

Maximum likelihood estimation of a multidimensional log-concave density

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Summary. Let X_1, \dots, X_n be independent and identically distributed random vectors with a log-concave (Lebesgue) density f . We first prove that, with probability one, there exists a unique maximum likelihood estimator \hat{f}_n of f . The use of this estimator is attractive because, unlike kernel density estimation, the method is fully automatic, with no smoothing parameters to choose. Although the existence proof is non-constructive, we are able to reformulate the issue of computing \hat{f}_n in terms of a non-differentiable convex optimisation problem, and thus combine techniques of computational geometry with Shor's r -algorithm to produce a sequence that converges to \hat{f}_n . For the moderate or large sample sizes in our simulations, the maximum likelihood estimator is shown to provide an improvement in performance compared with kernel-based methods, even when we allow the use of a theoretical, optimal fixed bandwidth for the kernel estimator that would not be available in practice. We also present a real data clustering example, which shows that our methodology can be used in conjunction with the Expectation–Maximisation (EM) algorithm to fit finite mixtures of log-concave densities. An R version of the algorithm is available in the package `LogConcDEAD` – Log-Concave Density Estimation in Arbitrary Dimensions.

Keywords: Computational geometry, log-concavity, maximum likelihood estimation, non-differentiable convex optimisation, nonparametric density estimation, Shor's r -algorithm

1. Introduction

Modern nonparametric density estimation began with the introduction of a kernel density estimator in the pioneering work of Fix and Hodges (1951), later republished as Fix and Hodges (1989). For independent and identically distributed real-valued observations, the appealing asymptotic theory of the mean integrated squared error was provided by Rosenblatt (1956) and Parzen (1962). This theory leads to an asymptotically optimal choice of the smoothing parameter, or bandwidth. Unfortunately, however, it depends on the unknown density f through the integral of the square of the second derivative of f . Considerable effort has therefore been focused on finding methods of automatic bandwidth selection (cf. Wand and Jones, 1995, Chapter 3, and the references therein). Although this has resulted in algorithms, e.g. Chiu (1992), that achieve the optimal rate of convergence of the relative error, namely $O_p(n^{-1/2})$, where n is the sample size, good finite sample performance is by no means guaranteed.

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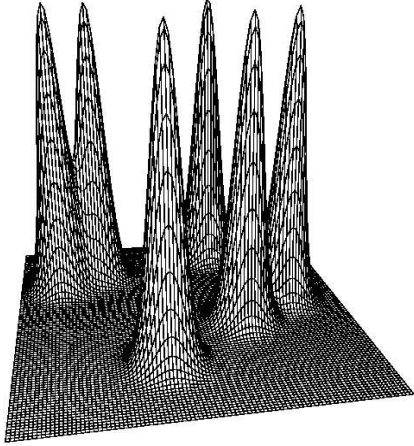


Fig. 1. Without any shape constraint on the class of densities, the likelihood function is unbounded, because we can take successively close approximations to a mixture of n ‘spikes’ (one on each X_i).

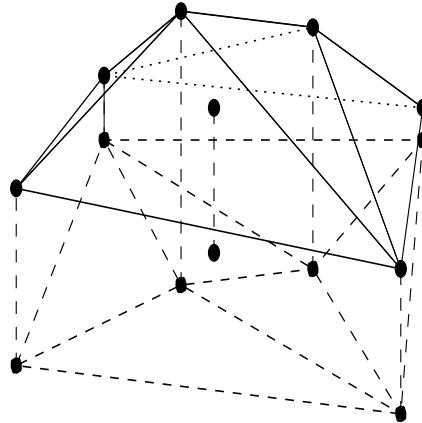


Fig. 2. The ‘tent-like’ structure of the graph of the logarithm of the maximum likelihood estimator for bivariate data.

This problem is compounded when the observations take values in \mathbb{R}^d , where the general kernel estimator (Deheuvels, 1977) requires the specification of a symmetric, positive definite $d \times d$ bandwidth matrix. The difficulties involved in making the $d(d+1)/2$ choices for its entries mean that attention is often restricted either to bandwidth matrices that are diagonal, or even to those that are scalar multiples of the identity matrix. Of course, practical issues of automatic bandwidth selection remain.

In this paper, we propose a fully automatic nonparametric estimator of f , with no tuning parameters to be chosen, under the condition that f is log-concave – that is, $\log f$ is a concave function. The class of log-concave densities has many attractive properties and has been well-studied, particularly in the economics, sampling and reliability theory literature. See Section 2 for further discussion of examples, applications and properties of log-concave densities.

In Section 3, we show that if X_1, \dots, X_n are independent and identically distributed random vectors with a log-concave density, then with probability one there exists a unique log-concave density \hat{f}_n that maximises the likelihood function,

$$L(f) = \prod_{i=1}^n f(X_i).$$

Before continuing, it is worth noting that without any shape constraints on the densities under consideration, the likelihood function is unbounded. To see this, we could define a sequence (f_n) of densities that represent successively close approximations to a mixture of n ‘spikes’ (one on each X_i), such as $f_n(x) = n^{-1} \sum_{i=1}^n \phi_{d, n^{-1}I}(x - X_i)$, where $\phi_{d, \Sigma}$ denotes the $N_d(0, \Sigma)$ density. This sequence satisfies $L(f_n) \rightarrow \infty$ as $n \rightarrow \infty$ (cf. Figure 1). In fact, a modification of this argument may be used to show that the likelihood function remains unbounded even if we restrict attention to unimodal densities.

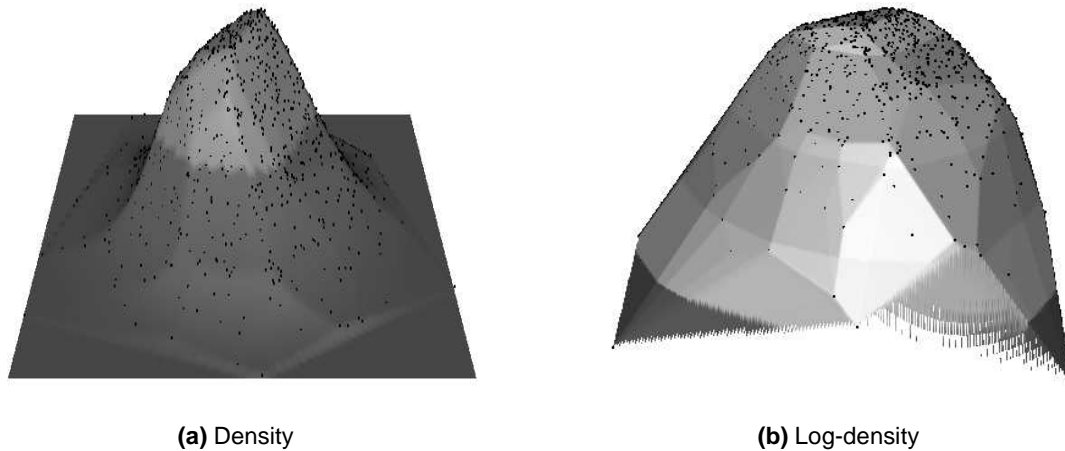


Fig. 3. Log-concave maximum likelihood estimates based on 1000 observations (plotted as dots) from a standard bivariate normal distribution.

Figure 2 gives a diagram illustrating the structure of the maximum likelihood estimator on the logarithmic scale. This structure is most easily visualised for two-dimensional data, where one can imagine associating a ‘tent pole’ with each observation, extending vertically out of the plane. For certain tent pole heights, the graph of the logarithm of the maximum likelihood estimator can be thought of as the roof of a taut tent stretched over the tent poles. The fact that the logarithm of the maximum likelihood estimator is of this ‘tent function’ form constitutes part of the proof of its existence and uniqueness.

In Section 4, we discuss the computational problem of how to adjust the n tent pole heights so that the corresponding tent functions converge to the logarithm of the maximum likelihood estimator. One reason that this computational problem is so challenging in more than one dimension is the fact that it is difficult to describe the set of tent pole heights that correspond to concave functions. The key observation, discussed in Section 4, is that it is possible to minimise a modified objective function that it is convex (though non-differentiable). This allows us to apply the powerful non-differentiable convex optimisation methodology of the subgradient method (Shor, 1985) and a variant called Shor’s r -algorithm, which has been implemented by Kappel and Kuntsevich (2000).

As an illustration of the estimates obtained, Figure 3 presents plots of the maximum likelihood estimator, and its logarithm, for 1000 observations from a standard bivariate normal distribution. These plots were created using the `LogConcDEAD` package (Cule *et al.*, 2008a) in R (R Development Core Team, 2008), which exploits the interactive surface-plotting software available in the `rgl` package (Adler and Murdoch, 2007).

In Section 5 we present simulations to compare the finite-sample performance of the maximum likelihood estimator with kernel-based methods. The results are striking: even when we use the theoretical, optimal bandwidth for the kernel estimator (or an asymptotic approximation to this when it is not available), we find that the maximum likelihood estimator has a rather smaller mean integrated squared error for moderate or large sample sizes, despite the fact that this optimal

bandwidth depends on properties of the density that would be unknown in practice. This suggests that the maximum likelihood estimator is able to adapt to the local smoothness of the underlying density automatically.

Nonparametric density estimation is a fundamental tool for the visualisation of structure in exploratory data analysis, and has an enormous literature that includes the monographs of Devroye and Györfi (1985), Silverman (1986), Scott (1992) and Wand and Jones (1995). Our proposed method may certainly be used for this purpose; however, it may also be used as an intermediary stage in more involved statistical procedures. For instance:

- (a) In classification problems, we have $p \geq 2$ populations of interest, and assume in this discussion that these have densities f_1, \dots, f_p on \mathbb{R}^d . We observe training data of the form $\{(X_i, Y_i) : i = 1, \dots, n\}$, where if $Y_i = j$, then X_i has density f_j . The aim is to classify a new observation $z \in \mathbb{R}^d$ as coming from one of the populations. Problems of this type occur in a huge variety of applications, including medical diagnosis, archaeology, ecology etc. – see Gordon (1981), Hand (1981) or Devroye *et al.* (1996) for further details and examples. A natural approach to classification problems is to construct density estimates $\hat{f}_1, \dots, \hat{f}_p$, where \hat{f}_j is based on the n_j observations, say, from the j th population, namely $\{X_i : Y_i = j\}$. We may then assign z to the j th population if $n_j \hat{f}_j(z) = \max\{n_1 \hat{f}_1(z), \dots, n_p \hat{f}_p(z)\}$. In this context, the use of kernel-based estimators in general requires the choice of p separate $d \times d$ bandwidth matrices, while the corresponding procedure based on the log-concave maximum likelihood estimates is again fully automatic.
- (b) Clustering problems are closely related to the classification problems described above. The difference is that, in the above notation, we do not observe Y_1, \dots, Y_n , and have to assign each of X_1, \dots, X_n to one of the p populations. A common technique is based on fitting a mixture density of the form $f(x) = \sum_{j=1}^p \pi_j f_j(x)$, where the mixture proportions π_1, \dots, π_p are positive and sum to one. Under the assumption that each of the component densities f_1, \dots, f_p is log-concave, we show in Section 6 that our methodology can be extended to fit such a finite mixture density, which need not itself be log-concave – cf. Section 2. We also illustrate this clustering algorithm on a Wisconsin breast cancer data set in Section 6, where the aim is to separate observations into benign and malignant component populations.
- (c) A functional of the true underlying density may be estimated by the corresponding functional of a density estimator, such as the log-concave maximum likelihood estimator. Examples of functionals of interest include probabilities, such as $\int_{\|x\| \geq 1} f(x) dx$, moments, e.g. $\int \|x\|^2 f(x) dx$, and the differential entropy, $-\int f(x) \log f(x) dx$. It may be possible to compute the plug-in estimator based on the log-concave maximum likelihood estimator analytically, but in Section 7, we show that even if this is not possible, in many cases of interest we can sample from the log-concave maximum likelihood estimator \hat{f}_n , and hence obtain a Monte Carlo estimate of the functional. This nice feature also means that the log-concave maximum likelihood estimator can be used in a Monte Carlo bootstrap procedure for assessing uncertainty in functional estimates – see Section 7 for further details.
- (d) The fitting of a nonparametric density estimate may give an indication of the validity of a particular smaller model (often parametric). Thus, a contour plot of the log-concave maximum likelihood estimator may provide evidence that the underlying density has elliptical contours, and thus suggest that a model that exploits this elliptical symmetry.
- (e) In the univariate case, Walther (2002) describes methodology based on log-concave density estimation for addressing the problem of detecting the presence of mixing in a distribution.

As an application, he cites the Pickering/Platt debate (Swales, 1985) on the issue of whether high blood pressure is a disease (in which case observed blood pressure measurements should follow a mixture distribution), or simply a label attached to people in the right tail of the blood pressure distribution. As a result of our algorithm for computing the multidimensional log-concave maximum likelihood estimator, this methodology extends immediately to more than one dimension.

There has been considerable recent interest in shape-restricted nonparametric density estimation, but most of it has been confined to the case of univariate densities, where the computational algorithms are more straightforward. Nevertheless, as was discussed above, it is in multivariate situations that the automatic nature of the maximum likelihood estimator is particularly valuable. Walther (2002), Dümbgen and Rufibach (2007) and Pal *et al.* (2007) have proved the existence and uniqueness of the log-concave maximum likelihood estimator in one dimension and Dümbgen and Rufibach (2007), Pal *et al.* (2007) and Balabdaoui *et al.* (2008) have studied its theoretical properties. Rufibach (2007) has compared different algorithms for computing the univariate estimator, including the iterative convex minorant algorithm (Groeneboom and Wellner, 1992; Jongbloed, 1998), and three others. Dümbgen *et al.* (2007) also present an Active Set algorithm, which has similarities with the vertex direction and vertex reduction algorithms described in Groeneboom *et al.* (2008). For univariate data, it is also well-known that there exist maximum likelihood estimators of a non-increasing density supported on $[0, \infty)$ (Grenander, 1956) and of a convex, decreasing density (Groeneboom *et al.*, 2001).

In Section 8, we give a brief concluding discussion, and suggest some directions for future research. Finally, we present in Appendix A a glossary of terms and results from convex analysis and computational geometry that appear in italics at their first occurrence in the main body of the paper; the references are Rockafellar (1997) and Lee (1997). Proofs are deferred to Appendix B, except that the beginning of the proof of Theorem 2 is given in the main text, as the ideas and notation introduced are needed in the remainder of the paper.

2. Log-concave densities: examples, applications and properties

Many of the most commonly-encountered parametric families of univariate distributions have *log-concave* densities, including the family of normal distributions, gamma distributions with shape parameter at least one, Beta(α, β) distributions with $\alpha, \beta \geq 1$, Weibull distributions with shape parameter at least one, Gumbel, logistic and Laplace densities; see Bagnoli and Bergstrom (1989) for other examples. Univariate log-concave densities are unimodal and have fairly light tails – it may help to think of the exponential distribution (where the logarithm of the density is a linear function on the positive half-axis) as a borderline case. Thus Cauchy, Pareto and lognormal densities, for instance, are not log-concave. Mixtures of log-concave densities may be log-concave, but in general they are not; for instance, for $p \in (0, 1)$, the location mixture of standard univariate normal densities $f(x) = p\phi(x) + (1 - p)\phi(x - \mu)$ is log-concave if and only if $\|\mu\| \leq 2$.

The assumption of log-concavity is a popular one in economics; Caplin and Naelbuff (1991b) show that in the theory of elections and under a log-concavity assumption, the proposal most preferred by the mean voter is unbeatable under a 64% majority rule. As another example, in the theory of imperfect competition, Caplin and Naelbuff (1991a) use log-concavity of the density of consumers'

utility parameters as a sufficient condition in their proof of the existence of a pure-strategy price equilibrium for any number of firms producing any set of products. See Bagnoli and Bergstrom (1989) for many other applications of log-concavity to economics. Brooks (1998) and Mengersen and Tweedie (1996) have exploited the properties of log-concave densities in studying the convergence of Markov chain Monte Carlo sampling procedures.

An (1998) lists many useful properties of log-concave densities. For instance, if f and g are (possibly multidimensional) log-concave densities, then their convolution $f * g$ is log-concave. In other words, if X and Y are independent and have log-concave densities, then their sum $X + Y$ has a log-concave density. The class of log-concave densities is also closed under the taking of pointwise limits. One-dimensional log-concave densities have increasing hazard functions, which is why they are of interest in reliability theory. Moreover, Ibragimov (1956) proved the following characterisation: a univariate density f is log-concave if and only if the convolution $f * g$ is unimodal for every unimodal density g . There is no natural generalisation of this result to higher dimensions.

As was mentioned in Section 1, this paper concerns multidimensional log-concave densities, for which fewer properties are known. It is therefore of interest to understand how the property of log-concavity in more than one dimension relates to the univariate notion. Our first proposition below is intended to give some insight into this issue. It is not formally required for the subsequent development of our methodology in Sections 3 and 4, although we did apply the result when designing our simulation study in Section 5. We assume throughout that log-concave densities are with respect to Lebesgue measure on the *affine hull* of their *support*, and ‘ X has a log-concave density’ means ‘there exists a version of the density of X that is log-concave’.

PROPOSITION 1. *Let X be a d -variate random vector having density f with respect to Lebesgue measure on \mathbb{R}^d . For a subspace V of \mathbb{R}^d , let $P_V(x)$ denote the orthogonal projection of x onto V . Then in order that f be log-concave, it is:*

- (a) *necessary that for any subspace V , the marginal density of $P_V(X)$ is log-concave and the conditional density $f_{X|P_V(X)}(\cdot|t)$ of X given $P_V(X) = t$ is log-concave for each t*
- (b) *sufficient that for every $(d-1)$ -dimensional subspace V , the conditional density $f_{X|P_V(X)}(\cdot|t)$ of X given $P_V(X) = t$ is log-concave for each t .*

The part of Proposition 1(a) concerning marginal densities is an immediate consequence of Theorem 6 of Prékopa (1973). One can regard Proposition 1(b) as saying that a multidimensional density is log-concave if the restriction of the density to any line is a (univariate) log-concave function.

It is interesting to compare the properties of log-concave densities presented in Proposition 1 with the corresponding properties of Gaussian densities. In fact, Proposition 1 remains true if we replace ‘log-concave’ with ‘Gaussian’ throughout (at least, provided that in part (b) we also assume there is a point at which f is twice differentiable). These shared properties suggest that the class of log-concave densities is a natural, infinite-dimensional generalisation of the class of Gaussian densities.

3. Existence, uniqueness and structure of the maximum likelihood estimator

Let \mathcal{F}_0 denote the class of log-concave densities on \mathbb{R}^d with d -dimensional support, and let $f_0 \in \mathcal{F}_0$. The degenerate case where the support is of dimension smaller than d can also be handled, but for simplicity of exposition we concentrate on the non-degenerate case. Suppose that X_1, \dots, X_n are a random sample from f_0 . We say that $\hat{f}_n = \hat{f}_n(X_1, \dots, X_n) \in \mathcal{F}_0$ is a (nonparametric) *maximum likelihood estimator* of f_0 if it maximises $\ell(f) = \sum_{i=1}^n \log f(X_i)$ over $f \in \mathcal{F}_0$.

THEOREM 2. *Suppose that $n \geq d + 1$. Then, with probability one, a nonparametric maximum likelihood estimator \hat{f}_n of f_0 exists and is unique.*

FIRST PART OF PROOF. We may assume that X_1, \dots, X_n are distinct and their convex hull, $C_n = \text{conv}(X_1, \dots, X_n)$, is a d -dimensional *polytope* (an event of probability one when $n \geq d + 1$). By a standard argument in convex analysis (Rockafellar, 1997, p. 37), for each $y = (y_1, \dots, y_n) \in \mathbb{R}^n$ there exists a function $\bar{h}_y : \mathbb{R}^d \rightarrow \mathbb{R}$ with the property that \bar{h}_y is the least concave function satisfying $\bar{h}_y(X_i) \geq y_i$ for all $i = 1, \dots, n$. Informally, \bar{h}_y is a ‘tent function’, and a typical example is depicted in Figure 2. Let $\mathcal{H} = \{\bar{h}_y : y \in \mathbb{R}^n\}$ denote ‘the class of tent functions’. Let \mathcal{F} denote the set of all log-concave functions on \mathbb{R}^d , and for $f \in \mathcal{F}$, define

$$\psi_n(f) = \frac{1}{n} \sum_{i=1}^n \log f(X_i) - \int_{\mathbb{R}^d} f(x) dx.$$

Suppose that f maximises $\psi_n(\cdot)$ over \mathcal{F} . The main part of the proof, which is completed in the Appendix, consists of showing that

- (i) $f(x) > 0$ for $x \in C_n$
- (ii) $f(x) = 0$ for $x \notin C_n$
- (iii) $\log f \in \mathcal{H}$
- (iv) $f \in \mathcal{F}_0$
- (v) there exists $M > 0$ such that if $\max_i |\bar{h}_y(X_i)| \geq M$, then $\psi_n(\exp(\bar{h}_y)) \leq \psi_n(f)$.

Although step (iii) above gives us a finite-dimensional class of functions to which $\log \hat{f}_n$ belongs, the proof of Theorem 2 gives no indication of how to find the member of this class that maximises the likelihood function. We therefore seek an iterative algorithm to compute the estimator, but first we describe the structure we see in Figure 2 in Section 1 more precisely. From now on, we assume:

(A1): $n \geq d + 1$, and every subset of $\{X_1, \dots, X_n\}$ of size $d + 1$ is *affinely independent*.

Note that when $n \geq d + 1$, the event in **(A1)** has probability one. From step (iii) in the proof of Theorem 2 above, there exists $y \in \mathbb{R}^n$ such that $\log \hat{f}_n = \bar{h}_y$. As illustrated in Figure 2, and justified formally by Corollary 17.1.3 and Corollary 19.1.2 of Rockafellar (1997), the convex hull of the data, C_n , may be *triangulated* in such a way that $\log \hat{f}_n$ coincides with an *affine function* on each *simplex* in the triangulation. In other words, if $j = (j_1, \dots, j_{d+1})$ is a $(d + 1)$ -tuple of distinct indices in $\{1, \dots, n\}$, and $C_{n,j} = \text{conv}(X_{j_1}, \dots, X_{j_{d+1}})$, then there exists a finite set J consisting of m such $(d + 1)$ -tuples, with the following three properties:

- (i) $\cup_{j \in J} C_{n,j} = C_n$
- (ii) the *relative interiors* of the sets $\{C_{n,j} : j \in J\}$ are pairwise disjoint
- (iii)

$$\log \hat{f}_n(x) = \begin{cases} \langle x, b_j \rangle - \beta_j & \text{if } x \in C_{n,j} \text{ for some } j \in J \\ -\infty & \text{if } x \notin C_n \end{cases}$$

for some $b_1, \dots, b_m \in \mathbb{R}^d$ and $\beta_1, \dots, \beta_m \in \mathbb{R}$. Here and below, $\langle \cdot, \cdot \rangle$ denotes the usual Euclidean inner product in \mathbb{R}^d .

In the iterative algorithm that we propose in Section 4 for computing the maximum likelihood estimator, we need to find convex hulls and triangulations at each iteration. Fortunately, these can be computed efficiently using the `Quickhull` algorithm of Barber *et al.* (1996).

4. Computation of the maximum likelihood estimator

4.1. Reformulation

As a first attempt to find an algorithm which produces a sequence that converges to the maximum likelihood estimator in Theorem 2, it is natural to try to minimise numerically the function

$$\tau(y_1, \dots, y_n) = -\frac{1}{n} \sum_{i=1}^n \bar{h}_y(X_i) + \int_{C_n} \exp\{\bar{h}_y(x)\} dx.$$

Although this approach might work in principle, one difficulty is that τ is not convex, so this approach is extremely computationally intensive, even with relatively few observations. Another reason for the numerical difficulties stems from the fact that the set of y -values on which τ attains its minimum is rather large: in general it may be possible to alter particular components y_i without changing \bar{h}_y . Of course, we could have defined τ as a function of \bar{h}_y rather than as a function of the vector of tent pole heights $y = (y_1, \dots, y_n)$. Our choice, however, motivates the following definition of a modified objective function:

$$\sigma(y_1, \dots, y_n) = -\frac{1}{n} \sum_{i=1}^n y_i + \int_{C_n} \exp\{\bar{h}_y(x)\} dx. \quad (4.1)$$

The great advantages of minimising σ rather than τ are seen by the following theorem.

THEOREM 3. *Assume (A1). The function σ is a convex function satisfying $\sigma \geq \tau$. It has a unique minimum at $y^* \in \mathbb{R}^n$, say, and $\log \hat{f}_n = \bar{h}_{y^*}$.*

Thus Theorem 3 shows that the unique minimum $y^* = (y_1^*, \dots, y_n^*)$ of σ belongs to the minimum set of τ . In fact, it corresponds to the element of the minimum set for which $\bar{h}_{y^*}(X_i) = y_i^*$ for $i = 1, \dots, n$. Informally, then, \bar{h}_{y^*} is ‘a tent function with all of the tent poles touching the tent’.

In order to compute the function σ at a generic point $y = (y_1, \dots, y_n) \in \mathbb{R}^n$, we need to be able to evaluate the integral in (4.1). In the notation of Section 3, we may write

$$\int_{C_n} \exp\{\bar{h}_y(x)\} dx = \sum_{j \in J} \int_{C_{n,j}} \exp\{\langle x, b_j \rangle - \beta_j\} dx.$$

For each $j = (j_1, \dots, j_{d+1}) \in J$, let A_j be the $d \times d$ matrix whose l th column is $X_{j_{l+1}} - X_{j_1}$ for $l = 1, \dots, d$, and let $\alpha_j = X_{j_1}$. Then the affine transformation $w \mapsto A_j w + \alpha_j$ takes the unit simplex $T_d = \{w = (w_1, \dots, w_d) : w_l \geq 0, \sum_{l=1}^d w_l \leq 1\}$ to $C_{n,j}$. Letting $z_{j,l} = y_{j_{l+1}} - y_{j_1}$, we can then establish by a simple change of variables and induction on d that if $z_{j,1}, \dots, z_{j,d}$ are non-zero and distinct, then

$$\int_{C_n} \exp\{\bar{h}_y(x)\} dx = \sum_{j \in J} |\det A_j| e^{y_{j_1}} \sum_{r=1}^d \frac{e^{z_{j,r}} - 1}{z_{j,r}} \prod_{\substack{1 \leq s \leq d \\ s \neq r}} \frac{1}{z_{j,r} - z_{j,s}}. \quad (4.2)$$

Further details of this calculation can be found in a longer version of this paper (Cule *et al.*, 2008b). The singularities that occur when some of $z_{j,1}, \dots, z_{j,d}$ may be zero or equal are removable. Thus, although (4.2) is a little complicated, it allows the computation of our objective function.

4.2. Nonsmooth optimisation

There is a vast literature on techniques of convex optimisation (cf. Boyd and Vandenberghe (2004), for example), including the method of steepest descent and Newton's method. Unfortunately, these methods rely on the differentiability of the objective function, and the function σ is not differentiable. This can be seen informally by studying the schematic diagram in Figure 2 again. If the i th tent pole, say, is touching but not critically supporting the tent, then decreasing the height of this tent pole does not change the tent function, and thus does not alter the integral in (4.1); on the other hand, increasing the height of the tent pole does alter the tent function and therefore the integral in (4.1). This argument may be used to show that at such a point, the i th partial derivative of σ does not exist.

The set of points at which σ is not differentiable constitute a set of Lebesgue measure zero, but the non-differentiability cannot be ignored in our optimisation procedure. Instead, it will be necessary to derive a *subgradient* of σ at each point $y \in \mathbb{R}^n$. This derivation, along with a more formal discussion of the non-differentiability of σ , can be found in the Appendix.

The theory of non-differentiable, convex optimisation is perhaps less well-known than its differentiable counterpart, but a fundamental contribution was made by Shor (1985) with his introduction of the subgradient method for minimising non-differentiable, convex functions defined on Euclidean spaces. A slightly specialised version of his Theorem 2.2 gives that if $\partial\sigma(y)$ is a subgradient of σ at y , then for any $y^{(0)} \in \mathbb{R}^n$, the sequence generated by the formula

$$y^{(\ell+1)} = y^{(\ell)} - h_{\ell+1} \frac{\partial\sigma(y^{(\ell)})}{\|\partial\sigma(y^{(\ell)})\|}$$

has the property that either there exists an index ℓ^* such that $y^{(\ell^*)} = y^*$, or $y^{(\ell)} \rightarrow y^*$ and $\sigma(y^{(\ell)}) \rightarrow \sigma(y^*)$ as $\ell \rightarrow \infty$, provided we choose the step lengths h_ℓ so that $h_\ell \rightarrow 0$ as $\ell \rightarrow \infty$, but $\sum_{\ell=1}^{\infty} h_\ell = \infty$.

Shor recognised, however, that the convergence of this algorithm could be slow in practice, and that although appropriate step size selection could improve matters somewhat, the convergence would never be better than linear (compared with quadratic convergence for Newton's method near the

optimum – see Boyd and Vandenberghe (2004, Section 9.5)). Slow convergence can be caused by taking at each stage a step in a direction nearly orthogonal to the direction towards the optimum, which means that simply adjusting the step size selection scheme will never produce the desired improvements in convergence rate.

One solution (Shor, 1985, Chapter 3) is to attempt to shrink the angle between the subgradient and the direction towards the minimum through a (necessarily nonorthogonal) linear transformation, and perform the subgradient step in the transformed space. By analogy with Newton’s method for smooth functions, an appropriate transformation would be an approximation to the inverse of the Hessian matrix at the optimum. This is not possible for nonsmooth problems, because the inverse might not even exist (and will not exist at points at which the function is not differentiable, which may include the optimum).

Instead, we perform a sequence of dilations in the direction of the difference between two successive subgradients, in the hope of improving convergence in the worst-case scenario of steps nearly perpendicular to the direction towards the minimiser. This variant, which has become known as Shor’s r -algorithm, has been implemented in Kappel and Kuntsevich (2000). Accompanying software `SolvOpt` is available from <http://www.uni-graz.at/imawww/kuntsevich/solvopt/>.

Although the formal convergence of the r -algorithm has not been proved, we agree with the authors’ claims that it is robust, efficient and accurate. Of course, it is clear that if we terminate the r -algorithm after any finite number of steps and apply the original Shor algorithm using our terminating value of y as the new starting value, then formal convergence is guaranteed. We have not found it necessary to run the original Shor algorithm after termination of the r -algorithm in practice.

If $(y^{(\ell)})$ denotes the sequence of vectors in \mathbb{R}^n produced by the r -algorithm, we terminate when

- $|\sigma(y^{(\ell+1)}) - \sigma(y^{(\ell)})| \leq \delta$
- $|y_i^{(\ell+1)} - y_i^{(\ell)}| \leq \epsilon$ for $i = 1, \dots, n$
- $|1 - \int \exp\{\bar{h}_{y^{(\ell)}}(x)\} dx| \leq \eta$

for some small δ, ϵ and $\eta > 0$. The first two termination criteria follow Kappel and Kuntsevich (2000), while the third is based on our knowledge that the true optimum corresponds to a density (Section 3). As default values, and throughout this paper, we took $\delta = 10^{-8}$ and $\epsilon = \eta = 10^{-4}$.

Table 1 gives approximate running times and number of iterations of Shor’s r -algorithm required for different sample sizes and dimensions on an ordinary desktop computer (1.8GHz, 2GB RAM). Unsurprisingly, the running time increases relatively quickly with the sample size, while the number of iterations increases approximately linearly with n . Each iteration takes longer as the dimension increases, though it is interesting to note that the number of iterations required for the algorithm to terminate decreases as the dimension increases. When $d = 1$, we recommend the Active Set algorithm of Dümbgen *et al.* (2007), which is implemented in the R package `logcondens` (Rufibach and Dümbgen, 2006).

Table 1. Approximate running times (with number of iterations in brackets) for computing the maximum likelihood estimator of a log-concave density

	$n = 100$	$n = 500$	$n = 1000$	$n = 2000$
$d = 2$	1.5 secs (260)	50 secs (1270)	4 mins (2540)	24 mins (5370)
$d = 3$	6 secs (170)	100 secs (820)	7 mins (1530)	44 mins (2740)
$d = 4$	23 secs (135)	670 secs (600)	37 mins (1100)	224 mins (2060)

5. Finite sample performance

Our simulation study considered, for $d = 2$ and 3, the following densities:

- (a) standard normal, $\phi_d \equiv \phi_{d,I}$
 - (b) dependent normal, $\phi_{d,\Sigma}$, with $\Sigma_{ij} = \mathbb{1}_{\{i=j\}} + 0.2\mathbb{1}_{\{i \neq j\}}$
 - (c) the joint density of independent $\Gamma(2, 1)$ components
 - (d-f) the normal location mixture $0.6\phi_d(\cdot) + 0.4\phi_d(\cdot - \mu)$ for (d) $\|\mu\| = 1$, (e) $\|\mu\| = 2$, (f) $\|\mu\| = 3$.
- An application of Proposition 1 gives that such a normal location mixture is log-concave if and only if $\|\mu\| \leq 2$.

In Tables 2 and 3 we present, for each density and for four different sample sizes, an estimate of the mean integrated squared error (MISE) of the nonparametric maximum likelihood estimator based on 100 Monte Carlo iterations. We also show the MISE for the kernel density estimates with a Gaussian kernel and, for all of the normal and mixture of normal examples, the choice of bandwidth that minimises the MISE. In the gamma example, exact MISE calculations are not possible, so we took the bandwidth that minimises the asymptotic mean integrated squared error (AMISE). These optimal bandwidths can be computed using the formulae in Wand and Jones (1995, Sections 4.3 and 4.4). As minimisation of the expressions for both the MISE and the AMISE requires knowledge of certain functionals of the true density that would be unknown in practice, we also provide a comparison with an empirical bandwidth selector based on least squares cross validation (LSCV) (Wand and Jones, 1995, Section 4.7). The LSCV bandwidths were computed using the `ks` package (Duong, 2007) in R, and we used the option of constraining the bandwidth matrices to be diagonal in cases (a) and (c) where the components are independent.

We see that in cases (a)-(e) the log-concave maximum likelihood estimator has a smaller MISE than the kernel estimate with bandwidth chosen by LSCV, and at least for moderate and large sample sizes, the difference is quite dramatic. Even more remarkably, in these cases the log-concave estimator also outperforms the kernel estimate with optimally chosen bandwidth when the sample size is not too small. It seems that for small sample sizes, the fact that the convex hull of the data is rather small hinders the performance of the log-concave estimator, but that this effect is reduced as the sample size increases. The log-concave estimator copes well with the dependence in case (b), and it also deals particularly impressively with case (c), where the true density decays to zero at the boundary of the positive orthant.

In case (f), where the log-concavity assumption is violated, the performance of our estimator is not as good as the kernel estimate with the optimally chosen bandwidth, but is still comparable in most cases with the LSCV method. One would not expect the MISE of \hat{f}_n to approach zero as $n \rightarrow \infty$ if log-concavity is violated, and in fact we conjecture that in this case the log-concave maximum

Table 2. Mean integrated squared error estimates (with standard errors in brackets where applicable; $d = 2$)

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00620(0.000222)	0.00431	0.00622(0.000383)
500	0.00161(0.0000514)	0.00164	0.00199(0.0000844)
1000	0.000983(0.0000289)	0.00106	0.00122(0.0000495)
2000	0.000599(0.0000155)	0.000686	0.000803(0.0000276)

(a) Independent Normal

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00607(0.000283)	0.00440	0.00827(0.000583)
500	0.00168(0.0000573)	0.00167	0.00240(0.000122)
1000	0.00100(0.0000295)	0.00108	0.00142(0.0000662)
2000	0.000608(0.0000154)	0.000700	0.000868(0.0000331)

(b) Dependent Normal

n	LogConcDEAD	Kernel (opt AMISE)	Kernel (LSCV)
100	0.00588(0.000222)	0.00644	0.00800(0.000339)
500	0.00143(0.0000478)	0.00220	0.00291(0.0000687)
1000	0.000802(0.0000236)	0.00139	0.00194(0.0000456)
2000	0.000451(0.0000110)	0.000874	0.00130(0.0000209)

(c) $\Gamma(2, 1)$ (independent components)

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00504(0.000206)	0.00384	0.00515(0.000195)
500	0.00136(0.0000745)	0.00145	0.00179(0.0000515)
1000	0.000747(0.0000622)	0.000945	0.00116(0.0000376)
2000	0.000543(0.0000553)	0.000610	0.000683(0.0000121)

(d) Normal location mixture, $\|\mu\| = 1$

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00434(0.00158)	0.00304	0.00514(0.000322)
500	0.000996(0.0000622)	0.00117	0.00146(0.000442)
1000	0.000640(0.0000502)	0.000760	0.000880(0.000176)
2000	0.000445(0.0000455)	0.000492	0.000583(0.0000192)

(e) Normal location mixture, $\|\mu\| = 2$

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00467(0.000139)	0.00326	0.00484(0.000244)
500	0.00173(0.0000522)	0.00126	0.00150(0.000363)
1000	0.00122(0.0000456)	0.000819	0.000925(0.0000131)
2000	0.00105(0.0000340)	0.000530	0.000577(0.0000651)

(f) Normal location mixture, $\|\mu\| = 3$

Table 3. Mean integrated squared error estimates (with standard errors in brackets where applicable; $d = 3$)

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00426(0.000131)	0.00240	0.00505(0.000279)
500	0.000835(0.0000302)	0.00106	0.00143(0.0000338)
1000	0.000442(0.0000236)	0.000737	0.000888(0.0000139)
2000	0.000304(0.0000238)	0.000508	0.000579(0.00000985)

(a) Independent Normal

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00467(0.000147)	0.00254	0.00550(0.000361)
500	0.000812(0.0000301)	0.00112	0.00152(0.0000367)
1000	0.000431(0.0000249)	0.000778	0.000922(0.0000145)
2000	0.000304(0.0000233)	0.000537	0.000603(0.00000676)

(b) Dependent Normal

n	LogConcDEAD	Kernel (opt AMISE)	Kernel (LSCV)
100	0.00365(0.000142)	0.00344	0.0741(0.00400)
500	0.000779(0.0000243)	0.00136	0.00192(0.0000518)
1000	0.000538(0.000104)	0.000922	0.00123(0.0000262)
2000	0.000292(0.0000414)	0.000622	0.000849(0.0000228)

(c) $\Gamma(2, 1)$ (independent components)

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00395(0.000124)	0.00214	0.00446(0.000242)
500	0.000743(0.0000272)	0.000946	0.00124(0.0000298)
1000	0.000446(0.0000218)	0.000656	0.000822(0.0000179)
2000	0.000265(0.0000202)	0.000452	0.000508(0.00000537)

(d) Normal location mixture, $\|\mu\| = 1$

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00319(0.000100)	0.00168	0.00371(0.000203)
500	0.000596(0.0000231)	0.000748	0.00103(0.0000340)
1000	0.000329(0.0000173)	0.000520	0.000656(0.0000160)
2000	0.000220(0.0000171)	0.000358	0.000410(0.00000519)

(e) Normal location mixture, $\|\mu\| = 2$

n	LogConcDEAD	Kernel (opt MISE)	Kernel (LSCV)
100	0.00328(0.0000930)	0.00166	0.00296(0.000120)
500	0.000803(0.0000184)	0.000751	0.000998(0.000254)
1000	0.000552(0.0000169)	0.000525	0.000613(0.0000892)
2000	0.000401(0.0000133)	0.000364	0.000404(0.00000488)

(f) Normal location mixture, $\|\mu\| = 3$

likelihood estimator will converge to the density f^* that minimises the Kullback–Leibler divergence $d(f_0 \| f) = \int f_0(x) \log \frac{f_0(x)}{f(x)} dx$ over $f \in \mathcal{F}_0$. Such a result would be interesting for robustness purposes, because it could be interpreted as saying that provided the underlying density does not violate the log-concavity assumption too seriously, the log-concave maximum likelihood estimator is still sensible.

6. Clustering example

In a recent paper, Chang and Walther (2008) introduced an algorithm which combines the univariate log-concave maximum likelihood estimator with the EM algorithm (Dempster *et al.*, 1977), to fit a finite mixture density of the form

$$f(x) = \sum_{j=1}^p \pi_j f_j(x), \quad (6.1)$$

where the mixture proportions π_1, \dots, π_p are positive and sum to one, and the component densities f_1, \dots, f_p are univariate and log-concave. The method is an extension of the standard Gaussian EM algorithm, e.g. Fraley and Raftery (2002), which assumes that each component density is normal. Once estimates $\hat{\pi}_1, \dots, \hat{\pi}_p, \hat{f}_1, \dots, \hat{f}_p$ have been obtained, clustering can be carried out by assigning to the j th cluster those observations X_i for which $j = \operatorname{argmax}_r \hat{\pi}_r \hat{f}_r(X_i)$. Chang and Walther (2008) show empirically that in cases where the true component densities are log-concave but not normal, their algorithm tends to make considerably fewer misclassifications and have smaller mean absolute error in the mixture proportion estimates than the Gaussian EM algorithm, with very similar performance in cases where the true component densities are normal.

Owing to the previous lack of an algorithm for computing the maximum likelihood estimator of a multidimensional log-concave density, Chang and Walther (2008) discuss an extension of the model in (6.1) to a multivariate context where the univariate marginal densities of each component in the mixture are assumed to be log-concave, and the dependence structure within each component density is modelled with a normal copula. Now that we are able to compute the maximum likelihood estimator of a multidimensional log-concave density, we can carry this method through to its natural conclusion. That is, in the finite mixture model (6.1) for a multidimensional log-concave density f , we simply assume that each of the component densities f_1, \dots, f_p is log-concave. An interesting problem that we do not address here that of finding appropriate conditions under which this model is identifiable – see Titterington *et al.* (1985, Section 3.1) for a nice discussion.

6.1. EM algorithm

An introduction to the EM algorithm can be found in McLachlan and Krishnan (1997). Briefly, given current estimates of the mixture proportions and component densities $\hat{\pi}_1^{(\ell)}, \dots, \hat{\pi}_p^{(\ell)}, \hat{f}_1^{(\ell)}, \dots, \hat{f}_p^{(\ell)}$ at the ℓ th iteration of the algorithm, we update the estimates of the mixture proportions by setting $\hat{\pi}_j^{(\ell+1)} = n^{-1} \sum_{i=1}^n \hat{\theta}_{i,j}^{(\ell)}$ for $j = 1, \dots, p$, where

$$\hat{\theta}_{i,j}^{(\ell)} = \frac{\hat{\pi}_j^{(\ell)} \hat{f}_j^{(\ell)}(X_i)}{\sum_{r=1}^p \hat{\pi}_r^{(\ell)} \hat{f}_r^{(\ell)}(X_i)}$$

is the current estimate of the posterior probability that the i th observation belongs to the j th component. We then update the estimates of the component densities in turn using the algorithm described in Section 4, choosing $\hat{f}_j^{(\ell+1)}$ to be the log-concave density f_j that maximises

$$\sum_{i=1}^n \hat{\theta}_{i,j}^{(\ell)} \log f_j(X_i).$$

The incorporation of the weights $\hat{\theta}_{1,j}^{(\ell)}, \dots, \hat{\theta}_{n,j}^{(\ell)}$ in the maximisation process presents no additional complication, as is easily seen by inspecting the proof of Theorem 2. As usual with methods based on the EM algorithm, although the likelihood increases at each iteration, there is no guarantee that the sequence converges to a global maximum. In fact, it can happen that the algorithm produces a sequence that approaches a degenerate solution, corresponding to a component concentrated on a single observation, so that the likelihood becomes arbitrarily high. The same issue can arise when fitting mixtures of Gaussian densities, and in this context Fraley and Raftery (2002) suggest that a Bayesian approach can alleviate the problem in these instances by effectively smoothing the likelihood. In general, it is standard practice to restart the algorithm from different initial values, taking the solution with the highest likelihood.

6.2. Breast cancer example

We illustrate the log-concave EM algorithm on the Wisconsin breast cancer data set of Street *et al.* (1993), available on the UCI Machine Learning Repository website (Asuncion and Newman, 2007):

<http://archive.ics.uci.edu/ml/datasets/Breast+Cancer+Wisconsin+%28Diagnostic%29>.

The data set was created by taking measurements from a digitised image of a fine needle aspirate of a breast mass, for each of 569 individuals, with 357 benign and 212 malignant instances. We study the problem of trying to diagnose (cluster) the individuals based on the standard errors of two of the measurements, namely the radius of the cell nucleus (mean of distances from center to points on the perimeter, X) and its texture (standard deviation of grey-scale values, Y). The data are presented in Figure 4(a). In fact, the full data set consists of 30 measurements for each patient, representing the mean, standard error and ‘worst’ (mean of the three largest values) of 10 different features computed for each cell nucleus in the image. Since one would reasonably expect the means of each feature to be approximately normally distributed, and hence the Gaussian EM algorithm to be appropriate, we took the standard errors of the first two measurements to illustrate the log-concave EM algorithm methodology.

It is important also to note that although for this particular data set we do know whether a particular instance is benign or malignant, we did not use this information in fitting our mixture model. Instead this information was only used afterwards to assess the performance of the method, as reported below. Thus we are studying a clustering (or unsupervised learning) problem, by taking a classification (or supervised learning) data set and ‘covering up the labels’ until it comes to performance assessment.

The skewness in the data suggests that the mixture of Gaussians model may be inadequate, and in Figure 4(b) we show the contour plot and misclassified instances from this model. The corresponding

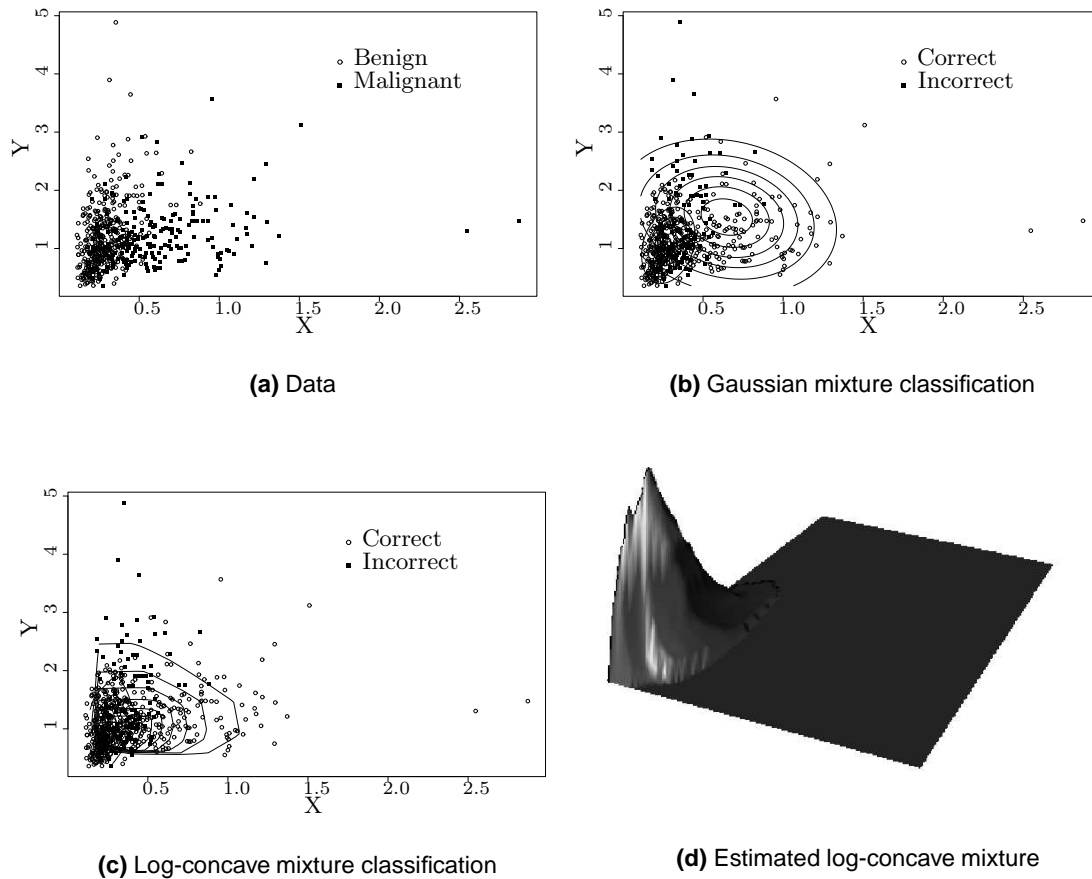


Fig. 4. Panel (a) plots the Wisconsin breast cancer data, with benign cases as solid squares and malignant ones as open circles. Panel (b) gives a contour plot together with the misclassified instances from the Gaussian EM algorithm, while the corresponding plot obtained from the log-concave EM algorithm is given in Panel (c). Panel (d) plots the fitted mixture distribution from the log-concave EM algorithm.

plot obtained from the log-concave EM algorithm is given in Figure 4(c), while Figure 4(d) plots the fitted mixture distribution from the log-concave EM algorithm. For this example, the number of misclassified instances is reduced from 144 with the Gaussian EM algorithm to 121 with the log-concave EM algorithm.

In some examples, it will be necessary to estimate p , the number of mixture components. In the general context of model-based clustering, Fraley and Raftery (2002) cite several possible approaches for this purpose, including methods based on resampling (McLachlan and Basford, 1988) and an information criterion (Bozdogan, 1994). Further research will be needed to ascertain which of these methods is most appropriate in the context of log-concave component densities.

7. Plug-in estimation of functionals, sampling and the bootstrap

Suppose X has density f . Often, we are less interested in estimating a density directly than in estimating some functional $\theta(f)$. Examples of functionals of interest (some of which were given in Section 1), include:

- (a) $\mathbb{P}(\|X\| \geq 1) = \int f(x) \mathbb{1}_{\{\|x\| \geq 1\}} dx$
- (b) Moments, such as $\mathbb{E}(X) = \int xf(x) dx$, or $\mathbb{E}(\|X\|^2) = \int \|x\|^2 f(x) dx$
- (c) The differential entropy of X (or f), defined by $H(f) = - \int f(x) \log f(x) dx$
- (d) The $100(1 - \alpha)\%$ highest density region, defined by $R_\alpha = \{x \in \mathbb{R}^d : f(x) \geq f_\alpha\}$, where f_α is the largest constant such that $\mathbb{P}(X \in R_\alpha) \geq 1 - \alpha$. Hyndman (1996) argues that this is an informative summary of a density; note that subject to a minor restriction on f , we have $\int f(x) \mathbb{1}_{\{f(x) \geq f_\alpha\}} dx = 1 - \alpha$.

Each of these may be estimated by the corresponding functional $\theta(\hat{f}_n)$ of the log-concave maximum likelihood estimator. In examples (a) and (b) above, $\theta(f)$ may also be written as a functional of the corresponding distribution function F , e.g. $\mathbb{P}(\|X\| \geq 1) = \int \mathbb{1}_{\{\|x\| \geq 1\}} dF(x)$. In such cases, it is more natural to use the plug-in estimator based on the empirical distribution function, \hat{F}_n , of the sample X_1, \dots, X_n , and indeed in our simulations we found that the log-concave plug-in estimator did not offer an improvement on this method. In the other examples, however, an empirical distribution function plug-in estimator is not available, and the log-concave plug-in estimator is a potentially attractive procedure.

7.1. Monte Carlo estimation of functionals and sampling from the density estimate

For some functionals we can compute $\hat{\theta} = \theta(\hat{f}_n)$ analytically. If this is not possible, but we can write $\theta(f) = \int f(x)g(x) dx$, we may approximate $\hat{\theta}$ by

$$\hat{\theta}_B = \frac{1}{B} \sum_{b=1}^B g(X_b^*),$$

for some (large) B , where X_1^*, \dots, X_B^* are independent samples from \hat{f}_n . Conditional on X_1, \dots, X_n , the strong law of large numbers gives that $\hat{\theta}_B \xrightarrow{a.s.} \hat{\theta}$ as $B \rightarrow \infty$. In practice, even when analytic calculation of $\hat{\theta}$ was possible, this method was found to be fast and accurate.

In order to use this Monte Carlo procedure, we must be able to sample from \hat{f}_n . Fortunately, this can be done efficiently using the following rejection sampling procedure. As in Section 4, for $j \in J$ let A_j be the $d \times d$ matrix whose l th column is $X_{j_{l+1}} - X_{j_l}$ for $l = 1, \dots, d$, and let $\alpha_j = X_{j_1}$, so that $w \mapsto A_j w + \alpha_j$ maps the unit simplex T_d to $C_{n,j}$. Recall that $\log \hat{f}_n(X_i) = y_i^*$, and let $z_j = (z_{j,1}, \dots, z_{j,d})$, where $z_{j,l} = y_{j_{l+1}}^* - y_{j_l}^*$ for $l = 1, \dots, d$. Write

$$q_j = \int_{C_{n,j}} \hat{f}_n(x) dx.$$

We may then draw an observation X^* from \hat{f}_n as follows:

- (i) Select $j^* \in J$, selecting $j^* = j$ with probability q_j
- (ii) Select $w \sim \text{Unif}(T_d)$ and $u \sim \text{Unif}([0, 1])$ independently. If

$$u < \frac{\exp(\langle w, z_{j^*} \rangle)}{\max_{v \in T_d} \exp(\langle v, z_{j^*} \rangle)},$$

accept the point and set $X^* = A_{j^*} w + \alpha_{j^*}$. Otherwise, repeat (ii).

7.2. Simulation study

In this section we illustrate some simple applications of this technique to functionals (c) and (d) above, using the Monte Carlo procedure and sampling scheme described in Section 7.1. Estimates are based on random samples from a $N_2(0, I)$ distribution, and we compare the performance of the `LogConcDEAD` estimate with that of a kernel-based plug-in estimate, where the bandwidth matrix was chosen using our knowledge of the underlying density to minimise the MISE.

Table 4(a) gives mean squared errors (with Monte Carlo standard errors) of the plug-in estimates of the differential entropy. In Table 4(b) we study the plug-in estimators \hat{R}_α of the highest density region R_α , and measure the quality of the estimation procedures through $\mathbb{E}\{\mu_f(\hat{R}_\alpha \Delta R_\alpha)\}$, where $\mu_f(A) = \int_A f(x) dx$ and Δ denotes set difference. Highest density regions can be computed once we have approximated the sample versions of f_α using the density quantile algorithm described in Hyndman (1996, Section 3.2).

For the differential entropy estimators, we find a similar pattern to that observed in Section 5: the log-concave plug-in estimator provides an improvement on the kernel-based estimator for the moderate and large sample sizes in our simulations. For the case of highest density regions, the relative performance of the log-concave estimator is better for the estimation of smaller density regions. In Figure 5, we illustrate the estimation of three highest density regions based on 500 points from a $N_2(0, I)$ distribution. For comparison, a kernel-based plug-in estimate (where the regions are not guaranteed to be convex) is also given.

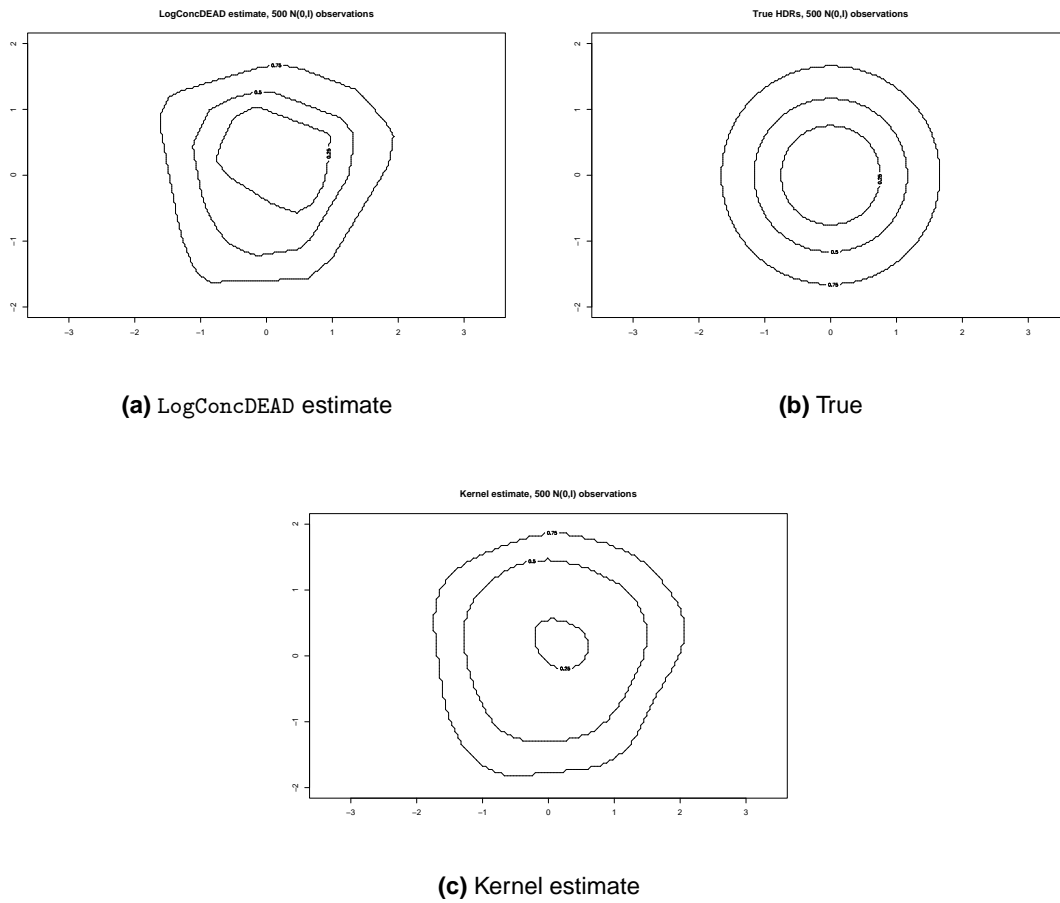


Fig. 5. Estimates of the 25%, 50% and 75% highest density region from 500 observations from the $N_2(0, I)$ distribution.

Table 4. (a) gives mean squared errors for estimating the differential entropy of the $N_2(0, I)$ distribution; (b) gives $\mathbb{E}\{\mu_f(\hat{R}_\alpha \triangle R_\alpha)\}$ when estimating highest density regions. The numbers in brackets are Monte Carlo standard errors.

n	LogConcDEAD	Kernel
100	0.0761(0.00629)	0.0457(0.00304)
500	0.00819(0.000653)	0.0137(0.000839)
1000	0.00378(0.000391)	0.00716(0.000581)
2000	0.00177(0.000232)	0.00427(0.000345)

(a) Differential entropy

n	LogConcDEAD	Kernel
100	0.0872(0.0024)/0.110(0.0033)/0.121(0.0047)	0.0753(0.0017)/0.0995(0.0028)/0.0959(0.0038)
500	0.0419(0.0010)/0.0587(0.0014)/0.0680(0.0022)	0.0467(0.0011)/0.0609(0.0013)/0.0637(0.0019)
1000	0.0311(0.00075)/0.0447(0.0011)/0.0536(0.0016)	0.0376(0.00095)/0.0476(0.0012)/0.0477(0.0015)
2000	0.0241(0.00054)/0.0363(0.00080)/0.0448(0.0013)	0.0322(0.00081)/0.0371(0.00098)/0.0399(0.0013)

(b) 25%/50%/75% highest density regions

In real data examples, we are unable to assess uncertainty in our functional estimates by taking repeated samples from the true underlying model. Nevertheless, the fact that we can sample from the log-concave maximum likelihood estimator does mean that we can apply standard bootstrap methodology to compute standard errors or confidence intervals, for example. Finally, we remark that the plug-in estimation procedure, sampling algorithm and bootstrap methodology extend in an obvious way to the case of a finite mixture of log-concave densities.

8. Concluding discussion

We have developed methodology that gives a fully automatic nonparametric density estimate under the condition that the density is log-concave, and shown how it may be extended to fit finite mixtures of log-concave densities. We have indicated a wide range of possible applications, including classification, clustering and functional estimation problems. The area of shape-constrained estimation is currently undergoing rapid growth, as evidenced by the many recent publications cited in the penultimate paragraph of Section 1, as well as recent workshops in Oberwolfach (November 2006), Eindhoven (October 2007) and Bristol (November 2007). We hope that this paper will stimulate further interest and research in the field.

As well as the continued development and refinement of the computational algorithms and graphical displays of estimates, and studies of theoretical performance, there remain many challenges and interesting directions for future research. These include:

- (i) Studying other shape constraints. These have received some attention for univariate data, dating back to Grenander (1956), but much less in the multivariate setting.
- (ii) Developing both formal and informal diagnostic tools for assessing the validity of shape con-

straints.

- (iii) Assessing the uncertainty in shape-constrained nonparametric density estimates, through confidence intervals/bands.
- (iv) Developing analogous methodology for discrete data from shape-constrained distributions.
- (v) Examining nonparametric shape constraints in regression problems.
- (vi) Studying methods for choosing the number of clusters in nonparametric, shape-constrained mixture models.

A. Glossary of terms and results from convex analysis and computational geometry

All of the definitions and results below can be found in Rockafellar (1997) and Lee (1997). The *epigraph* of a function $f : \mathbb{R}^k \rightarrow [-\infty, \infty)$ is the set

$$\text{epi}(f) = \{(x, \mu) : x \in \mathbb{R}^k, \mu \in \mathbb{R}, \mu \leq f(x)\}.$$

We say f is *concave* if its epigraph is non-empty and convex as a subset of \mathbb{R}^{k+1} ; note that this agrees with the terminology of Barndorff-Nielsen (1978), but is what Rockafellar (1997) calls a *proper concave* function. If C is a convex subset of \mathbb{R}^k then provided $f : C \rightarrow [-\infty, \infty)$ is not identically $-\infty$, it is *concave* if and only if

$$f(tx + (1-t)y) \geq tf(x) + (1-t)f(y)$$

for $x, y \in C$ and $t \in (0, 1)$. A non-negative function f is *log-concave* if $\log f$ is concave, with the convention that $\log 0 = -\infty$. The *support* of a log-concave function f is $\{x \in \mathbb{R}^k : \log f(x) > -\infty\}$, a convex subset of \mathbb{R}^k .

A subset M of \mathbb{R}^k is *affine* if $tx + (1-t)y \in M$ for all $x, y \in M$ and $t \in \mathbb{R}$. The *affine hull* of M , denoted $\text{aff}(M)$, is the smallest affine set containing M . Every non-empty affine set M in \mathbb{R}^k is *parallel* to a unique subspace of \mathbb{R}^k , meaning that there is a unique subspace L of \mathbb{R}^k such that $M = L + a$, for some $a \in \mathbb{R}^k$. The *dimension* of M is the dimension of this subspace, and more generally the dimension of a non-empty convex set is the dimension of its affine hull. A finite set of points $M = \{x_0, x_1, \dots, x_d\}$ is *affinely independent* if $\text{aff}(M)$ is d -dimensional. The *relative interior* of a convex set C is the interior which results when we regard C as a subset of its affine hull. The *relative boundary* of C is the set difference between its closure and its relative interior. If M is an affine set in \mathbb{R}^k , then an *affine transformation* (or *affine function*) is a function $T : M \rightarrow \mathbb{R}^k$ such that $T(tx + (1-t)y) = tT(x) + (1-t)T(y)$ for all $x, y \in M$ and $t \in \mathbb{R}$.

The *closure* of a concave function g on \mathbb{R}^d , denoted $\text{cl}(g)$, is the function whose epigraph is the closure in \mathbb{R}^{d+1} of $\text{epi}(g)$. It is the least upper semi-continuous, concave function satisfying $\text{cl}(g) \geq g$. The function g is *closed* if $\text{cl}(g) = g$. An arbitrary function h on \mathbb{R}^d is *continuous relative* to a subset S of \mathbb{R}^d if its restriction to S is a continuous function. A non-zero vector $z \in \mathbb{R}^d$ is a *direction of increase* of h on \mathbb{R}^d if $t \mapsto h(x + tz)$ is non-decreasing for every $x \in \mathbb{R}^d$.

The convex hull of finitely many points is called a *polytope*. The convex hull of $d + 1$ affinely independent points is called a *d -dimensional simplex* (pl. *simplices*). If C is a convex set in \mathbb{R}^d , then a *supporting half-space* to C is a closed half-space which contains C and has a point of C in its boundary. A *supporting hyperplane* H to C is a hyperplane which is the boundary of a supporting

half-space to C . Thus $H = \{x \in \mathbb{R}^d : \langle x, b \rangle = \beta\}$, for some $b \in \mathbb{R}^d$ and $\beta \in \mathbb{R}$ such that $\langle x, b \rangle \leq \beta$ for all $x \in C$ with equality for at least one $x \in C$.

If V is a finite set of points in \mathbb{R}^d such that $P = \text{conv}(V)$ is a d -dimensional polytope in \mathbb{R}^d , then a *face* of P is a set of the form $P \cap H$, where H is a supporting hyperplane to P . The *vertex set* of P , denoted $\text{vert}(P)$, is the set of 0-dimensional faces (*vertices*) of P . A *subdivision* of P is a finite set of d -dimensional polytopes $\{S_1, \dots, S_t\}$ such that P is the union of S_1, \dots, S_t and the intersection of any two distinct polytopes in the subdivision is a face of both of them. If $S = \{S_1, \dots, S_t\}$ and $\tilde{S} = \{\tilde{S}_1, \dots, \tilde{S}_{t'}\}$ are two subdivisions of P , then \tilde{S} is a *refinement* of S if each S_l is contained in some $\tilde{S}_{l'}$. The *trivial subdivision* of P is $\{P\}$. A *triangulation* of P is a subdivision of P in which each polytope is a simplex.

If P is a d -dimensional polytope in \mathbb{R}^d , F is a $(d-1)$ -dimensional face of P and $v \in \mathbb{R}^d$, then there is a unique supporting hyperplane H to P containing F . The polytope P is contained in exactly one of the closed half-spaces determined by H , and if v is in the opposite open half-space, then F is *visible* from v . If V is a finite set in \mathbb{R}^d such that $P = \text{conv}(V)$, if $v \in V$ and $S = \{S_1, \dots, S_t\}$ is a subdivision of P , then the result of *pushing* v is the subdivision \tilde{S} of P obtained by modifying each $S_l \in S$ as follows:

- (i) If $v \notin S_l$, then $S_l \in \tilde{S}$
- (ii) If $v \in S_l$ and $\text{conv}(\text{vert}(S_l) \setminus \{v\})$ is $(d-1)$ -dimensional, then $S_l \in \tilde{S}$
- (iii) If $v \in S_l$ and $S'_l = \text{conv}(\text{vert}(S_l) \setminus \{v\})$ is d -dimensional, then $S'_l \in \tilde{S}$. Also, if F is any $(d-1)$ -dimensional face of S'_l that is visible from v , then $\text{conv}(F \cup \{v\}) \in \tilde{S}$.

If σ is a convex function on \mathbb{R}^n , then $y' \in \mathbb{R}^n$ is a *subgradient* of σ at y if

$$\sigma(z) \geq \sigma(y) + \langle y', z - y \rangle$$

for all $z \in \mathbb{R}^n$. If σ is differentiable at y , then $\nabla\sigma(y)$ is the unique subgradient to σ at y ; otherwise the set of subgradients at y has more than one element. The *one-sided directional derivative* of σ at y with respect to $z \in \mathbb{R}^n$ is

$$\sigma'(y; z) = \lim_{t \searrow 0} \frac{\sigma(y + tz) - \sigma(y)}{t},$$

which always exists (allowing $-\infty$ and ∞ as limits) provided $\sigma(y)$ is finite.

B. Proofs

PROOF OF PROPOSITION 1

(a) If f is log-concave, then for $x \in \mathbb{R}^d$, we can write

$$f_{X|P_V(X)}(x|t) \propto f(x) \mathbb{1}_{\{P_V(x)=t\}},$$

a product of log-concave functions. Thus $f_{X|P_V(X)}(\cdot|t)$ is log-concave for each t .

(b) Let $x_1, x_2 \in \mathbb{R}^d$ be distinct and let $\lambda \in (0, 1)$. Let V be the $(d-1)$ -dimensional subspace of \mathbb{R}^d whose orthogonal complement is *parallel* to the *affine hull* of $\{x_1, x_2\}$ (i.e. the line through x_1 and

x_2). Writing $f_{P_V(X)}$ for the marginal density of $P_V(X)$ and t for the common value of $P_V(x_1)$ and $P_V(x_2)$, the density of X at $x \in \mathbb{R}^d$ is

$$f(x) = f_{X|P_V(X)}(x|t)f_{P_V(X)}(t).$$

Thus f is log-concave, as required. \square

COMPLETION OF THE PROOF OF THEOREM 2

We prove each of the steps (i)–(v) outlined in Section 3 in turn. First note that if $x_0 \in C_n$, then by Carathéodory's theorem (Theorem 17.1 of Rockafellar (1997)), there exist distinct indices i_1, \dots, i_r with $r \leq d + 1$, such that $x_0 = \sum_{l=1}^r \lambda_l X_{i_l}$ with each $\lambda_l > 0$ and $\sum_{l=1}^r \lambda_l = 1$. Thus, if $f(x_0) = 0$, then by Jensen's inequality,

$$-\infty = \log f(x_0) \geq \sum_{l=1}^r \lambda_l \log f(X_{i_l}),$$

so $f(X_i) = 0$ for some i . But then $\psi_n(f) = -\infty$. This proves (i).

Now suppose $f(x_0) > 0$ for some $x_0 \notin C_n$. Then $\{x : f(x) > 0\}$ is a convex set containing $C_n \cup \{x_0\}$, a set which has strictly larger d -dimensional Lebesgue measure than that of C_n . We therefore have $\psi_n(f) < \psi_n(f\mathbb{1}_{C_n})$, which proves (ii).

To prove (iii), we first show that $\log f$ is *closed*. Suppose that $\log f(X_i) = y_i$ for $i = 1, \dots, n$ but that $\log f \neq \bar{h}_y$. Then since $\log f(x) \geq \bar{h}_y(x)$ for all $x \in \mathbb{R}^d$, we may assume that there exists $x_0 \in C_n$ such that $\log f(x_0) > \bar{h}_y(x_0)$. If x_0 is in the *relative interior* of C_n , then since $\log f$ and \bar{h}_y are continuous at x_0 (by Theorem 10.1 of Rockafellar (1997)), we must have

$$\psi_n(f) < \psi_n(\exp(\bar{h}_y)).$$

The only remaining possibility is that x_0 is on the *relative boundary* of C_n . But \bar{h}_y is closed by Corollary 17.2.1 of Rockafellar (1997), so writing $\text{cl}(g)$ for the *closure* of a concave function g , we have $\bar{h}_y = \text{cl}(\bar{h}_y) = \text{cl}(\log f) \geq \log f$, where we have used Corollary 7.3.4 of Rockafellar (1997) to obtain the middle equality. It follows that $\log f$ is closed and $\log f = \bar{h}_y$, which proves (iii).

Note that $\log f$ has no *direction of increase*, because if $x \in C_n$, z is a non-zero vector and $t > 0$ is large enough that $x + tz \notin C_n$, then $-\infty = \log f(x + tz) < \log f(x)$. It follows by Theorem 27.2 of Rockafellar (1997) that the supremum of f is finite (and is attained). Using properties (i) and (ii) as well, we may write $\int f(x) dx = c$, say, where $c \in (0, \infty)$. Thus $f(x) = cf(x)$, for some $\bar{f} \in \mathcal{F}_0$. But then

$$\psi_n(\bar{f}) - \psi_n(f) = -1 - \log c + c \geq 0,$$

with equality only if $c = 1$. This proves (iv).

To prove (v), we may assume by (iv) that $\exp(\bar{h}_y)$ is a density. Let $\max_i \bar{h}_y(X_i) = M$ and $\min_i \bar{h}_y(X_i) = m$. We show that when M is large, in order for $\exp(\bar{h}_y)$ to be a density, m must be negative with $|m|$ so large that $\psi_n(\exp(\bar{h}_y)) \leq \psi_n(f)$. First observe that if $x \in C_n$ and $\bar{h}_y(X_i) = M$, then for M sufficiently large we must have $M - m > 1$, and then

$$\begin{aligned} \bar{h}_y\left(X_i + \frac{1}{M-m}(x - X_i)\right) &\geq \frac{1}{M-m}\bar{h}_y(x) + \frac{M-m-1}{M-m}\bar{h}_y(X_i) \\ &\geq \frac{m}{M-m} + \frac{(M-m-1)M}{M-m} = M-1. \end{aligned}$$

(The fact that $\bar{h}_y(x) \geq m$ follows by Jensen's inequality.) Hence, denoting Lebesgue measure on \mathbb{R}^d by μ , we have

$$\mu(\{x : \bar{h}_y(x) \geq M - 1\}) \geq \mu\left(\left\{X_i + \frac{1}{M - m}(C_n - X_i)\right\}\right) = \frac{\mu(C_n)}{(M - m)^d}.$$

Thus

$$\int_{\mathbb{R}^d} \exp\{\bar{h}_y(x)\} dx \geq e^{M-1} \frac{\mu(C_n)}{(M - m)^d}.$$

For $\exp(\bar{h}_y)$ to be a density, then, we require $m \leq -\frac{1}{2}e^{(M-1)/d}\mu(C_n)^{1/d}$ when M is large. But then

$$\psi_n(\exp(\bar{h}_y)) \leq \frac{(n-1)M}{n} - \frac{1}{2n}e^{(M-1)/d}\mu(C_n)^{1/d} \leq \psi_n(f)$$

when M is sufficiently large. This proves (v).

It is not hard to see that for any $M > 0$, the function $y \mapsto \psi_n(\exp(\bar{h}_y))$ is continuous on the compact set $[-M, M]^n$, and thus the proof of the existence of a maximum likelihood estimator is complete. To prove uniqueness, suppose that $f_1, f_2 \in \mathcal{F}$ and both f_1 and f_2 maximise $\psi_n(f)$. We may assume $f_1, f_2 \in \mathcal{F}_0$, $\log f_1, \log f_2 \in \mathcal{H}$ and f_1 and f_2 are supported on C_n . Then the normalised geometric mean

$$g(x) = \frac{\{f_1(x)f_2(x)\}^{1/2}}{\int_{C_n} \{f_1(y)f_2(y)\}^{1/2} dy},$$

is a log-concave density, with

$$\begin{aligned} \psi_n(g) &= \frac{1}{2n} \sum_{i=1}^n \log f_1(X_i) + \frac{1}{2n} \sum_{i=1}^n \log f_2(X_i) - \log \int_{C_n} \{f_1(y)f_2(y)\}^{1/2} dy - 1 \\ &= \psi_n(f_1) - \log \int_{C_n} \{f_1(y)f_2(y)\}^{1/2} dy. \end{aligned}$$

However, by Cauchy-Schwarz, $\int_{C_n} \{f_1(y)f_2(y)\}^{1/2} dy \leq 1$, so $\psi_n(g) \geq \psi_n(f_1)$. Equality is obtained if and only if $f_1 = f_2$ almost everywhere, but since f_1 and f_2 are *continuous relative* to C_n (Theorem 10.2 of Rockafellar (1997)), this implies that $f_1 = f_2$. An alternative way of proving the uniqueness of the maximum likelihood estimator may be based on the fact that $\psi_n(tf_1 + (1-t)f_2) > t\psi_n(f_1) + (1-t)\psi_n(f_2)$ for all $t \in (0, 1)$, provided f_1 and f_2 are distinct elements of \mathcal{F} . \square

PROOF OF THEOREM 3

For $t \in (0, 1)$ and $y^{(1)}, y^{(2)} \in \mathbb{R}^n$, the function $\bar{h}_{ty^{(1)}+(1-t)y^{(2)}}$ is the least concave function satisfying $\bar{h}_{ty^{(1)}+(1-t)y^{(2)}}(X_i) \geq ty_i^{(1)} + (1-t)y_i^{(2)}$ for $i = 1, \dots, n$, so $\bar{h}_{ty^{(1)}+(1-t)y^{(2)}} \leq t\bar{h}_{y^{(1)}} + (1-t)\bar{h}_{y^{(2)}}$. The convexity of σ follows from this and the convexity of the exponential function. It is clear that $\sigma \geq \tau$, since $\bar{h}_y(X_i) \geq y_i$ for $i = 1, \dots, n$.

From Theorem 2, we can find $y^* \in \mathbb{R}^n$ such that $\log \hat{f}_n = \bar{h}_{y^*}$ with $\bar{h}_{y^*}(X_i) = y_i^*$ for $i = 1, \dots, n$, and this y^* minimises τ . For any other $y \in \mathbb{R}^n$ which minimises τ , by the uniqueness part of Theorem 2 we must have $\bar{h}_y = \bar{h}_{y^*}$, so $\sigma(y) > \sigma(y^*) = \tau(y^*)$. \square

B.1. Non-differentiability of σ and computation of subgradients

In this section, we find explicitly the set of points at which the function σ defined in (4.1) is differentiable, and compute a subgradient of σ at each point. For $i = 1, \dots, n$, define

$$J_i = \{j = (j_1, \dots, j_{d+1}) \in J : i = j_l \text{ for some } l = 1, \dots, d+1\}.$$

The set J_i is the index set of those simplices $C_{n,j}$ that have X_i as a vertex. Let \mathcal{Y} denote the set of vectors $y = (y_1, \dots, y_n) \in \mathbb{R}^n$ with the property that for each $j = (j_1, \dots, j_{d+1}) \in J$, if $i \neq j_l$ for any l then

$$\{(X_i, y_i), (X_{j_1}, y_{j_1}), \dots, (X_{j_{d+1}}, y_{j_{d+1}})\}$$

is affinely independent in \mathbb{R}^{d+1} . This is the set of points for which no tent pole is touching but not critically supporting the tent. Notice that the complement of \mathcal{Y} has zero Lebesgue measure in \mathbb{R}^n . For $y \in \mathbb{R}^n$ and $i = 1, \dots, n$, and in the notation of Section 4, let

$$\partial_i(y) = -\frac{1}{n} + \sum_{j \in J_i} |\det A_j| \int_{T_d} e^{\langle w, z_j \rangle + y_{j_1}} \left\{ \left(1 - \sum_{l=1}^d w_l\right) \mathbb{1}_{\{j_1=i\}} + \sum_{l=1}^d w_l \mathbb{1}_{\{j_{l+1}=i\}} \right\} dw.$$

PROPOSITION 4. *Assume (A1). (a) For $y \in \mathcal{Y}$, the function σ is differentiable at y and for $i = 1, \dots, n$ satisfies*

$$\frac{\partial \sigma}{\partial y_i}(y) = \partial_i(y).$$

(b) For $y \in \mathcal{Y}^c$, the function σ is not differentiable at y , but the vector $(\partial_1(y), \dots, \partial_n(y))$ is a subgradient of σ at y .

PROOF. By Theorem 25.2 of Rockafellar (1997), it suffices to show that for $y \in \mathcal{Y}$, all of the partial derivatives exist and are given by the expression in the statement of the proposition. For $i = 1, \dots, n$ and $t \in \mathbb{R}$, let $y^{(t)} = y + te_i^n$, where e_i^n denotes the i th unit coordinate vector in \mathbb{R}^n . For sufficiently small values of $|t|$, we may write

$$\bar{h}_{y^{(t)}}(x) = \begin{cases} \langle x, b_j^{(t)} \rangle - \beta_j^{(t)} & \text{if } x \in C_{n,j} \text{ for some } j \in J \\ -\infty & \text{if } x \notin C_n, \end{cases}$$

for certain values of $b_1^{(t)}, \dots, b_m^{(t)} \in \mathbb{R}^d$ and $\beta_1^{(t)}, \dots, \beta_m^{(t)} \in \mathbb{R}$. If $j \notin J_i$, then $b_j^{(t)} = b_j$ and $\beta_j^{(t)} = \beta_j$ for sufficiently small $|t|$. On the other hand, if $j \in J_i$, then there are two cases to consider:

- (i) If $j_1 = i$, then for sufficiently small t , we have $z_j^{(t)} = z_j - t1_d$, where 1_d denotes a d -vector of ones, so that $b_j^{(t)} = b_j - t(A_j^T)^{-1}1_d$ and $\beta_j^{(t)} = \beta_j - t(1 + \langle A_j^{-1}\alpha_j, 1_d \rangle)$
- (ii) If $j_{l+1} = i$ for some $l \in \{1, \dots, d\}$, then for sufficiently small t , we have $z_j^{(t)} = z_j + te_l^d$, so that $b_j^{(t)} = b_j + t(A_j^T)^{-1}e_l^d$ and $\beta_j^{(t)} = \beta_j + t\langle A_j^{-1}\alpha_j, e_l^d \rangle$.

It follows that

$$\begin{aligned}
\frac{\partial \sigma}{\partial y_i}(y) &= -\frac{1}{n} + \lim_{t \rightarrow 0} \frac{1}{t} \sum_{j \in J_i} \int_{C_{n,j}} \exp\{\langle x, b_j^{(t)} \rangle - \beta_j^{(t)}\} - \exp\{\langle x, b_j \rangle - \beta_j\} dx \\
&= -\frac{1}{n} + \lim_{t \rightarrow 0} \frac{1}{t} \sum_{j \in J_i} \left[\int_{C_{n,j}} e^{\langle x, b_j \rangle - \beta_j} \{e^{t(1 - \langle A_j^{-1}(x - \alpha_j), 1_d \rangle)} - 1\} dx \mathbb{1}_{\{j_1=i\}} \right. \\
&\quad \left. + \sum_{l=1}^d \int_{C_{n,j}} e^{\langle x, b_j \rangle - \beta_j} \{e^{t \langle A_j^{-1}(x - \alpha_j), e_l^d \rangle} - 1\} dx \mathbb{1}_{\{j_{l+1}=i\}} \right] \\
&= \partial_i(y),
\end{aligned}$$

where to obtain the final line we have made the substitution $x = A_j w + \alpha_j$, after taking the limit as $t \rightarrow 0$.

(b) If $y \in \mathcal{Y}^c$, then it can be shown that there exists a unit coordinate vector e_i^n in \mathbb{R}^n such that the *one-sided directional derivative* at y with respect to e_i^n , denoted $\sigma'(y; e_i^n)$, satisfies $\sigma'(y; e_i^n) > -\sigma'(y; -e_i^n)$. Thus σ is not differentiable at y . To show that $\partial(y) = (\partial_1(y), \dots, \partial_n(y))$ is a subgradient of σ at y , it is enough by Theorem 25.6 of Rockafellar (1997) to find, for each $\epsilon > 0$, a point $\tilde{y} \in \mathbb{R}^n$ such that $\|\tilde{y} - y\| < \epsilon$ and such that σ is differentiable at \tilde{y} with $\|\nabla \sigma(\tilde{y}) - \partial(y)\| < \epsilon$. This can be done by sequentially making small adjustments to the components of y in the same order as that in which the vertices were *pushed* in constructing the triangulation. \square

A subgradient of σ at any $y \in \mathbb{R}^n$ may be computed using Proposition 4, (B.1) and (4.2) once we have a formula for

$$\tilde{I}_{d,u}(z) = \int_{T_d} w_u \exp\left(\sum_{r=1}^d z_r w_r\right) dw,$$

when z_1, \dots, z_d are non-zero and distinct. In Cule *et al.* (2008b), it is shown that the required formula is

$$\begin{aligned}
\tilde{I}_{d,u}(z) &= \sum_{\substack{1 \leq r \leq d \\ r \neq u}} \frac{e^{z_r}}{z_r(z_r - z_u)} \prod_{\substack{1 \leq s \leq d \\ s \neq r}} \frac{1}{(z_r - z_s)} - \sum_{\substack{1 \leq r \leq d \\ r \neq u}} \frac{e^{z_u}}{z_r(z_r - z_u)} \prod_{\substack{1 \leq s \leq d \\ s \neq r}} \frac{1}{(z_r - z_s)} \\
&\quad + \frac{(-1)^d (e^{z_u} - 1)}{z_u \prod_{s=1}^d z_s} + \frac{e^{z_u}}{z_u} \prod_{\substack{1 \leq s \leq d \\ s \neq u}} \frac{1}{(z_u - z_s)}. \quad (\text{B.1})
\end{aligned}$$

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Further technical arguments

We first establish (4.2). We have

$$\begin{aligned} \int_{C_n} \exp\{\bar{h}_y(x)\} dx &= \sum_{j \in J} |\det A_j| e^{-\beta_j} \int_{T_d} \exp\{\langle A_j w + \alpha_j, b_j \rangle\} dw \\ &= \sum_{j \in J} |\det A_j| e^{y_{j1}} \int_{T_d} \exp\left(\sum_{r=1}^d z_{j,r} w_r\right) dw. \end{aligned}$$

For ease of notation, we drop the j subscript, and consider

$$I_d(z) = \int_{T_d} \exp\left(\sum_{r=1}^d z_r w_r\right) dw.$$

Observe that $I_1(z_1) = z_1^{-1}(e^{z_1} - 1)$, in agreement with (4.2). Assume the result for $d - 1$ as an inductive hypothesis. Then

$$\begin{aligned} I_d(z) &= \frac{1}{z_d} e^{z_d} \int_{T_{d-1}} \exp\left(\sum_{r=1}^{d-1} (z_r - z_d) w_r\right) dw - \frac{1}{z_d} \int_{T_{d-1}} \exp\left(\sum_{r=1}^{d-1} z_r w_r\right) dw \\ &= \frac{1}{z_d} e^{z_d} I_{d-1}(z_1 - z_d, \dots, z_{d-1} - z_d) - \frac{1}{z_d} I_{d-1}(z_1, \dots, z_d) \\ &= \frac{1}{z_d} \sum_{r=1}^{d-1} \frac{e^{z_r} - e^{z_d}}{(z_r - z_d)} \prod_{\substack{1 \leq s \leq d-1 \\ s \neq r}} \frac{1}{(z_r - z_s)} - \frac{1}{z_d} \sum_{r=1}^{d-1} \frac{e^{z_r} - 1}{z_r} \prod_{\substack{1 \leq s \leq d-1 \\ s \neq r}} \frac{1}{(z_r - z_s)} \\ &= \sum_{r=1}^{d-1} \frac{e^{z_r}}{z_r} \prod_{\substack{1 \leq s \leq d \\ s \neq r}} \frac{1}{(z_r - z_s)} - \frac{e^{z_d}}{z_d} \sum_{r=1}^{d-1} \prod_{\substack{1 \leq s \leq d \\ s \neq r}} \frac{1}{(z_r - z_s)} + \frac{1}{z_d} \sum_{r=1}^{d-1} \frac{1}{z_r} \prod_{\substack{1 \leq s \leq d-1 \\ s \neq r}} \frac{1}{(z_r - z_s)}. \quad (\text{B.2}) \end{aligned}$$

To deal with the middle term in the last line of (B.2) above, define a polynomial

$$P_d(t) = \sum_{r=1}^d \prod_{\substack{1 \leq s \leq d \\ s \neq r}} \left(\frac{t - z_s}{z_r - z_s} \right) - 1.$$

This polynomial $P_d(t)$ is of degree at most $d - 1$, but has roots z_1, \dots, z_d , so is identically zero. Examining the coefficient of t^{d-1} in this polynomial, we find

$$\sum_{r=1}^d \prod_{\substack{1 \leq s \leq d \\ s \neq r}} \frac{1}{(z_r - z_s)} = 0.$$

To deal with the final term in (B.2), observe that

$$\begin{aligned} \frac{1}{z_d} \sum_{r=1}^{d-1} \frac{1}{z_r} \prod_{\substack{1 \leq s \leq d-1 \\ s \neq r}} \frac{1}{(z_r - z_s)} &= \frac{1}{z_1 z_2 \dots z_d} \sum_{r=1}^{d-1} \prod_{\substack{1 \leq s \leq d-1 \\ s \neq r}} \frac{z_s}{(z_r - z_s)} = \frac{(-1)^d}{z_1 z_2 \dots z_d} \{P_{d-1}(0) + 1\} \\ &= \frac{(-1)^d}{z_1 z_2 \dots z_d} = - \sum_{r=1}^d \frac{1}{z_r} \prod_{\substack{1 \leq s \leq d \\ s \neq r}} \frac{1}{(z_r - z_s)}. \end{aligned}$$

Substituting these expressions into (B.2) yields (4.2). \square

Our final task is to establish (B.1). To this end, observe that for $u = 1$, we have the recurrence relation

$$\tilde{I}_{d,1}(z) = \sum_{r=0}^R \frac{e^{z_{d-r}} \tilde{I}_{d-R-1,1}(z_1 - z_{d-r}, \dots, z_{d-R-1} - z_{d-r})}{z_{d-r} \prod_{\substack{1 \leq s \leq R \\ s \neq r}} (z_{d-r} - z_{d-s})} + \frac{(-1)^{R+1}}{\prod_{r=0}^R z_{d-r}} \tilde{I}_{d-R-1,1}(z_1, \dots, z_{d-R-1}),$$

which holds for $R = 0, \dots, d-2$, and may be proved by induction on R . The formulae for other values of u may be deduced by symmetry. The formula for $\tilde{I}_{d,u}(z)$ in (B.1) is found by using this expression with $R = d-2$ together with the fact that for $z \neq 0$,

$$\int_0^1 w \exp(zw) dw = \frac{1}{z} \left(1 - \frac{1}{z}\right) e^z + \frac{1}{z^2}.$$

\square