k + 1$ -th iteration. Namely, for a pre-specified value e say, if $\tau_{ij}^{k+1} > e$, then τ_{ij}^{k+1} itself will be used for the next iteration; otherwise, e will be used as the weight for the next iteration as is used in Wei (2012). In our simulation study, $e = 10^{-6}$ is adopted.

Similar to the traditional M-estimate for linear regression and Wei (2012)'s mixture regression by a t -distribution, the above EM algorithm based on the Laplace distribution is robust against outliers along the y -direction, but not in the x -direction. This is also confirmed by the sensitivity study conducted in Section 3. As a consequence, if there are any high leverage points in the data sets, then the proposed EM algorithm might fail our aims, and certain modifications would be necessary. An obvious modification is first to identify these high leverage points, then just discard them. A commonly used method is to calculate the leverage value for each observation using the formula $h_{jj} = n^{-1} + (n - 1)^{-1} MD_j$, where $MD_j = (X_j - \bar{X})' S^{-1} (X_j - \bar{X})$, and \bar{X}, S are the sample mean and sample covariance matrix of X_j 's, respectively. The j -th observation will be identified as a high leverage point if $h_{jj} > 2p/n$, where p is the dimension of X . To avoid the masking effect caused by using \bar{X} and S in detecting the high leverage points, some robust estimation of the population mean and covariance matrix of X can be used instead of the sample mean and sample covariance. Wei (2012) adopted the minimum

covariance determinant (MCD) estimators for the population mean and covariance matrix, which is implemented by the Fast MCD algorithm developed in Rousseeuw and Van Driessen (1999). Certainly, other robust estimates of the population mean and covariance matrix could be also used for this purpose, for example, the Stahel–Donoho (SD) estimator from Stahel (1981) and Donoho (1982). The j -th observation will be considered as a high leverage point if the resulting MD_j exceeds the threshold $\chi_{p,0.975}^2$, as proposed by Pison et al. (2002). In the simulation studies, we apply the proposed EM algorithm based on the Laplace distribution after removing the observations with $MD_j > \chi_{p,0.975}^2$ using both the MCD estimator and the SD estimator to calculate MD_j .

3. Numerical studies

To assess the finite sample performance of the proposed robust estimation procedure, we conducted an extensive simulation study. It is well known that label switching is always an issue when evaluating different estimation methods in mixture models, and there are no widely accepted labeling standards. In our simulation, as in Wei (2012), we simply choose the labels by minimizing the distance to the true parameter values. The effects of labeling schemes on the comparison of different estimation procedures deserves further study.

3.1. Simulation studies

In the simulation study, we chose equal variance for all components. The reason for doing this is that, if the variances are not same, the log-likelihood function (5) is unbounded and goes to infinity if one observation lies exactly on one component line and the corresponding variance goes to 0, which makes the simulation very unstable.

To compare our method with some existing estimation procedures, we generated sample data $(X_{j1}, X_{j2}, Y_j)_{j=1}^n$ from the following two-component mixture regression models, following Wei (2012):

$$Y = \begin{cases} 0 + X_1 + X_2 + \varepsilon_1, & \text{if } Z = 1, \\ 0 - X_1 - X_2 + \varepsilon_2, & \text{if } Z = 2, \end{cases}$$

where Z is the component indicator. That is, the data were generated from a two-component mixture linear regression model with $\beta_1 = (\beta_{10}, \beta_{11}, \beta_{12})' = (0, 1, 1)'$, and $\beta_2 = (\beta_{20}, \beta_{21}, \beta_{22})' = (0, -1, -1)'$. The predictors $X_1 \sim N(0, 1)$ and $X_2 \sim N(0, 1)$ are independent. The random error ε_1 and ε_2 are independent and have the same distribution as ε . To see the effects of different distributions of ε and high leverage outliers in x -direction on various estimation methods, we considered the following six cases: (1), $\varepsilon \sim N(0, 1)$; (2), $\varepsilon \sim$ Laplace distribution with mean 0 and variance 1; (3), $\varepsilon \sim t_1$, t -distribution with 1 degree of freedom or the Cauchy distribution; (4), $\varepsilon \sim t_3$, t -distribution with 3 degrees of freedom. (5), $\varepsilon \sim 0.95N(0, 1) + 0.05N(0, 25)$, a mixture of two normal distributions; (6), $\varepsilon \sim N(0, 1)$ with 5% high leverage outliers being $X_1 = X_2 = 20$ and $Y = 100$.

Case 1 is often used to evaluate the efficiency of different estimation methods compared to the traditional MLE when the error is exactly normally distributed and there are no outliers. For Case 2, the estimation methods proposed in this paper will provide the MLE of unknown parameters, which, as in the first case, would serve as a baseline to evaluate the performance of other estimation procedures. Both Case 3 and 4 are heavy tailed distributions and are often used in the literature to mimic the outlier situations. Case 5 would produce 5% data likely to be low leverage outliers, and in Case 6, 5% of the observations are replicated serving as the high leverage outliers, which will be used to check the robustness of estimation procedures against the high leverage outliers.

Nine estimation methods were compared in the simulation study: (1), the maximum likelihood method based on the normality assumption (MLE); (2), the trimmed likelihood estimator (TLE) proposed by Neykov et al. (2007); (3), the robust modified EM algorithm based on bisquare (Bisquare) proposed by Bai et al. (2012); (4), robust mixture regression based on the t -distribution (Mixreg t) proposed by Wei (2012); (5), the trimmed mixture regression based on t -distribution (Mixreg t -MCD), with MCD trimming method; (6), the trimmed mixture regression method based on t -distribution (Mixreg t -SD), with SD trimming method; (7), the proposed robust EM mixture regression method based on the Laplace-distribution (MixregL); (8), the trimmed mixture regression method based on the Laplace-distribution (MixregL-MCD), with MCD trimming method, and (9), the trimmed mixture regression method based on the Laplace-distribution (MixregL-SD), with the SD trimming method. In all simulation studies, the iteration is terminated when the change in the likelihood function is less than 10^{-6} .

The simulation results are present in Tables 1 and 2. From the simulation studies, we can see that if the true distribution of ε is normal, the MSEs of MLE procedure are slightly larger than our proposed method for the first regression component when the sample size is 100, but the superiority of MLE over all other methods becomes clear when the sample size gets larger. For other cases when the distribution of ε has a heavier tail, contaminated by some outliers, or there are high leverage outliers in the data set, MLE fails to provide reasonable estimates.

The performance of TLE and Bisquare is satisfying when ε has a lighter tail. See the simulation results for all cases except Case III, where ε has a t -distribution with 1 degree of freedom. The overall performance of the Mixreg t proposed by Wei (2012) is also satisfying when the sample size gets larger except for Case VI when high leverage points are present in the data set, but this disadvantage is remedied by the modified procedure Mixreg t -MCD.

Table 1
MSE (bias) of point estimates for $n = 100$.

	MLE	TLE	Bisquare	Mixregt	Mixregt-MCD	MixregL	MixregL-MCD
Case I: $\varepsilon \sim N(0, 1)$							
β_{10}	0.130(0.011)	0.139(0.033)	0.143(0.011)	0.124(0.021)	0.163(0.029)	0.093(0.079)	0.090(0.069)
β_{11}	0.160(−0.025)	0.212(−0.195)	0.157(−0.022)	0.130(−0.032)	0.175(−0.115)	0.094(−0.015)	0.113(−0.103)
β_{12}	0.135(−0.034)	0.248(−0.195)	0.171(−0.048)	0.123(−0.004)	0.247(−0.031)	0.088(0.008)	0.165(−0.039)
β_{20}	0.018(−0.003)	0.038(−0.004)	0.021(−0.001)	0.022(−0.012)	0.022(0.008)	0.028(−0.026)	0.027(−0.001)
β_{21}	0.021(−0.016)	0.030(0.011)	0.023(−0.017)	0.021(−0.006)	0.029(−0.011)	0.027(−0.001)	0.035(−0.021)
β_{22}	0.018(0.009)	0.024(0.034)	0.019(0.014)	0.021(−0.010)	0.030(−0.020)	0.026(−0.010)	0.042(−0.017)
π_1	0.005(0.003)	0.007(0.025)	0.005(0.005)	0.005(0.013)	0.007(0.016)	0.005(0.017)	0.007(0.022)
Case II: $\varepsilon \sim \text{Laplace}(1)$							
β_{10}	0.177(−0.006)	0.075(−0.007)	0.137(−0.016)	0.085(0.012)	0.123(−0.001)	0.058(0.022)	0.060(0.020)
β_{11}	0.145(−0.040)	0.097(−0.107)	0.142(−0.054)	0.084(−0.029)	0.150(−0.033)	0.050(−0.024)	0.080(−0.033)
β_{12}	0.152(0.009)	0.084(−0.077)	0.126(0.000)	0.080(−0.021)	0.150(−0.026)	0.055(−0.006)	0.063(−0.020)
β_{20}	0.016(−0.002)	0.013(0.004)	0.013(0.002)	0.011(−0.007)	0.016(−0.019)	0.010(−0.010)	0.015(−0.026)
β_{21}	0.021(−0.017)	0.013(0.007)	0.014(−0.019)	0.012(−0.008)	0.018(−0.030)	0.011(−0.004)	0.019(−0.020)
β_{22}	0.016(−0.006)	0.013(0.019)	0.013(−0.002)	0.012(−0.002)	0.020(0.009)	0.012(0.003)	0.026(0.018)
π_1	0.004(−0.004)	0.004(0.019)	0.004(0.016)	0.004(0.015)	0.005(0.012)	0.003(0.013)	0.005(0.009)
Case III: $\varepsilon \sim t_1$							
β_{10}	242.992(−0.120)	3.200(−0.150)	1.683(−0.116)	1.708(−0.026)	0.945(−0.075)	0.163(0.061)	0.122(0.034)
β_{11}	174.666(−1.568)	1.886(−0.170)	1.571(−0.347)	1.990(−0.252)	1.621(−0.535)	0.521(−0.377)	0.561(−0.430)
β_{12}	148.108(−1.770)	1.797(−0.033)	1.642(−0.306)	2.410(−0.447)	1.538(−0.360)	0.548(−0.412)	0.418(−0.405)
β_{20}	244.822(0.172)	1.526(0.065)	0.910(0.024)	0.113(−0.020)	3.237(−0.173)	0.032(−0.024)	0.025(−0.038)
β_{21}	175.583(−1.080)	0.774(−0.129)	0.489(−0.088)	0.079(−0.041)	0.949(−0.102)	0.032(0.052)	0.047(0.081)
β_{22}	142.861(−0.454)	0.773(−0.065)	0.580(−0.116)	0.112(−0.049)	0.968(−0.028)	0.037(0.052)	0.048(0.054)
π_1	0.084(0.213)	0.039(0.060)	0.047(0.105)	0.023(0.093)	0.028(0.108)	0.022(0.070)	0.023(0.083)
Case IV: $\varepsilon \sim t_3$							
β_{10}	1.568(−0.129)	0.238(0.007)	0.460(0.006)	0.529(0.031)	0.475(0.126)	0.131(0.065)	0.130(0.108)
β_{11}	0.997(−0.234)	0.264(−0.135)	0.341(−0.041)	0.361(0.010)	0.772(−0.109)	0.176(−0.021)	0.183(−0.041)
β_{12}	1.240(−0.024)	0.239(−0.096)	0.375(−0.058)	0.394(−0.010)	0.804(−0.040)	0.132(0.013)	0.186(−0.046)
β_{20}	0.723(−0.029)	0.038(−0.008)	0.063(0.013)	0.034(0.002)	0.077(−0.018)	0.032(−0.005)	0.030(−0.009)
β_{21}	0.188(0.028)	0.034(0.010)	0.085(−0.034)	0.037(−0.005)	0.062(−0.014)	0.042(0.004)	0.052(−0.018)
β_{22}	0.115(0.031)	0.026(0.010)	0.041(−0.013)	0.029(−0.018)	0.166(−0.027)	0.035(−0.015)	0.048(0.003)
π_1	0.028(0.025)	0.007(0.037)	0.009(0.030)	0.006(0.011)	0.014(0.035)	0.007(0.012)	0.007(0.021)
Case V: $\varepsilon \sim 0.95N(0, 1) + 0.05N(0, 25)$							
β_{10}	2.243(−0.020)	0.124(0.046)	0.202(0.042)	0.152(0.015)	0.350(0.037)	0.097(0.034)	0.098(0.042)
β_{11}	1.366(0.054)	0.282(−0.209)	0.225(−0.037)	0.153(−0.029)	0.528(−0.106)	0.100(−0.008)	0.160(−0.056)
β_{12}	2.117(−0.113)	0.221(−0.190)	0.217(−0.056)	0.163(−0.050)	0.705(0.094)	0.099(−0.030)	0.175(0.023)
β_{20}	1.767(0.159)	0.030(0.013)	0.021(0.011)	0.026(0.020)	0.028(−0.004)	0.029(0.008)	0.035(−0.003)
β_{21}	1.277(−0.122)	0.034(0.001)	0.028(−0.023)	0.022(−0.009)	0.035(0.010)	0.026(−0.005)	0.040(0.008)
β_{22}	0.284(0.006)	0.027(0.011)	0.029(−0.009)	0.120(−0.036)	0.038(−0.017)	0.027(−0.006)	0.044(−0.020)
π_1	0.040(0.015)	0.010(0.034)	0.008(0.020)	0.007(0.015)	0.009(0.012)	0.005(0.006)	0.009(0.013)
Case VI: $\varepsilon \sim N(0, 1)$ with 5% high leverage outliers							
β_{10}	18.364(−2.878)	0.173(0.002)	0.152(0.015)	2.456(0.169)	0.175(−0.032)	0.036(0.080)	0.111(0.092)
β_{11}	5.876(1.422)	0.248(−0.209)	0.200(−0.068)	3.444(1.473)	0.219(−0.055)	0.056(−0.037)	0.133(−0.012)
β_{12}	6.520(1.641)	0.219(−0.168)	0.227(−0.091)	3.589(1.517)	0.262(0.006)	0.042(−0.014)	0.153(−0.046)
β_{20}	11.938(2.451)	0.036(−0.002)	0.023(−0.011)	0.023(0.002)	0.027(0.019)	0.015(−0.058)	0.032(0.011)
β_{21}	12.578(3.316)	0.028(0.000)	0.025(−0.014)	0.053(0.139)	0.027(0.010)	0.013(0.033)	0.042(0.000)
β_{22}	12.561(3.315)	0.022(0.025)	0.020(0.019)	0.053(0.136)	0.023(−0.017)	0.012(0.021)	0.046(0.004)
π_1	0.113(0.165)	0.007(0.017)	0.007(0.003)	0.007(−0.074)	0.006(0.005)	0.005(0.030)	0.006(0.011)

The simulation results clearly show that the proposed method in the paper outperforms or is at least comparable to any other methods. It is rather unexpected that our proposed method performs better than the Mixregt and Mixregt-MCD procedures even when ε has a t -distribution. The larger MSEs in the latter two procedures might have resulted from the extra step involved in the algorithm, the selection of v , the degrees of freedom of the t -distribution.

The MCD estimator is used in Mixregt-MCD and MixregL-MCD to remove high leverage outliers. In the simulation study, the SD estimator is also used to remove high leverage outliers. The simulation results are similar to those from Mixregt-MCD and MixregL-MCD, hence are omitted here for the sake of brevity.

A common criticism about the EM algorithm is its slow convergence. This can also be seen from our simulation study. The average numbers of iterations to achieve convergence using the MixregL procedure are 96, 97, 78, 98, 102, 10 for Case I to Case VI, respectively, when the sample size is 100; and 124, 111, 41, 123, 119, 17 when the sample size is 400. The number of iterations not only depends on the choice of the stopping rule, but also depends on different specifications of the EM algorithm, such as the selection of initial values or other parameters for the latent distribution being placed.

Table 2
MSE (bias) of point estimates for $n = 400$.

	MLE	TLE	Bisquare	Mixregt	Mixregt-MCD	MixregL	MixregL-MCD
Case I: $\varepsilon \sim N(0, 1)$							
β_{10}	0.018(−0.006)	0.041(0.012)	0.020(−0.005)	0.019(0.004)	0.027(0.008)	0.025(0.018)	0.031(0.014)
β_{11}	0.020(0.002)	0.108(−0.178)	0.021(−0.001)	0.018(−0.014)	0.028(−0.014)	0.024(−0.028)	0.034(−0.029)
β_{12}	0.018(−0.006)	0.096(−0.171)	0.020(0.000)	0.016(0.008)	0.031(0.012)	0.029(−0.001)	0.042(−0.012)
β_{20}	0.004(0.003)	0.009(0.002)	0.004(0.002)	0.005(−0.006)	0.005(0.012)	0.008(−0.010)	0.008(0.014)
β_{21}	0.004(0.004)	0.007(0.020)	0.004(0.002)	0.004(−0.009)	0.006(−0.002)	0.006(−0.005)	0.009(0.002)
β_{22}	0.004(−0.005)	0.006(0.013)	0.004(−0.006)	0.005(−0.004)	0.006(0.000)	0.007(0.003)	0.008(0.009)
π_1	0.001(0.000)	0.002(−0.001)	0.001(0.002)	0.001(0.001)	0.002(0.005)	0.001(0.000)	0.002(0.006)
Case II: $\varepsilon \sim \text{Laplace}(1)$							
β_{10}	0.022(−0.005)	0.012(0.012)	0.015(−0.003)	0.012(−0.004)	0.013(0.003)	0.010(0.007)	0.012(0.010)
β_{11}	0.014(0.008)	0.013(−0.041)	0.010(0.005)	0.012(0.003)	0.018(−0.013)	0.011(0.005)	0.017(−0.007)
β_{12}	0.016(−0.006)	0.017(−0.050)	0.012(−0.004)	0.011(−0.013)	0.016(0.000)	0.008(−0.007)	0.014(0.005)
β_{20}	0.004(−0.003)	0.003(−0.003)	0.003(−0.003)	0.002(0.001)	0.002(0.000)	0.002(0.002)	0.002(−0.001)
β_{21}	0.004(−0.013)	0.003(0.005)	0.003(−0.015)	0.003(−0.009)	0.004(−0.003)	0.003(−0.004)	0.004(−0.001)
β_{22}	0.004(−0.011)	0.004(0.012)	0.003(−0.009)	0.003(−0.003)	0.004(−0.006)	0.002(−0.003)	0.003(−0.003)
π_1	0.001(0.002)	0.001(0.016)	0.001(0.022)	0.001(0.004)	0.001(0.006)	0.001(0.001)	0.001(0.004)
Case III: $\varepsilon \sim t_1$							
β_{10}	313.757(−0.917)	0.735(−0.040)	0.631(−0.083)	0.147(0.019)	0.154(0.002)	0.016(0.073)	0.017(0.076)
β_{11}	278.219(−3.135)	0.398(0.097)	0.607(−0.187)	0.458(−0.191)	0.485(−0.257)	0.194(−0.352)	0.322(−0.454)
β_{12}	455.172(−1.369)	0.399(0.059)	0.716(−0.146)	0.351(−0.177)	0.484(−0.200)	0.197(−0.361)	0.351(−0.462)
β_{20}	313.757(−0.917)	0.021(−0.001)	0.514(−0.052)	0.023(−0.008)	0.021(−0.002)	0.008(−0.061)	0.008(−0.067)
β_{21}	269.680(−1.135)	0.032(0.003)	0.047(0.034)	0.014(0.006)	0.022(−0.003)	0.011(0.092)	0.015(0.099)
β_{22}	453.695(0.630)	0.093(−0.009)	0.083(0.014)	0.017(0.009)	0.020(−0.002)	0.012(0.094)	0.016(0.102)
π_1	0.061(0.247)	0.008(0.003)	0.016(0.062)	0.009(0.031)	0.008(0.037)	0.037(0.160)	0.038(0.161)
Case IV: $\varepsilon \sim t_3$							
β_{10}	0.301(0.020)	0.037(−0.008)	0.038(−0.010)	0.039(−0.014)	0.059(−0.016)	0.033(0.002)	0.044(0.005)
β_{11}	0.210(−0.046)	0.039(−0.070)	0.044(0.049)	0.034(−0.013)	0.071(−0.008)	0.028(−0.019)	0.049(−0.033)
β_{12}	0.227(−0.049)	0.037(−0.081)	0.034(0.021)	0.046(0.000)	0.045(0.009)	0.031(0.008)	0.048(−0.043)
β_{20}	0.066(0.018)	0.008(−0.017)	0.007(−0.007)	0.006(−0.007)	0.006(0.011)	0.008(−0.011)	0.006(0.008)
β_{21}	0.069(0.055)	0.007(0.001)	0.006(−0.025)	0.007(−0.005)	0.009(−0.008)	0.007(0.003)	0.010(0.005)
β_{22}	0.069(0.055)	0.009(0.006)	0.008(−0.025)	0.008(0.009)	0.010(−0.001)	0.008(0.011)	0.012(0.003)
π_1	0.010(−0.017)	0.002(0.023)	0.002(0.023)	0.002(0.004)	0.003(0.007)	0.002(−0.001)	0.003(0.003)
Case V: $\varepsilon \sim 0.95N(0, 1) + 0.05N(0, 25)$							
β_{10}	0.098(0.000)	0.041(0.005)	0.024(0.004)	0.029(−0.007)	0.038(0.015)	0.034(0.009)	0.042(0.028)
β_{11}	0.394(0.028)	0.048(−0.095)	0.021(0.027)	0.022(0.011)	0.044(−0.012)	0.025(0.003)	0.040(−0.012)
β_{12}	0.081(−0.050)	0.051(−0.119)	0.022(0.014)	0.026(0.001)	0.045(0.012)	0.032(0.000)	0.048(−0.001)
β_{20}	0.041(0.015)	0.006(0.003)	0.005(0.002)	0.006(−0.002)	0.006(0.006)	0.008(−0.006)	0.008(0.003)
β_{21}	0.088(0.046)	0.006(0.010)	0.005(−0.008)	0.006(0.006)	0.009(0.004)	0.008(0.009)	0.011(0.009)
β_{22}	0.135(0.041)	0.007(0.024)	0.004(0.000)	0.005(0.002)	0.008(0.000)	0.007(0.008)	0.011(0.007)
π_1	0.007(−0.033)	0.001(0.003)	0.001(0.006)	0.001(0.000)	0.002(−0.002)	0.002(−0.003)	0.002(−0.007)
Case VI: $\varepsilon \sim N(0, 1)$ with 5% high leverage outliers							
β_{10}	9.355(−1.688)	0.033(0.010)	0.020(−0.010)	1.375(0.246)	0.021(−0.014)	0.013(0.065)	0.029(0.002)
β_{11}	5.188(1.667)	0.049(−0.102)	0.023(−0.011)	2.505(1.479)	0.027(−0.002)	0.014(−0.049)	0.033(−0.037)
β_{12}	4.187(1.307)	0.039(−0.098)	0.021(−0.007)	2.594(1.507)	0.029(0.007)	0.017(−0.034)	0.031(−0.015)
β_{20}	11.697(2.305)	0.005(0.002)	0.004(0.003)	0.005(0.005)	0.005(0.004)	0.007(−0.047)	0.007(−0.002)
β_{21}	11.586(3.309)	0.006(0.011)	0.005(0.012)	0.021(0.125)	0.006(0.004)	0.004(0.026)	0.009(0.005)
β_{22}	12.442(3.437)	0.006(0.003)	0.005(0.003)	0.020(0.122)	0.006(−0.005)	0.005(0.028)	0.010(0.000)
π_1	0.140(0.204)	0.002(0.004)	0.001(−0.006)	0.008(−0.089)	0.001(0.005)	0.004(0.020)	0.001(0.002)

3.2. Sensitivity study based on a real data

In this section, we describe a sensitivity study based on a real data set to compare how outliers affect various different estimation procedures. A typical real data set suitable for mixture regression modeling is the tone data collected in a tone perception experiment of Cohen (1984). In the experiment, a pure fundamental tone was played to a trained musician and electronically generated overtones were added, determined by a stretching ratio (stretchratio). A value of 2 for the stretch ratio corresponds to the harmonic pattern usually heard in traditional definite pitched instruments. The musician was asked to tune an adjustable tone to the octave above the fundamental tone, and a measurement called “tuned” gives the ratio of the adjusted tone to the fundamental. 150 pairs of (tuned, stretchratio) values are obtained for the same musician. The variable “stretchratio” is treated as a response variable and “tuned” as a predictor. The setup of the experiment indicates two mixture components in the model, and the scatter plot of the data collected from the experiment confirms this point. To see the impact of different types of outliers on various procedures, we first add 5 identical pairs, (3, 4), to the original data set

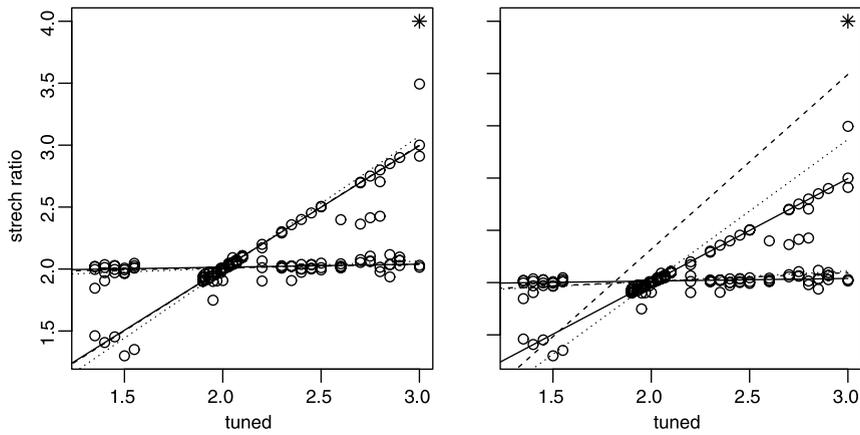


Fig. 1. Mixture linear fitting with outlier (3, 4). Left panel: solid line—Bisquare, dashed line—MixregL, dotted line—Mixregt, Right panel: solid line—Bisquare, dashed—TLE, dotted line—MLE.

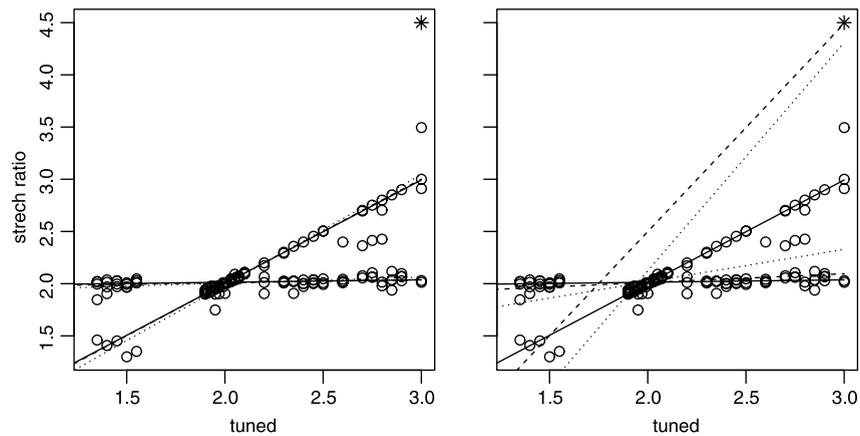


Fig. 2. Mixture linear fitting with outlier (3, 4.5). Left panel: solid line—Bisquare, dashed line—MixregL, dotted line—Mixregt, Right panel: solid line—Bisquare, dashed—TLE, dotted line—MLE.

as outliers in the y -direction. The circles in all plots denote the original data points, and the star denotes the outliers. The right-hand plots in all the figures below have the same y -scales as in the left-hand plots.

The left-hand plot in Fig. 1 clearly shows that the fitting by MixregL and Bisquare are almost identical, and Mixregt also provides a very good fit. For comparison, The Bisquare fit is also drawn in the right-hand plot in Fig. 1, it is quite obvious that the TLE and MLE are affected severely by the outliers. 43 iterations are used for MixregL in this set up to achieve the convergence.

We increase the y -value of the outliers from 4 to 4.5, and the fitting results as seen in Fig. 2 clearly show that the performance of Bisquare, MixregL and Mixregt procedures are quite stable, while the TLE and MLE procedure fits are dragged more severely towards the outliers. For this case, 43 iterations are used for MixregL to achieve convergence.

Then we add 10 identical pairs, (0, 3), to the original data set as high leverage outliers. The left-hand plot in Fig. 3 shows that both Bisquare and MixregL give a reasonable fit, but surprisingly the Mixregt performs less satisfactorily, implying that Mixregt is not quite robust to x -outliers. From the right-hand plot in Fig. 3, we see that MLE has inferior performance against the outliers, and TLE works better. Here, 77 iterations are needed to get convergence for the MixregL procedure.

Finally 10 identical pairs (0, 4) were added to the original data set as both outliers in x and y -direction. The left-hand plot in Fig. 4 shows that Bisquare continues to provide a robust fit, MixregL barely keeps a vague two-line structure, and Mixregt is affected severely by the outliers. The right-hand plot in Fig. 4 shows that MLE is still the worst, and TLE works fine. 22 iterations were used for MixregL to achieve convergence.

In all the scenarios, the Bisquare performed uniformly better than other fitting procedures, although the simulation studies show that Bisquare is less satisfying in some cases, such as in $\varepsilon \sim t$ -distributions. Instead of modifying the log likelihood objective function, the Bisquare procedure tries to modify the existing EM algorithm for mixture regression models by replacing the least squares criterion with a robust criteria in the M step. See Bai et al. (2012) for detail. Generally MixregL performed better than Mixregt, but both procedures are not quite robust to the high leverage outliers. We also applied

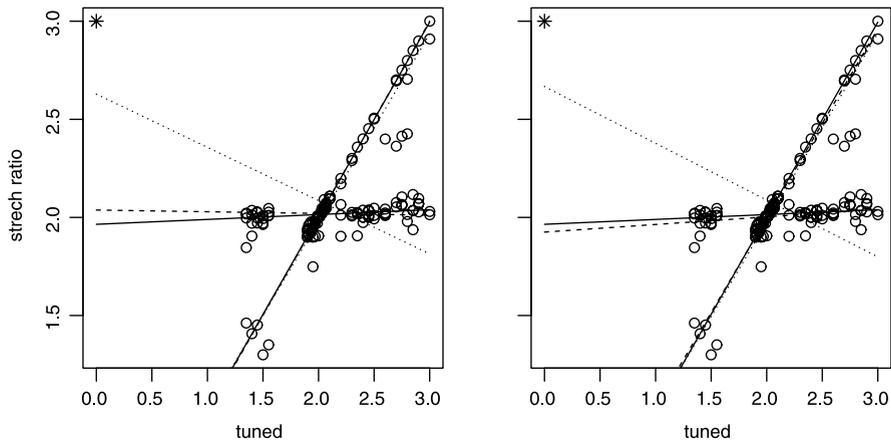


Fig. 3. Mixture linear fitting with outlier (0, 3). Left panel: solid line—Bisquare, dashed line—MixregL, dotted line—Mixregt, Right panel: solid line—Bisquare, dashed—TLE, dotted line—MLE.

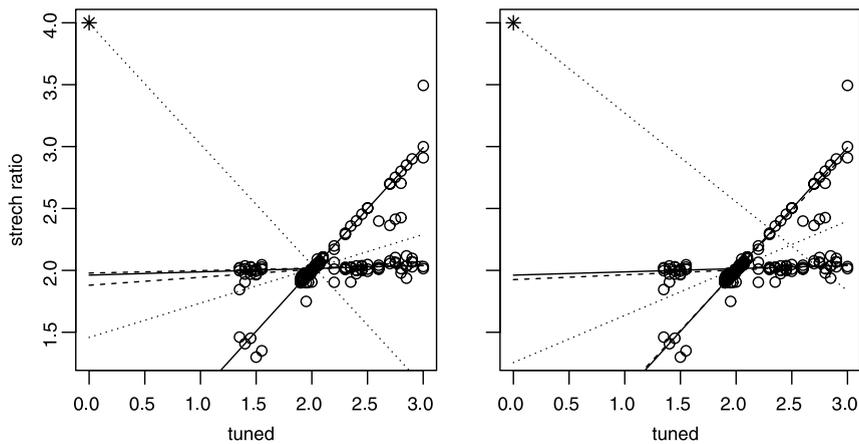


Fig. 4. Mixture linear fitting with outlier (0, 4). Left panel: solid line—Bisquare, dashed line—MixregL, dotted line—Mixregt, Right panel: solid line—Bisquare, dashed—TLE, dotted line—MLE.

Mixregt-MCD and MixregL-MCD to the data set. Both procedures can successfully remove the high leverage outliers and give similar results to the Bisquare.

4. Conclusion

A new robust estimation procedure tailored to mixture linear regression models is proposed by assuming the random error has a Laplace distribution. The robustness is achieved essentially by the LAD procedure, and is implemented with an EM algorithm. Efficiency and effectiveness of the proposed EM algorithm relies upon the fact that the Laplace distribution is a scale mixture of a normal distribution and a distribution related to exponential. The simulation study shows that the proposed method is superior to or comparable to existing robust estimation procedures in all simulation setups. However, the real data example shows that when high leverage outliers exist, the trimmed version of the proposed procedure should be used.

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