

Fitting circles to scattered data: parameter estimates have no moments

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Abstract

We study a nonlinear regression problem of fitting a circle (or a circular arc) to scattered data. We prove that under any standard assumptions on the statistical distribution of errors that are commonly adopted in the literature, the estimates of the circle center and radius have infinite moments. We also discuss methodological implications of this fact.

Keywords: orthogonal regression, errors-in-variables, least squares fit, circle fitting, moments of estimates.

1 Introduction

Regression models in which all variables are subject to errors are known as error-in-variables (EIV) models. The EIV regression problem is quite different (and far more difficult) than the classical regression where the independent variable is assumed to be error-free. The EIV regression, even in the linear case, presents extremely challenging questions and leads to some counterintuitive results (some of them are mentioned below).

This work is devoted to a nonlinear EIV model where one fits a circle to scattered data. This is one of the basic tasks in pattern recognition and

computer vision. The need of fitting circles to planar images also arises in biology and medicine, nuclear physics, archeology, industry, and other areas of human practice.

The most popular method used to solve this problem is orthogonal least squares, i.e. the minimization of the sum of squares of the distances from the data points to the fitting contour. This method is often called *geometric fit* or *orthogonal distance regression* (ODR).

Fitting a circle to observed points $(x_1, y_1), \dots, (x_n, y_n)$ amounts to minimizing the objective function

$$(1) \quad \mathcal{F}(a, b, R) = \sum_{i=1}^n [\sqrt{(x_i - a)^2 + (y_i - b)^2} - R]^2,$$

where (a, b) denotes the center and R the radius of the circle. Then the parameters of the best fitting circle are defined by

$$(2) \quad (\hat{a}, \hat{b}, \hat{R}) = \operatorname{argmin} \mathcal{F}(a, b, R).$$

To explore the statistical properties of the estimates $\hat{a}, \hat{b}, \hat{R}$ one needs to make assumptions on the probability distribution of the data points. It is commonly assumed that each (x_i, y_i) is a noisy observation of some *true point* (x_i^*, y_i^*) , i.e.

$$(3) \quad x_i = x_i^* + \delta_i, \quad y_i = y_i^* + \varepsilon_i, \quad i = 1, \dots, n,$$

where $(\delta_1, \varepsilon_1), \dots, (\delta_n, \varepsilon_n)$ are n independent random vectors, usually with zero mean.

A standard assumption is that each $(\delta_i, \varepsilon_i)$ is a normal (Gaussian) vector with some covariance matrix \mathbf{C}_i . The simplest choice is $\mathbf{C}_i = \sigma^2 \mathbf{I}$, in which case all errors ε_i 's and δ_i 's are i.i.d. normal random variables with zero mean and a common variance σ^2 . In that case the geometric fit (2) coincides with the maximum likelihood estimate (MLE), see Chan 1965.

The true points (x_i^*, y_i^*) are supposed to lie on a ‘true circle’, i.e. satisfy

$$(4) \quad (x_i^* - a^*)^2 + (y_i^* - b^*)^2 = (R^*)^2, \quad i = 1, \dots, n,$$

where (a^*, b^*, R^*) denote the ‘true’ (unknown) parameters. Therefore

$$x_i^* = a^* + R^* \cos \varphi_i, \quad y_i^* = b^* + R^* \sin \varphi_i,$$

where $\varphi_1, \dots, \varphi_n$ specify the location of the true points on the true circle.

The angles $\varphi_1, \dots, \varphi_n$ can be regarded as fixed unknowns, then they have to be treated as additional parameters of the model (often called *incidental* or *latent* parameters). This setup is known as a *functional model*, see Chan 1965.

Alternatively, $\varphi_1, \dots, \varphi_n$ can be regarded as independent realizations of a random variable with a certain probability distribution on $[0, 2\pi]$; then one gets the so called *structural model*, see Anderson 1981 or Berman and Culpin 1986. Both models are widely used in the literature.

Many authors study the distribution of the estimates $\hat{a}, \hat{b}, \hat{R}$ under the above assumptions and try to evaluate their biases and covariance matrix. Our main result is

Theorem 1. *If the probability distribution of each vector $(\delta_i, \varepsilon_i)$ has a continuous strictly positive density, then $\hat{a}, \hat{b}, \hat{R}$ do not have moments, i.e.*

$$E(|\hat{a}|) = E(|\hat{b}|) = E(\hat{R}) = \infty.$$

Thus the estimates $\hat{a}, \hat{b}, \hat{R}$ have no mean values or variances.

Our assumptions include (but are not limited to) normally distributed errors. The distribution of $(\delta_i, \varepsilon_i)$ need not be the same for different i 's, it may depend on i , but the vectors $(\delta_i, \varepsilon_i)$ must be independent. The mean value of $(\delta_i, \varepsilon_i)$ need not be zero. The theorem is valid for every $n \geq 3$.

2 Historical remarks

Our result is not entirely surprising as a similar theorem has been proven for orthogonal least squares lines by Anderson 1976. Suppose one fits a line $y = \alpha + \beta x$ to data points $(x_1, y_1), \dots, (x_n, y_n)$ by minimizing the sum of squares of (orthogonal) distances, i.e. the estimates are defined by

$$(\hat{\alpha}, \hat{\beta}) = \operatorname{argmin} \frac{1}{1 + \beta^2} \sum_{i=1}^n (y_i - \alpha - \beta x_i)^2.$$

Again the observed points are random perturbations of some true points, in the sense of (3), which lie on an unknown true line, i.e. satisfy

$$(5) \quad y_i^* = \alpha^* + \beta^* x_i^*, \quad i = 1, \dots, n.$$

The true points are either fixed parameters (making it a functional model), or randomly sampled on the true line (structural model).

Theorem 2 (Anderson 1976). *If the errors δ_i 's and ε_i 's are i.i.d. normal random variables with zero mean and a common variance $\sigma^2 > 0$, then $\hat{\alpha}$ and $\hat{\beta}$ do not have moments, i.e. $E(|\hat{\alpha}|) = E(|\hat{\beta}|) = \infty$.*

Until Anderson's discovery, statisticians were used to employ Taylor expansion to derive some 'approximate' formulas for the moments of the estimates $\hat{\alpha}$ and $\hat{\beta}$ (including their means and variances). Anderson demonstrated that all those formulas should be regarded as *moments of some approximations*, rather than 'approximate moments'.

Anderson's result was rather sensational at the time, it was followed by heated discussions and a period of acute interest in the linear EIV regression. It also created methodological problems which we discuss in the next section.

Anderson proved his theorem by using an explicit formula for the density function of $\hat{\beta}$ (that formula was mentioned but not given in his paper; it appeared in a later paper by Anderson and Sawa 1982. Anderson also remarked that his result can be 'intuitively seen' from a well known formula for $\hat{\beta}$:

$$(6) \quad \hat{\beta} = \frac{s_{yy} - s_{xx} + \sqrt{(s_{yy} - s_{xx})^2 + 4s_{xy}^2}}{2s_{xy}},$$

where standard statistical notation are used: $s_{xx} = \sum_{i=1}^n (x_i - \bar{x})^2$, $s_{yy} = \sum_{i=1}^n (y_i - \bar{y})^2$, $s_{xy} = \sum_{i=1}^n (x_i - \bar{x})(y_i - \bar{y})$, and $\bar{x} = \frac{1}{n} \sum_{i=1}^n x_i$, $\bar{y} = \frac{1}{n} \sum_{i=1}^n y_i$. Anderson 1976 says that the (continuous) density of s_{xx} , s_{yy} , and s_{xy} for which the numerator in (6) is different from 0 and the denominator is equal to 0 is positive, hence the integral of the product of \hat{b} and its density diverges.

This 'intuitive' explanation can be easily converted into a rigorous proof, and then one readily extends Anderson's theorem to arbitrary distributions of errors as long as they have continuous strictly positive densities, like in our Theorem 1. (Alternatively, one can easily modify our constructions below to achieve this goal; this is all fairly straightforward, so we omit details.)

We note that Anderson's result was recently extended to some other estimates of the linear parameters α and β , see Chen and Kukush 2006 and an example in Zelniker and Clarkson 2006.

The problem of fitting circles (as well as other nonlinear curves) to data is technically much more difficult than that of fitting lines. In particular,

there are no explicit formulas for the estimates \hat{a} , \hat{b} or \hat{R} , analogous to (6), let alone explicit formulas for their probability densities. All the methods of computing the estimates \hat{a} , \hat{b} or \hat{R} are based on iterative numerical schemes.

All this makes the problem of fitting circles (ellipses, etc.) so much different from that of fitting lines that Anderson's result apparently passed unnoticed by the 'curve fitting community'. It is still commonly believed that the curve's parameters have finite moments; thus many researchers try to minimize their bias and variances or compute Cramer-Rao lower bounds on the covariance matrix, see e.g. Kanatani 1998 or Chernov and Lesort 2004.

Our theorem shows that the true moments are infinite in the case of fitting circles. We believe this result holds for ellipses and other types of curves, and we plan to investigate this issue.

We note that Kukush et al. 2004 recently modified the geometric fitting of ellipses to data in order to ensure the consistency of the parameter estimates (as the orthogonal regression estimates are inconsistent). They noted that their modified estimates had infinite moments, which at the time seemed to be a price to pay for consistency. It is now clear that the lack of moments is a rather general property of estimates under the EIV model.

3 Methodological issues

When Anderson proved his Theorem 2, it immediately lead to fundamental methodological questions: can one trust a statistical estimate that has an infinite mean square error (not to mention infinite bias)? Can such an estimate be better than others which have finite moments?

Fitting lines. To explore this issue, Anderson 1976 and 1984, Kunitomo 1980, and others compared the MLE estimate $\hat{\beta}$ given by (6) with the classical estimate $\hat{\beta} = s_{xy}/s_{xx}$ of the slope of the regression line that is known to be optimal when x_i 's are error-free (i.e., $\delta_i = 0$). They denote the former by $\hat{\beta}_M$ (Maximum likelihood) and the latter by $\hat{\beta}_L$ (Least squares); of course, both estimates were studied in the framework of the EIV model described in Section 2. Their results can be summarized in two seemingly conflicting verdicts:

- (a) The mean square error of $\hat{\beta}_M$ is infinite, and that of $\hat{\beta}_L$ is finite (whenever $n \geq 4$), thus $\hat{\beta}_L$ appears (infinitely!) more accurate;

- (b) The estimate $\hat{\beta}_M$ is consistent and asymptotically unbiased, while $\hat{\beta}_L$ is inconsistent and asymptotically biased (unless $\beta = 0$).

Besides, Anderson 1976 shows that if $\beta \neq 0$, then

$$P(|\hat{\beta}_M - \beta| > t) < P(|\hat{\beta}_L - \beta| > t)$$

for all $t > 0$ of practical interest, i.e. the accuracy of $\hat{\beta}_M$ dominates that of $\hat{\beta}_L$ everywhere, except for very large deviations (large t). It is the heavy tails of $\hat{\beta}_M$ that make its mean square error infinite, otherwise it tends to be closer to β than its rival $\hat{\beta}_L$.

Anderson 1976 remarks that this situation, in its extreme, resembles the following dilemma: suppose we are estimating a parameter θ whose true value is $\theta^* \approx 0$, and we have to choose between two estimates: one, $\hat{\theta}_1$, has Cauchy distribution, and the other, $\hat{\theta}_2$, has a normal distribution with mean 100 and variance 1. Would anyone prefer $\hat{\theta}_2$ only because it has finite moments?

Thus Anderson and others build a very strong case supporting the MLE estimate $\hat{\beta}_M$, despite its infinite moments. Furthermore, Gleser 1983 proves that the MLE estimate $\hat{\beta}_M$ is the best possible in a certain formal sense, we refer the reader to Chen and Van Ness 1994 for a detailed survey.

Fitting circles. Now we return to the circle fitting problem. It allows an alternative approach: instead of minimizing geometric distances (1)–(2) one can minimize the so-called ‘algebraic distances’:

$$(7) \quad (\hat{a}_0, \hat{b}_0, \hat{R}_0) = \operatorname{argmin} \sum_{i=1}^n [(x_i - a)^2 + (y_i - b)^2 - R^2]^2.$$

By changing parameters $A = -2a$, $B = -2b$, and $C = a^2 + b^2 - R^2$ one reduces (7) to a *linear* least squares problem

$$(8) \quad (\hat{a}_0, \hat{b}_0, \hat{R}_0) = \operatorname{argmin} \sum_{i=1}^n [x_i^2 + y_i^2 + Ax_i + By_i + C]^2$$

which has a unique and explicit solution. This approach is known as a *simple algebraic fit* (see Chernov and Lesort 2005) or Delogne-Kása method (Zelniker and Clarkson 2006); it was introduced in the 1970s. It has an obvious advantage of simplicity over the geometric fit, which requires iterative numerical schemes.

The competition between the geometric and algebraic circle fits is now over 30 years old, and so far it was focused on simplicity versus accuracy. Geometric estimates $(\hat{a}, \hat{b}, \hat{R})$ are widely known to be extremely accurate in practical applications, despite their slight tendency to overestimate the circle's radius (the latter was pointed out by Berman 1989). On the other hand, the Delogne-Kåsa estimates are heavily biased toward smaller circles, see Chernov and Lesort 2004 and 2005 and references therein, and generally much less accurate than the geometric estimates.

Now this competition acquires a new, purely statistical momentum. Recently Zelniker and Clarkson 2006 proved that the Delogne-Kåsa estimates $(\hat{a}_0, \hat{b}_0, \hat{R}_0)$ have finite mean values whenever $n > 3$ and finite variances whenever $n > 4$. Our work shows that the geometric estimates $(\hat{a}, \hat{b}, \hat{R})$ have infinite moments.

This competition very much resembles the one described above between the two line slope estimates: the MLE $\hat{\beta}_M$ and the 'classical least squares' $\hat{\beta}_L$. It would be interesting to further compare the two circle fits along the lines of the cited works by Anderson, Kunitomo, Gleser, and others, but this is perhaps a research program for distant future.

Alternative parametrizations. One can also say that non-existence of moments is an artifact of a poorly chosen parametrization, and the problem is easily remedied by changing parameters. In the case of lines, one can replace its slope β with the angle θ the line makes with, say, the y -axis. Then the line can be described as $x \cos \theta + y \sin \theta + d = 0$. Now it is easy to check that the estimates of θ and d have finite moments. These parameters are commonly used after Anderson's work in 1976.

In the case of fitting circles, the radius R can be replaced with the curvature $\rho = 1/R$. It is easy to check that the estimate of ρ has finite moments (up to the order $2n - 3$). The center coordinates (a, b) can be replaced by, say, $c = a/R$ and $d = b/R$, which would also have finite moments. Alternatively one can replace them with (q, θ) defined by $a = q^{-1} \cos \theta$ and $b = q^{-1} \sin \theta$ (V. Clakson, private communication). All these new parameters have finite moments.

Alternatively, one can describe circles by equation

$$A(x^2 + y^2) + Bx + Cy + D = 0$$

subject to constraint $B^2 + C^2 - 4AD = 1$; this was proposed by Pratt 1987. Now the parameters (A, B, C, D) are defined uniquely; and using the results of Chernov and Lesort 2005 it is easy to check that they have finite moments.

4 Proof of Theorem 1

It is enough to prove our theorem for the functional model. Indeed, then in the context of the structural model the conditional expectations of $|\hat{a}|$, $|\hat{b}|$, and \hat{R} for every given realization of $\varphi_1, \dots, \varphi_n$ will be infinite, thus their unconditional expectations will be infinite, too.

Next we need to make a few general remarks. First, the objective function (1) may not have a minimum. For example, if the data points are collinear, then $\inf \mathcal{F}(a, b, R) = 0$, but there is no circle that would interpolate $n > 2$ distinct collinear points, hence $\mathcal{F}(a, b, R) > 0$ for all a, b, R . In that case the best fit is achieved by a line, which can be regarded as a ‘degenerate circular arc with infinite radius’.

It is proved in Chernov and Lesort 2005 that if one poses the circle fitting problem in this ‘extended sense’, i.e. as finding a circle *or a line* which minimizes the sum of squares of distances to the given data points, then the problem always has a solution. That is, the best fitting contour (a circle or a line) always exists. The solution may not be unique, though, as the global minimum of the objective function (1) can be attained simultaneously on several distinct circles, examples are given in Chernov and Lesort 2005 and Zelniker and Clarkson 2006.

In the case of multiple solutions, any one can be selected, our theorem remains valid for any selection. If the best fit is a line, rather than a circle (for example, if the data are collinear), then we can set $\hat{a} = \hat{b} = \hat{R} = \infty$.

This fact by itself does not prove our theorem, of course, as the probability of such an exceptional event is zero. It shows, however, that in nearly collinear cases the estimates $\hat{a}, \hat{b}, \hat{R}$ tend to take arbitrarily large values, and we will explore this tendency thoroughly.

Simple case $n = 3$. Our argument is particularly simple if $n = 3$, and this case also illustrates our main idea.

Let 3 data points be located at $(0, 0)$, $(0, -1)$ and $(x, 1+y)$ where x and y are small, say $\max\{|x|, |y|\} \leq h = 10^{-9}$. Note that for $n = 3$ the best fitting circle simply interpolates the three given points, so by elementary geometry $\hat{a} = (2 + 3y + y^2 + 2x^2)/(4x)$, in particular $|\hat{a}| \geq 1/(3|x|)$. Since the density of $(x_3, y_3) = (x, 1+y)$ is continuous and positive, it has a minimum value $p_0 > 0$ in the rectangle $|x_3| \leq h$, $|y_3 - 1| \leq h$. Therefore the conditional

expectation of $|\hat{a}|$, when the other two points are fixed, is

$$E(|\hat{a}|/B) \geq p_0 \int_{1-h}^{1+h} \int_{-h}^h \frac{1}{3|x|} dx dy = \infty,$$

where $B = \{(x_1, y_1) = (0, 0), (x_2, y_2) = (0, -1)\}$.

A similar estimate holds if the points $(0, 0)$ and $(0, -1)$ are perturbed slightly, say within a little square of size h^2 around their initial positions. Now the densities of (x_1, y_1) and (x_2, y_2) are also positive, so a direct integration yields $E(|\hat{a}|) = \infty$. It is also clear that $E(\hat{R}) = \infty$. Rotating our construction, say by $\pi/2$, we obtain $E(|\hat{b}|) = \infty$, too.

General case $n > 3$. We modify our previous construction as follows. Let $h = 10^{-9}n^{-2}$ (here 10^{-9} may be replaced with any sufficiently small constant).

We place our first point (x_1, y_1) in the ‘lower’ square $[-h^2, h^2] \times [-1 - h^2, -1 + h^2]$, then $n - 2$ points (x_i, y_i) , $i = 2, \dots, n - 1$, in the ‘central’ square $[-h^2, h^2] \times [-h^2, h^2]$, and the last point (x_n, y_n) in the (horizontally extended) ‘upper’ rectangle $[-h, h] \times [1 - h^2, 1 + h^2]$.

For every fixed positions of the first $n - 1$ points and the fixed y -coordinate y_n of the last point, we will examine how the best fitting circle changes as the x -coordinate $x = x_n$ of the last point changes from $-h$ to h . Let $\hat{a}(x)$ denote the first coordinate of the circle’s center (we suppress its dependence on the other x_i and y_i coordinates). If the best fit is a line (and that line is clearly almost vertical), we set $\hat{a} = \infty$. Since \hat{a} is large, it is more convenient to work with $\zeta(x) = 1/\hat{a}(x)$, which is always finite and small.

Observe that all our points (x_i, y_i) , $1 \leq i \leq n$, are located in the h^2 -vicinity of three points: $(0, 0)$, $(0, -1)$, and $(x, 1)$, thus the best fitting circle (or line) passes through the h^2 -vicinity of these three points, too. By elementary geometry, if $x = h$, then $\hat{a}(x) > 1/(2h)$, hence $\zeta(h) \in (0, 2h)$. Similarly, $\hat{a}(-h) < -1/(2h)$, hence $\zeta(-h) \in (-2h, 0)$.

As $x = x_n$ changes from $-h$ to h , the $\zeta(x)$ function moves from the negative interval $(-2h, 0)$ into the positive interval $(0, 2h)$, and it stays between $-2h$ and $2h$. All we need now is that $\zeta(x)$ behave regularly in the following sense:

Lemma 3 (Regularity). *For any fixed values (x_i, y_i) , $1 \leq i \leq n - 1$, and y_n , as above, the function $\zeta(x)$ is differentiable and its derivative is bounded, i.e. $|\zeta'(x)| \leq D$ for some constant $D > 0$. Here D may depend on n and h but not on the fixed coordinates (x_i, y_i) .*

The proof of Lemma is rather technical; it is given in Appendix.

Proof of Theorem 1. Due to the regularity lemma, the function $\zeta(x)$ is continuous, hence $\zeta(x_0) = 0$ for some $x_0 \in (-h, h)$. The boundedness of the derivative $\zeta'(x)$ implies that for any $\varepsilon > 0$ if $|x - x_0| < \varepsilon$, then $|\zeta(x)| < D\varepsilon$, hence $|\hat{a}(x)| > 1/(D\varepsilon)$. The conditional probability of this event (when (x_i, y_i) for $i = 1, \dots, n-1$ and y_n are fixed) is $\geq p_0\varepsilon$ with some constant $p_0 > 0$, due to the positivity of the density of (x_n, y_n) . Therefore again, as in the $n = 3$ case, the conditional expectation of $|\hat{a}|$ is infinite, hence so is the unconditional expectation due to the positivity of the densities of (x_n, y_n) , $i = 1, \dots, n-1$.

Our analysis also implies $E(\hat{R}) = \infty$. Rotating our construction by $\pi/2$ gives $E(|\hat{b}|) = \infty$. \square

Radial model. Our theorem can be extended to another interesting model for the circle fitting problem proposed by Berman and Culpin 1986 and further studied by Chernov and Lesort 2004. In this model the error vector $\mathbf{e}_i = (\delta_i, \varepsilon_i)$, cf. (3), satisfies $\mathbf{e}_i = \xi_i \mathbf{n}_i$, where \mathbf{n}_i is a unit normal vector to the true circle at the true point (x_i^*, y_i^*) and ξ_i 's are independent normally distributed random variables with zero mean. In other words, the noise $(\delta_i, \varepsilon_i)$ is normal but restricted to the radial direction (perpendicular to the circle).

To extend our theorem to this model we need to assume that there are at least three distinct true points (x_i^*, y_i^*) on the true circumference. We outline the modifications in our argument needed to cover this new case.

Clearly it is possible that all the data points are collinear, i.e. there is a line L such that the probability that all the data points lie in the h -vicinity of L is positive for any $h > 0$. Also, for at least one data point its radial direction (on which its distribution is concentrated) must be transversal to L . Let that point be (x_n, y_n) . Now we can repeat our construction by moving (x_n, y_n) across L and keeping all the other points fixed in a tiny vicinity of L . The technical analysis only requires minor modifications in this new case, so we omit details.

Acknowledgement. The author is partially supported by NSF grant DMS-0652896.

Appendix

Here we prove our regularity Lemma 3. First we eliminate R from the picture. The objective function (1) is a quadratic polynomial in R , hence it has a unique global minimum in R when the other two variables a and b are kept fixed, and it is attained at

$$(9) \quad R = R(a, b) = \frac{1}{n} \sum_{i=1}^n \sqrt{(x_i - a)^2 + (y_i - b)^2}.$$

This allows us to express \mathcal{F} as a function of a and b only:

$$(10) \quad \begin{aligned} \mathcal{F}(a, b) &= \sum_{i=1}^n [\sqrt{(x_i - a)^2 + (y_i - b)^2} - R(a, b)]^2 \\ &= n[\bar{z} - 2a\bar{x} - 2b\bar{y} + a^2 + b^2] - n[R(a, b)]^2, \end{aligned}$$

where for brevity we denote $z_i = x_i^2 + y_i^2$; here and on we use standard statistical ‘sample mean’ notation $\bar{z} = \frac{1}{n} \sum z_i$, $\bar{x} = \frac{1}{n} \sum x_i$, etc.

Next we switch to polar coordinates $a = \rho \cos \theta$ and $b = \rho \sin \theta$ in which (10) takes form

$$(11) \quad \begin{aligned} \frac{1}{n} \mathcal{F}(\rho, \theta) &= \bar{z} - 2\rho(\bar{x} \cos \theta + \bar{y} \sin \theta) + \rho^2 \\ &\quad - \left[\frac{1}{n} \sum \sqrt{z_i - 2\rho(x_i \cos \theta + y_i \sin \theta) + \rho^2} \right]^2. \end{aligned}$$

Note that \mathcal{F} in (10) and (11) denotes the same function, though expressed in different sets of variables. We introduce more convenient notation

$$u_i = x_i \cos \theta + y_i \sin \theta \quad \text{and} \quad v_i = -x_i \sin \theta + y_i \cos \theta$$

(observe that $u_i^2 + v_i^2 = z_i$), so that (11) becomes shorter:

$$(12) \quad \frac{1}{n} \mathcal{F}(\rho, \theta) = \bar{z} - 2\rho\bar{u} + \rho^2 - \left[\frac{1}{n} \sum \sqrt{z_i - 2\rho u_i + \rho^2} \right]^2.$$

Now we introduce another variable

$$(13) \quad \begin{aligned} w_i &= \rho[\sqrt{z_i - 2\rho u_i + \rho^2} - (\rho - u_i)] \\ &= \frac{v_i^2}{\sqrt{1 - 2u_i\rho^{-1} + z_i\rho^{-2}} + 1 - u_i\rho^{-1}}. \end{aligned}$$

From (13) we have $\sqrt{z_i - 2\rho u_i + \rho^2} = \rho - u_i + w_i \rho^{-1}$, hence

$$\frac{1}{n} \sum \sqrt{z_i - 2\rho u_i + \rho^2} = \rho - \bar{u} + \bar{w} \rho^{-1}.$$

Now (12) takes form

$$(14) \quad \frac{1}{n} \mathcal{F}(\rho, \theta) = \bar{z} - \bar{u}^2 - 2\bar{w} + 2\bar{u}\bar{w}\rho^{-1} - \bar{w}^2\rho^{-2}.$$

By elementary geometry, the (averaged) objective function $\frac{1}{n} \mathcal{F}$ takes all its small values (say, all values less than $h^2/10$) on circles and lines that pass in the h -vicinity of the three basic points: $(0, 0)$, $(0, -1)$ and $(0, 1)$. These circles and lines have parameters restricted to the region where $\rho > 1/(100h)$ and $|\sin \theta| < 100h$.

We replace the large parameter ρ with its reciprocal $\delta = \rho^{-1}$ and obtain

$$(15) \quad \frac{1}{n} \mathcal{F}(\delta, \theta) = \bar{z} - \bar{u}^2 - 2\bar{w} + 2\bar{u}\bar{w}\delta - \bar{w}^2\delta^2,$$

where

$$w_i = \frac{v_i^2}{\sqrt{1 - 2u_i\delta + z_i\delta^2 + 1 - u_i\delta}}$$

(recall that u_i 's and v_i 's depend on θ but not on ρ).

Observe that the transformation $\theta \mapsto \theta + \pi$ and $\delta \mapsto -\delta$ leaves w_i 's and $\mathcal{F}(\delta, \theta)$ unchanged; thus we can let δ take (small) negative values but keep θ close to 0. More precisely, we can restrict our analysis to the region

$$(16) \quad \Omega = \{|\delta| \leq 100h \quad \text{and} \quad |\theta| \leq 100h\}.$$

Now one can easily see that the function $\mathcal{F}(\delta, \theta)$ in Ω is regular in the following sense: it is continuous and has bounded first and second derivatives (including partial derivatives) with respect to its variables δ and θ and with respect to $x = x_n$. We denote the first derivatives by \mathcal{F}_δ , \mathcal{F}_θ , \mathcal{F}_x and second derivatives by $\mathcal{F}_{\delta\delta}$, $\mathcal{F}_{\delta\theta}$, etc.

All these derivatives are uniformly bounded by a constant $M > 0$ that may depend on n and h but not on the other point coordinates.

By direct differentiation of $\mathcal{F}(\delta, \theta)$ we see that

$$(17) \quad \mathcal{F}_{\delta\delta} = 1 - \frac{2}{n} + \chi_1, \quad \mathcal{F}_{\theta\theta} = 4 + \chi_2, \quad \mathcal{F}_{\delta\theta} = \chi_3,$$

where χ_i are various small quantities (that can be made as small as we please by further decreasing h). Thus, \mathcal{F} is a convex function that has exactly one minimum in Ω and no other critical points.

Let $(\hat{\delta}, \hat{\theta})$ denote that unique minimum. Differentiating equations

$$\mathcal{F}_{\delta}(\hat{\delta}, \hat{\theta}) = 0 \quad \text{and} \quad \mathcal{F}_{\theta}(\hat{\delta}, \hat{\theta}) = 0$$

with respect to x gives

$$\begin{aligned} \mathcal{F}_{\delta\delta}(\hat{\delta}, \hat{\theta}) \hat{\delta}' + \mathcal{F}_{\delta\theta}(\hat{\delta}, \hat{\theta}) \hat{\theta}' + \mathcal{F}_{\delta x}(\hat{\delta}, \hat{\theta}) &= 0 \\ \mathcal{F}_{\theta\delta}(\hat{\delta}, \hat{\theta}) \hat{\delta}' + \mathcal{F}_{\theta\theta}(\hat{\delta}, \hat{\theta}) \hat{\theta}' + \mathcal{F}_{\theta x}(\hat{\delta}, \hat{\theta}) &= 0, \end{aligned}$$

where $\hat{\delta}'$ and $\hat{\theta}'$ denote the derivatives with respect to x .

Since all partial derivatives are uniformly bounded by M and the determinant is $\approx 4 - \frac{8}{n}$ due to (17), we have that $|\hat{\delta}'| \leq 2M$ and $|\hat{\theta}'| \leq 2M$. Lastly, recall that $\zeta = 1/\hat{a} = \hat{\delta}/\cos\hat{\theta}$, hence

$$|\zeta'| = \left| \frac{\hat{\theta}' \sin \hat{\theta}}{\cos^2 \hat{\theta}} \hat{\delta} + \frac{\hat{\delta}'}{\cos \hat{\theta}} \right| \leq 4M,$$

which proves the lemma with $D = 4M$. □

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