

Comparative testing of ellipse-fitting algorithms: implications for analysis of strain and curvature

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Abstract

Several types of geological problem involve fitting an ellipse to sparse data in order to define a property such as strain or curvature. The sensitivity of ellipse-fitting algorithms to noise in input geological data is often poorly documented. Here we compare the performance of some well-known approaches to this problem in geology against each other and an algorithm developed for machine vision. The specific methods tested here are an analytical method, a Mohr diagram method, two least squares methods and a constrained ellipse approach. The algorithms were tested on artificial datasets of known elliptical and noise properties. These results allow the selection of an ellipse fitting method in a variety of geological applications and also allow an assessment of the absolute and relative accuracy of a chosen method for various combinations of sample numbers and noise levels. For the most accurate semi axis magnitude and orientation estimates with five or more input data, the ‘mean object’ least squares approach is recommended. However, the other least squares method also yields good results and is also suitable for three or four data points. Where curvature data is being assessed, the least squares method is preferred as it can handle negative principal curvature values.

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1. Introduction

Determination of principal vectors, or maximum and minimum magnitudes and directions from sparse or scattered data, is a feature of several types of problem in structural geology. Examples are determining two-dimensional strain in classic ‘stretched belemnite’-type problems (Lisle and Ragan, 1988 and references therein) and estimation of principal surface curvatures (Lisle and Robinson, 1995; Belfield, 2000). With these applications in mind, in this paper we investigate the sensitivity of prevailing algorithms to quality of input data.

Strain (in two dimensions) and curvature can be represented by second order tensors (Ramsay and Huber, 1983). Tensors allow the description of the variation of a property independently of the co-ordinate system used to

define the locations of the property. However, they are linear operators that send a vector to a vector so a description of the co-ordinate system is required for practical purposes. In two-dimensional Cartesian space, four components are generally required to fully define a tensor. In the examples mentioned here, only three components are needed since strain and surface curvature can be represented by symmetric second order tensors with mutually orthogonal principal axes (Ramsay and Huber, 1983; Lisle and Robinson, 1995). These properties of symmetrical second order tensors allow them to be represented by an ellipse, which are defined by the magnitudes of two mutually orthogonal principal axes and an orientation of the major axis. Strain or curvature data usually have to be measured in directions that are likely to be arbitrary with respect to the principal axes. Therefore, the principal axes have to be derived from these arbitrarily oriented measurements by fitting ellipses to the sampled data (Lisle and Ragan, 1988; Lisle and Robinson, 1995; Stewart and Podolski, 1998).

Many methods are available for estimating principal axes from arbitrarily oriented measurements in geological

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situations; these methods are generally based upon numerical ellipse fitting or an equivalent method such as use of Mohr circles (Ramsay and Huber, 1983; Erslev and Ge, 1990; Lisle, 1994; Lisle and Robinson, 1995; Belfield, 2000; Stewart and Wynn, 2000). However, barring a few comments in key publications (DePaor, 1988; Robin, 2002), there are no published indications or metrics of relative performance of these algorithms with respect to geological data. So, a geoscientist approaching a problem of this nature is faced with choosing an algorithm without comparative information. The choice is made more difficult by the fact that several algorithms used in geological situations are designed to give a precise answer from a restricted number of measurements (DePaor, 1988; Lisle and Ragan, 1988), rather than yielding a best fit from a larger number of measurements that include some level of noise (Erslev and Ge, 1990; McNaught, 2002). It should be noted that the algorithms tested here all belong to the group of methods classified as boundary based methods. These methods all attempt to fit ellipses to a set of points that represent the boundary of an elliptical region (Mulchrone and Choudhury, 2004). An alternative group of methods are the region based methods that use the moments of a shape to derive a best fit ellipse that describes the shape (Hu, 1962; Teague, 1980). The advantage of region based methods is that they are less sensitive to boundary irregularities than the boundary based methods (Mulchrone and Choudhury, 2004). However, the purpose of this paper is to compare the performance of algorithms within the class of boundary based methods that are common in geological applications.

In this study we set out to test key methods against each other using data sampled from a wide range of purpose-built, ‘noisy’ ellipses. Rather than a detailed critique of the theoretical merits of each algorithm, the intent of this study was to calibrate their performance for practical application, so that a preferred algorithm could be chosen for ellipse estimation without a priori knowledge of dataset parameters such as noise characteristics and principal axis orientations.

2. Ellipse fitting algorithms

Published methods previously applied to geological data can be grouped into ‘families’ of essentially similar approaches, namely:

- (1) Analytical, or algebraic methods
- (2) Mohr circle methods
- (3) Least squares methods (algebraic and geometric distances)

In this study, we chose a representative method or methods from each group to go forward into the testing process. The rationale behind the choice of a representative from each ‘family’ follows, together with a more detailed description of each selected algorithm. Within the group of

least squares methods, algebraic and geometric distances refer to the parameters being minimised within the second order polynomial equation representing the ellipse (see Fitzgibbon et al. (1999) for more details). Algebraic least squares solutions are linear and relatively simple to solve. However, geometric fitting gives a more meaningful parameter to minimise with respect to curved lines such as ellipses. Unfortunately, geometric distance minimisation methods are non-linear, requiring complex iterative solutions. It should be noted that hereafter in the paper the term ‘least squares’ refers to an algebraic distance, linear least squares minimisation. It should also be noted that all the methods require ellipse radii and their orientations as input and return estimates of the principal semi axes magnitudes and orientations.

2.1. Analytical method (AN)

A number of algebraic solutions for evaluating the strain ellipse from variously oriented longitudinal strain measurements have been published (Sanderson (1977), Robin (1983), Ragan (1987), DePaor (1988) and review by Lisle (1994)). The approach of DePaor (1988) was selected here because it is an exact solution and is computationally efficient. The method was designed to yield an ‘exact’ estimate of the principal axes of strain from three reciprocal quadratic stretches (λ'_a , λ'_b and λ'_c) in arbitrary directions, using three simultaneous quadratic equations. First, the angle of the major principal axis is found and this is then used to find the principal semi axes of the ellipse. This algorithm was published with strain analysis in mind but other tensor data, such as curvatures, can be used instead of stretches resulting in estimates of the principal reciprocal quadratic curvatures. A limitation of the method is the possibility of non-real solutions because all possible combinations of three radii of a quadratic do not necessarily lie on an ellipse and may be better represented by a hyperbola (DePaor, 1988). It is also the only method to use only three input points and is therefore not directly comparable with the other methods tested here. However, it is a common method and will be useful where data is very sparse and a quick computation is required.

2.2. Mohr circle method (MC)

Mohr circle methods have been widely used in strain and curvature problems (Means, 1982; Treagus, 1987; Lisle and Ragan, 1988; Lisle and Robinson, 1995). The formulation for the estimation of principal curvatures from gridded map data described by Lisle and Robinson (1995) is adopted here since it is designed specifically for computer implementation and deals with large digital datasets. The form of the construction is shown in Fig. 1 and is described briefly below for the case with four input points although the construction is valid for more points. The reader is referred

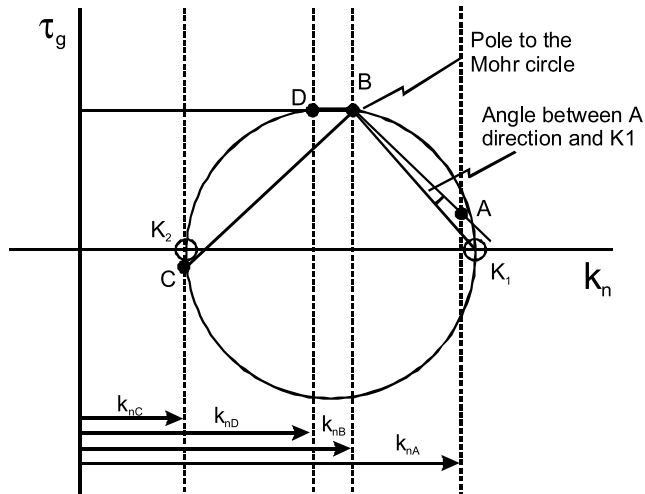


Fig. 1. Mohr diagram with pole construction marked used for the Mohr method. K_n is normal surface curvature, τ_g is surface curve torsion. See text for details.

to Lisle and Robinson (1995) for a more detailed explanation:

1. Obtain the data to plot on the horizontal axis of a Mohr diagram (e.g. normal curvature K_n or longitudinal strain e). In the case of curvature derived from gridded data, a node and its eight surrounding neighbours is chosen. A circle is fitted to the central node and each pair of opposing neighbours giving four circles in directions A–D, which are 45° apart. The curvature of these circular arcs (K_{nA} – K_{nD}) is found by reciprocating the fitted radii.
2. Use a graphical construction (Lisle and Ragan (1988) after Murphy (1945)) to generate the pole to the Mohr circle and plot values on the vertical axis (e.g. surface torsion τ_g or shear strain τ). Fig. 1 shows the τ_g values plotted at points A–D.
3. Fit a least squares circle to the data (Lisle, 1992), scale the Mohr diagram vertical axis and obtain the principal semi axis magnitudes (K_1 and K_2). Use the pole of the Mohr circle to obtain the principal semi axis orientations (Lisle and Ragan, 1988).

2.3. General least squares methods (MO and LS)

A further family of methods involves minimising ellipse-fitting errors using summed least squares of the shortest distance from input data to the fitted elliptical curve. Various forms of least squares ellipse fitting are employed in geological problems (Erslev and Ge, 1990; Hart and Rudman, 1997; Belfield, 2000), though algorithm design for least squares ellipse fitting has probably been studied in more detail in the context of pattern recognition (Rosin, 1993, 1996; Ahn and Rauh, 1999). Belfield (2000) used a least squares ellipse fit to measure principal curvatures of geological surfaces, adapting a general technique outlined

by Nye (1957) for calculating second order tensor principal axes. Erslev and Ge (1990) reported the use of a least squares algorithm for ellipse fitting of geological strain data and also implemented an estimate of the closeness of the data to the best fit ellipse using the ‘radial error’ parameter. The radial error is the absolute difference in distance between the data point and the ellipse margin on a line intersecting the ellipse centre, the ellipse margin and the data point, divided by the ellipse radius on that line. Therefore, radial error is measuring the difference between a data point and the fitted ellipse. However, in this study we are assessing the differences between the fitted ellipses and the original input ellipses before the addition of any synthetic noise. Therefore, the radial error approach was deemed inappropriate for use in this study.

The steps employed by Belfield (2000) for curvature analysis with four input points are described in more detail in Nye (1957). The approach is the same for n input points where $n \geq 3$. Obtain the normal surface curvature data A_1, A_2, \dots, A_n and their orientations $\theta_1, \theta_2, \dots, \theta_n$. Convert to reciprocal root curvatures. Let A equal the $n \times 1$ vector of normal curvature measurements and θ equal the $n \times 3$ matrix of sine and cosine functions of the orientations and α equal the 3×1 matrix of coefficients of the second rank tensor (Eq. (1)):

$$\alpha = (\theta_i \theta)^{-1} \theta^t A \quad (1)$$

Note that the 3×1 α matrix is derived from a symmetrical 3×3 matrix with the zero values and repeated elements removed. The superscript ‘t’ denotes the transpose and the superscript ‘-1’ the inverse. In this case, A and θ are known and α is unknown. In this example, α_{11} and α_{33} represent normal surface curvature estimates and α_{31} a surface torsion estimate. In strain analysis α_{11} and α_{33} represent longitudinal strain estimates and α_{31} a shear strain estimate. The matrix α represents the best estimates of tensor coefficients with the available input data. These values are then used in a Mohr circle construction to calculate the principal semi axis magnitudes and the orientation of the principal axis. The principal difference between this method and the Lisle and Robinson (1995) Mohr circle method is that it uses a least squares method on the original data points and calculates three best fit parameters for use in a simple Mohr circle construction.

The least squares method outlined in Erslev and Ge (1990) used different equations to those outlined above because it uses different coordinate systems but the methods are essentially equivalent. This method calculates the A, B and C coefficients of a centred conic. For given A, B and C values, a point (x_a, y_a) will satisfy the centred conic equation with a residual value k_a . If $k_a = 0$ and $B^2 - 4AC < 0$ then the point falls on an ellipse defined by the coefficients A, B and C . Standard least squares analysis can be used to minimise the deviations of the points from the best fit ellipse. The resultant partial differential equations can be simplified to three formulae with eight unique coefficients

that can be rearranged to yield the centred conic equation A , B and C values. To obtain data on the principal semi axis orientations and magnitudes, a further three formulae are required (see [Erslev and Ge, 1990](#) for more details).

Throughout this text the Erslev and Ge method is referred to as the mean object ellipse (MO) method. Strictly speaking, the method as it is described here is a least squares algorithm that is used as input to a mean object ellipse technique. However, the term is used to differentiate it from the least squares method outlined in [Belfield \(2000\)](#), which is termed the least squares (LS) method in this paper.

2.4. Constrained ellipse least squares method (CE)

This method was developed by image processing researchers as a significant refinement of least squares ellipse fitting ([Fitzgibbon et al., 1999](#)). The method is the first to utilise linear least squares error minimisation whilst also constraining the fitted conic to be an ellipse. The method was tested against other widely used approaches in pattern recognition applications (those of [Bookstein \(1979\)](#), [Taubin \(1991\)](#) and [Gander et al. \(1994\)](#)) by [Fitzgibbon et al. \(1999\)](#). These tests focused on the addition of Gaussian noise to the data and indicated that the constrained ellipse was the most robust approach. Based on this performance, we selected the constrained ellipse approach for comparison with the methods described so far. The method as presented in [Fitzgibbon et al. \(1999\)](#) is described briefly below. To fit ellipses directly whilst retaining the efficiency of a linear least squares approach, it is required that the appropriate constraint is always $b^2 - 4ac < 0$. In practice, this constrained problem is difficult to solve. This is circumvented by scaling the parameters to obtain the equality constraint $4ac - b^2 = 1$. This is a quadratic constraint, which may be expressed in matrix form. Although the minimisation only yields one elliptical solution, [Fitzgibbon et al. \(1999\)](#) noted that this ellipse is inherently biased toward low eccentricity. Eccentricity is defined in Eq. (2):

$$e = \sqrt{1 - \frac{b^2}{a^2}} \quad (2)$$

where $0 < e < 1$. It should be noted that e is not used in this study because it becomes asymptotic to 1 above axial ratios of 10 and the testing described here used axial ratios up to 60. However, lower eccentricities correlate with lower axial ratios.

3. Testing procedure

Each of the algorithms described above was fed the same data from an ellipse generating program and the estimated principal semi axis magnitudes (Ea , Eb) and orientations ($E\theta$) were compared with the known input values (a , b , θ).

This procedure was repeated with incremental variations in the following parameters of the input ellipse:

1. Ratio of the principal axes (a/b).
2. Position of the sampled points on the perimeter of the ellipse with respect to the location of the principal axes (θ).
3. Noise added to the points defining the perimeter of the ellipse.

Each ellipse is defined by 720 data points (0–359.5° at 0.5° intervals). These datasets were then sampled with 30 sample points 360 or 720 times corresponding to half or full ellipse sampling over 180 or 360°, respectively. These sample points had the same angular spacings to provide input to the algorithms under investigation. Equal angular spacing was employed to reduce potential bias from arbitrary sample directions and also to simulate the automatic application of these algorithms to large gridded datasets or image processing applications relevant to deformed objects on outcrops or thin sections. An example of a source ellipse with and without noise is shown in [Fig. 2](#).

The noise distribution model used here was Gaussian, though other distributions could have been chosen (uniform, Poisson, etc.). A symmetrical distribution was desired to avoid masking consistent error trends in the algorithms. It could be argued that a deformed Gaussian distribution for noise should have been used to represent the effects of strain on a 'natural' noise distribution present in the original data.

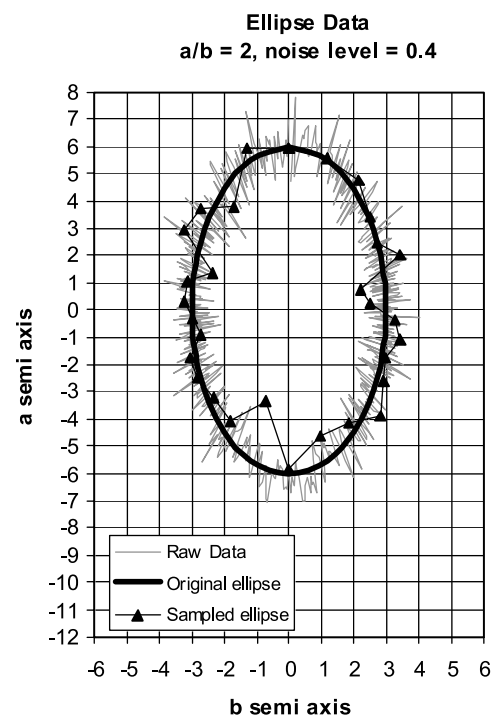


Fig. 2. Input data for the ellipse algorithms. Raw data (red curve) has added noise with a normal distribution with a mean of 0 and a maximum difference of $\pm 40\%$ of ellipse radii magnitude. Original data = blue curve, sampled ellipse = yellow curve.

However, this would only be suitable to strain measurements and would also imply that no noise or error existed in the sample measurement techniques. The source Gaussian noise distribution used had a mean of zero and standard deviation of ± 1 . This initial noise distribution was then normalised by the largest absolute value of the minimum and maximum to be $-1 < n < 1$. This noise distribution was then scaled again as a proportion (0.4) of the sampled ellipse radii length. The noise was then added to the original ellipse radii (see Fig. 2). The formula used is shown in Eq. (3):

$$r = r + ens \quad (3)$$

where r = ellipse radii, e = ellipse radii length, n = Gaussian random noise with mean of 0 and range between -1 and 1 and s = noise level scaling value of 0 or 0.4. This noise model is static in the sense that the added noise remains the same for each radius even when the ellipse sampling is rotated. This has the effect of the noise distribution only being sampled over a window width equal to $180/n$ or $360/n$ where n is the number of sample directions. Whilst this should be expected to produce a repeating pattern in the results every $180/n$ or $360/n$ degrees, it does not always do so.

It should be noted that for curvature data, one or both of the principal curvatures can be negative. For the least squares and Mohr circle algorithms this is not a problem as they both utilise a Mohr circle construction that can return negative principal semi axis magnitudes. However, the analytical, mean object and constrained ellipse methods can only calculate curvature where both values are positive.

4. Results

The size of the dataset output in this study precludes exhaustive reporting here. Instead we restrict discussion of results to key correlations and findings that might influence choice of algorithm. Throughout this discussion, a and b are the real major and minor ellipse semi axes (i.e. semi axis magnitudes of noise-free source ellipse; Fig. 2) and Ea and Eb are the corresponding estimates of major and minor semi axis yielded by each algorithm. In this section the terms a axis and b axis are often used on plots for conciseness but all plots and text in this paper refer to the estimated or actual ellipse semi axis magnitudes. $E\theta$ is the azimuth of the estimated a semi axis relative to the vertical or y axis in the local coordinate system (Fig. 2) and θ is the angle of the original ellipse a semi axis and the vertical or y axis. a/b is the axial ratio of the noise free source ellipse. Axial ratios up to 60 were examined in these experiments, since high axial ratios are relevant in very high strain shear zones and also represent a tendency towards cylindrical surface structures relevant in curvature analysis. However, the principle reason behind testing to very high axial ratios was to see at what point the algorithm performances degraded

noticeably. The a/b axis is displayed as a log scale in order to emphasise data in the $a/b < 10$ range, which are particularly relevant to strain analysis.

For the testing reported here, 30 input points were used for all the algorithms spread over 360° . This was to ensure full sampling of each ellipse and to avoid sampling problems in the least squares fitting methods. The exception to this is the analytical algorithm (DePaor, 1988), which is designed for use with only three input data. In addition, the Mohr circle algorithm was further tested with 30 points with 6° sample angles over 180° because of severe problems when using the 12° sample spacing (see below). The results examine algorithm performance by assessing the axial ratios of the estimated ellipses, and comparing estimates of each semi axis with the actual semi axis magnitude with noise levels of 0.4. Finally, the accuracy of estimated ellipse orientation is reported. It should be noted that the ellipse major and minor semi axis estimates for zero noise were perfect for most algorithms within numerical rounding errors. However, the analytical and Mohr circle methods were susceptible to some biases and problems with zero noise and these are discussed below.

The analytical method displayed artefacts in the Ea , Eb and $E\theta$ results at a semi axis angles of 60 , 150 , 240 and 330° . These directions correspond to the positions where the a semi axis direction bisects two of the sample directions and two of the three points have identical values. Therefore, the algorithm does not have three independent input values and cannot compute an ellipse. The Mohr circle method when used with the 12° sample spacing produced artefacts with zero noise levels. These appeared to worsen within certain a semi axis angle ranges that change systematically when the pole to the Mohr circle is changed. Fig. 3 shows the results obtained at a and b semi axes values of 6 and 3, respectively, using sample directions of 0 and 180° for the pole to the Mohr circle. The results are worse for the Ea values but the Eb values also show perturbations. With the addition of noise the results became unusable with many non-numeric values for the semi axis magnitudes and $E\theta$ values clustered around 45 and 135° . Therefore, further testing of the Mohr circle method was undertaken using 30 input points with 6° spacing over 180° . At zero noise, this produced perfect values of Ea and Eb , within numerical rounding errors, for all θ values. However, artefacts did occur with the addition of noise, which are discussed below.

4.1. Ellipse axial ratio variations

Fig. 4 shows estimated vs. actual ellipse eccentricity (Ea/Eb vs. a/b) for all the methods with noise levels of 0.4. It can be seen that for the least squares, constrained ellipse and mean object methods there are relatively small underestimates or overestimates of Ea/Eb up to values of 8 and then relatively large underestimates at larger axial ratios for the least squares and constrained ellipse methods. The analytical and Mohr circle methods display less

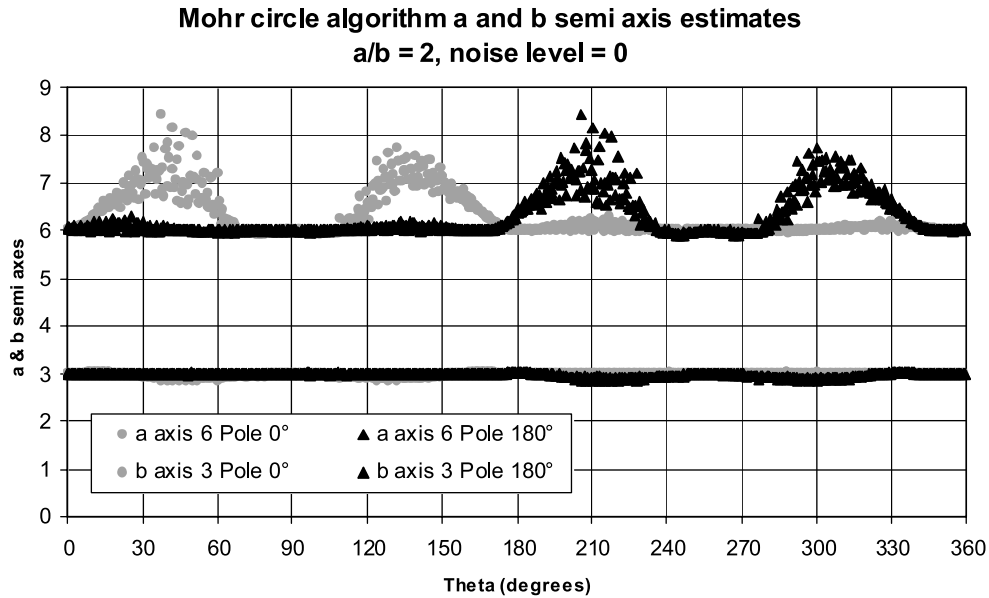


Fig. 3. Mohr circle algorithm output with pole directions of 0 and 180° and zero added noise. Note the repetition of artefacts within each set of data.

accuracy at all axial ratios. Most algorithms seem to yield a step change to underestimates at a/b values of 8. This is probably related to the noise level and the a/b value at which it occurs would be expected to change with different noise levels.

4.2. Estimated semi axis parameter variations with increasing a/b

On the results presented in Figs. 5–7 each plotted

ordinate value represents the mean of 720 separate estimates. The error bars on the plots represent one standard deviation of the ordinate parameter. This assumes that the distribution of estimates is Gaussian, though this may not necessarily be the case and they should only be used as a guide to the precision. Perfect estimates of each semi axis are represented by Ea/a or $Eb/b=1$, underestimates by Ea/a or $Eb/b < 1$ and overestimates by Ea/a or $Eb/b > 1$.

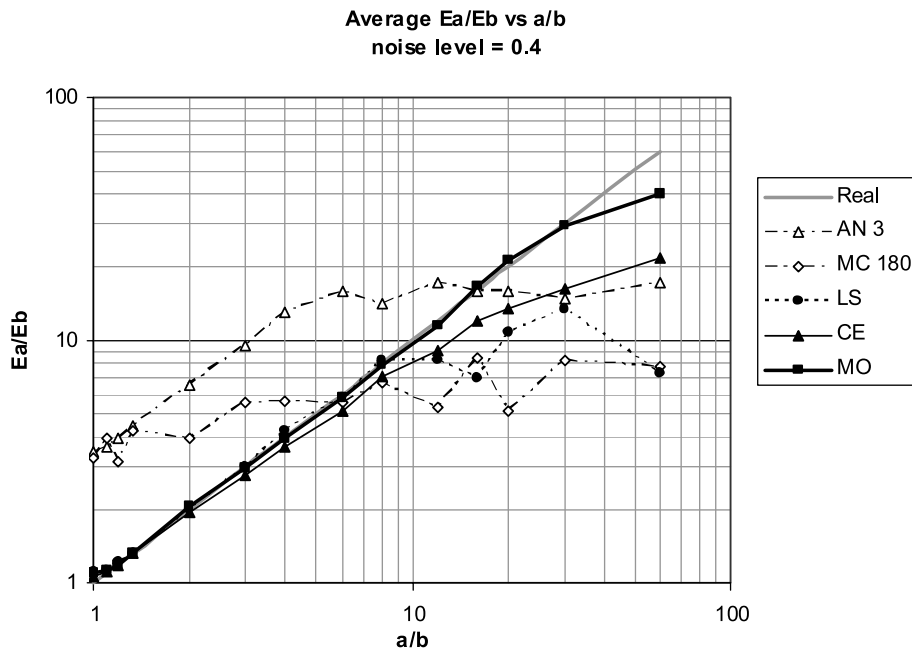


Fig. 4. Ea/Eb vs. a/b with noise levels of 0.4. See text for details. Key for all figures: AN=analytical, MC 180=Mohr circle with 180° sampling, LS=least squares, CE=constrained ellipse, MO=mean object ellipse. All methods use 30 input points at 12° increments apart from the AN algorithm with 3 input points and 120° sample increments and MC 180 with 30 input points and 6° sample increments.

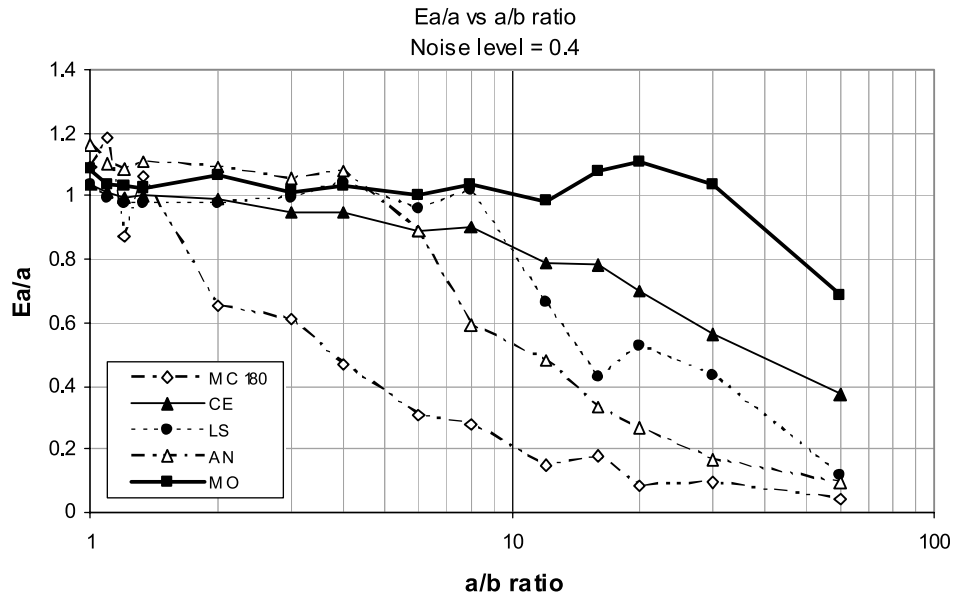


Fig. 5. Ea/a vs. a/b with noise levels of 0.4. See text for details.

4.2.1. Major semi axis

Fig. 5 shows variations in Ea/a vs. a/b at noise levels of 0.4. The mean object ellipse method is the most accurate and gives Ea/a values very close to 1 for a/b values up to 30 and then falls to 0.7 up to a/b ratios of 60. The Mohr circle method is the least accurate and shows values around 1 for axial ratios up to 1.3 and then progressively decreases to approach zero at axial ratios of >20 . The other three algorithms show behaviour intermediate between the mean object and Mohr circle methods.

4.2.2. Minor semi axis

Fig. 6 shows variations in Eb/b vs. a/b at noise levels of

0.4. Apart from the Mohr circle algorithm, all the algorithms show greater accuracy and precision than the Ea/a vs. a/b plot and generally vary between 0.95 and 1.05. The Mohr circle method shows underestimates at all axial ratios.

4.2.3. Major semi axis orientation

Fig. 7 shows the variations in $E\theta$ with a/b . For all algorithms there is a general increase in accuracy and precision with increasing a/b . This is expected as θ is less sensitive to added noise when a/b is larger. The mean object and least squares algorithms produce the best results and are very similar, both with average differences of $<5^\circ$ above a/b ratios of 2. The constrained ellipse algorithm shows the

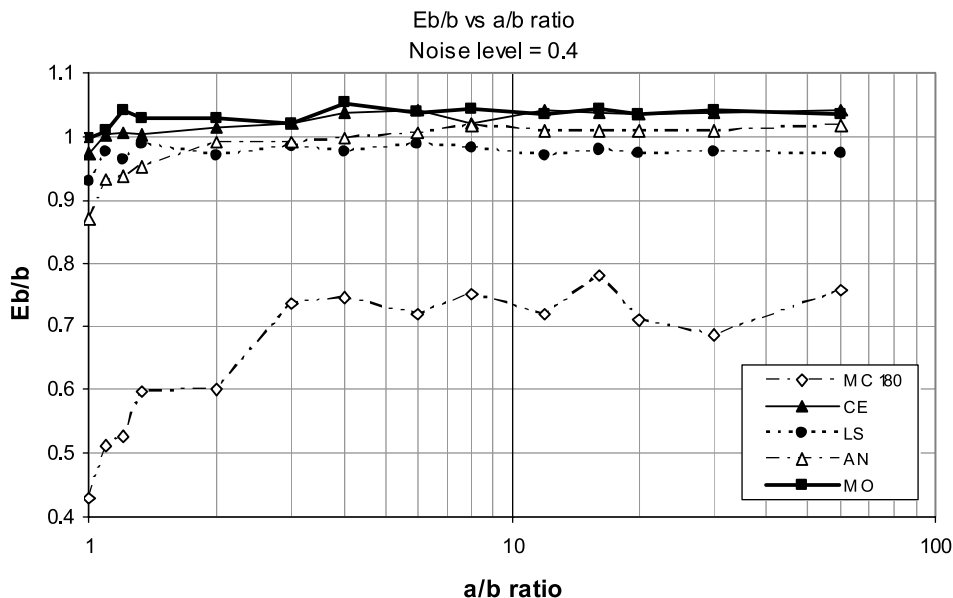


Fig. 6. Eb/b vs. a/b with noise levels of 0.4. See text for details.

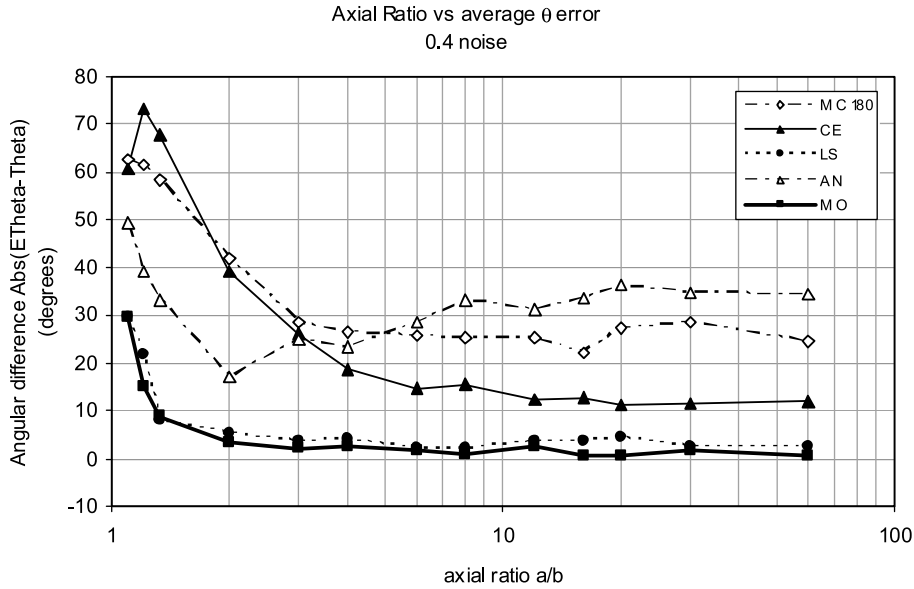


Fig. 7. $E\theta$ vs. a/b with noise levels of 0.4. See text for details.

worst results at a/b ratios less than 2 and then steadily increases in precision. The Mohr circle algorithm is similar to the constrained ellipse algorithm but shows higher average differences above a/b ratios of 4. The analytical algorithm is the only one that displays decreasing precision above a/b ratios of 4. This is probably related to the small number of sample directions (3), which will be much more sensitive to noise than the other algorithms.

The general increase in accuracy and precision for the Eb/b estimates compared with the Ea/a estimates can be attributed to the clustering of ellipse radii estimates around the b axis. Fig. 8 shows the variations in ellipse radii magnitudes with increasing angle away from the a axis. Fig. 8 shows that although there is a factor of 30 difference

between the largest and smallest ellipse major semi axes (180 and 6) there is only a factor of 3 difference between ellipse radii on these two ellipses at 10° from the long axis direction. Therefore, added noise will have a greater impact on the smaller number of points that are at or close to the a axis direction.

4.3. Estimated semi axis parameter variations with increasing theta

In contrast to previous plots that showed the averages of the algorithm performances for different a/b ratios and noise levels, most of the following plots (Figs. 9–15) show Ea vs. θ , Eb vs. θ and $E\theta$ vs. θ values for a/b ratios of 2 and 6 and

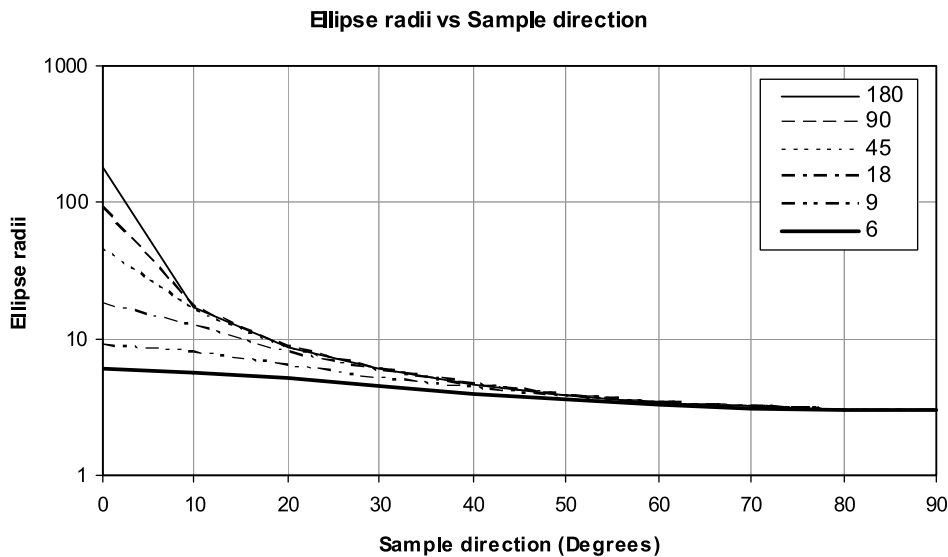


Fig. 8. Variations in ellipse radii magnitude with ellipse radii orientation for a/b ratios of 6 to 180. Note that significant differences with respect to the b semi axis magnitude only occur at ellipse radii orientations greater than $\pm 40^\circ$ from the b semi axis direction.

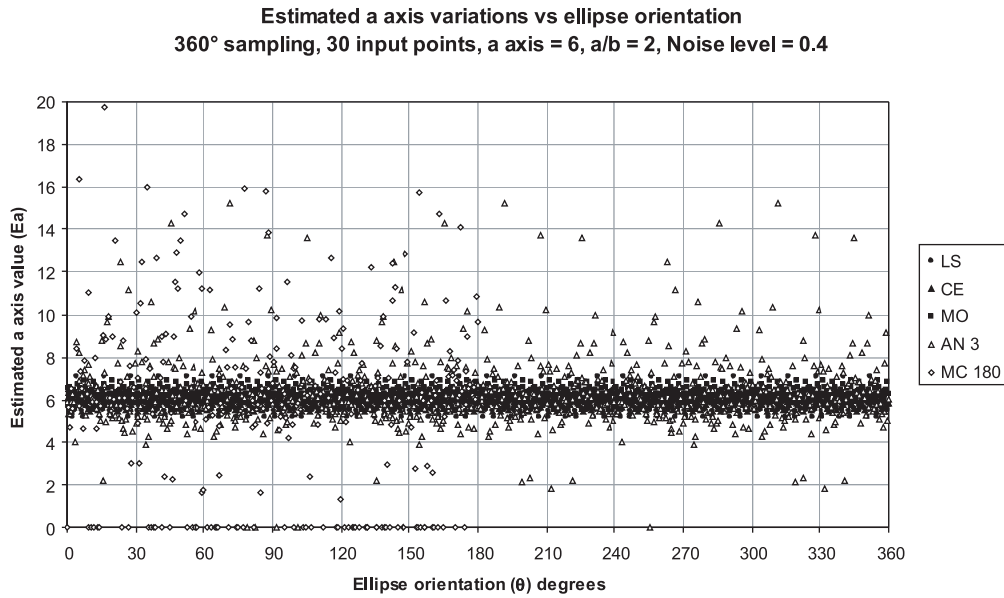


Fig. 9. E_a vs. θ , $a/b=2$. See text for details.

noise levels of 0.4. Effectively these plots are showing the population of points that make up the averages and standard deviations for the estimated values on Figs. 4–7.

4.3.1. Major semi axis

Figs. 9 and 10 show the variation in E_a values with ellipse orientation (θ) for a/b ratios of 2 and 6, respectively. Plots of this type should display repeating patterns of noise in the estimated values every 12° . This is because the ellipse a semi axis rotates through a sample direction at 12° intervals. These repetition intervals are expected to be 120° and 6° for the analytical and Mohr circle methods, respectively. The Mohr circle and analytical algorithms

show the worst accuracy and precision with a lot of scatter and many null values (plotting on the abscissa) on both plots. The analytical algorithm shows the repeating noise pattern but the Mohr circle algorithm does not. Fig. 11 shows the output from the Mohr circle algorithm only using pole directions of 0 and 90° . The 0° data is the same as that in Figs. 9 and 10. The data from the pole at 90° is the same as that for 0° but offset by either 84 or 96° . This indicates that the algorithm is not detecting the repeated noise pattern but the biased sampling is being repeated with different pole directions. This may be because it is a geometrical construction that uses the sample directions in conjunction with the abscissa data ($1/(\text{ellipse radii})^2$) to calculate the

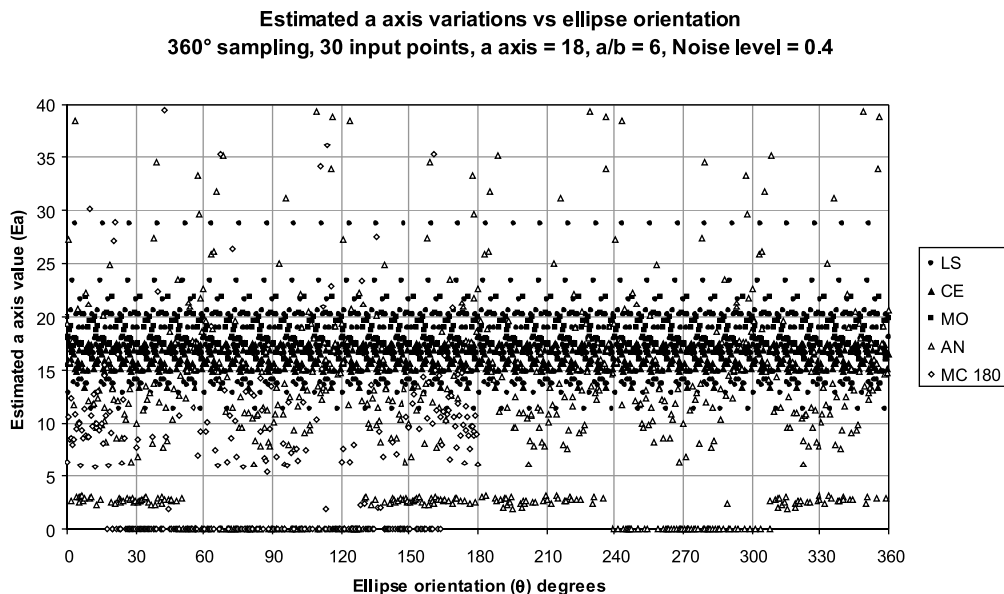


Fig. 10. E_a vs. θ , $a/b=6$. See text for details.

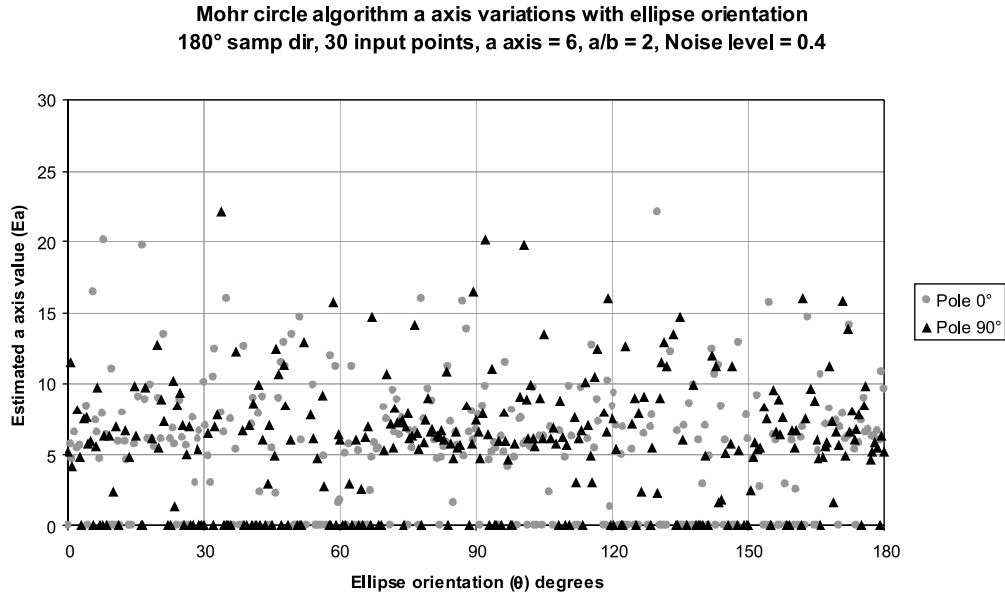


Fig. 11. Mohr circle method E_a vs. θ . $a/b=2$. Pole directions of 0 and 90°. See text for details.

ordinate data. When the sample directions are subvertical and the noisy input data plots on the abscissa slightly out of sequence compared with the pole location, the ordinate values will be much larger or smaller than if the pole was in a slightly different position on the abscissa. These errors are non-linear with respect to changes in the sample direction and reduce the effectiveness of the least squares circle fitting to noisy data on a Mohr diagram.

For the $a/b=2$ case in Fig. 9, the mean object, least squares and constrained ellipse algorithms display similar accuracies and precision with most points clustering around 6 ± 1 . The repeating pattern of noise is also evident in all three algorithms indicating no bias with respect to the

sample directions. For the $a/b=6$ case in Fig. 10 the data scatter is wider with most points clustering around 17 ± 5 . The least squares algorithm also displays a few more outliers than the constrained and mean object algorithms.

4.3.2. Minor semi axis

Figs. 12 and 13 show the variability in E_b vs. θ for a/b ratios of 2 and 6, respectively. There is less scatter in the results overall and greater similarity between the $a/b=2$ and $a/b=6$ results compared with the E_a vs. θ data. This indicates that the b semi axis estimates are generally more precise and accurate and less sensitive to axial ratio than the E_a values. This is probably due to the increased number of

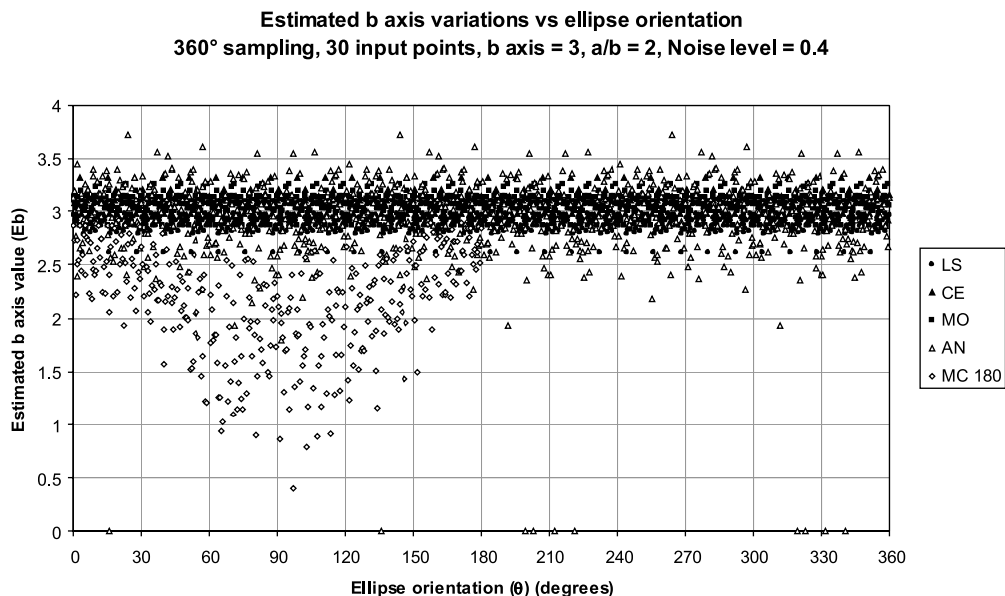


Fig. 12. E_b vs. θ , $a/b=2$. See text for details.

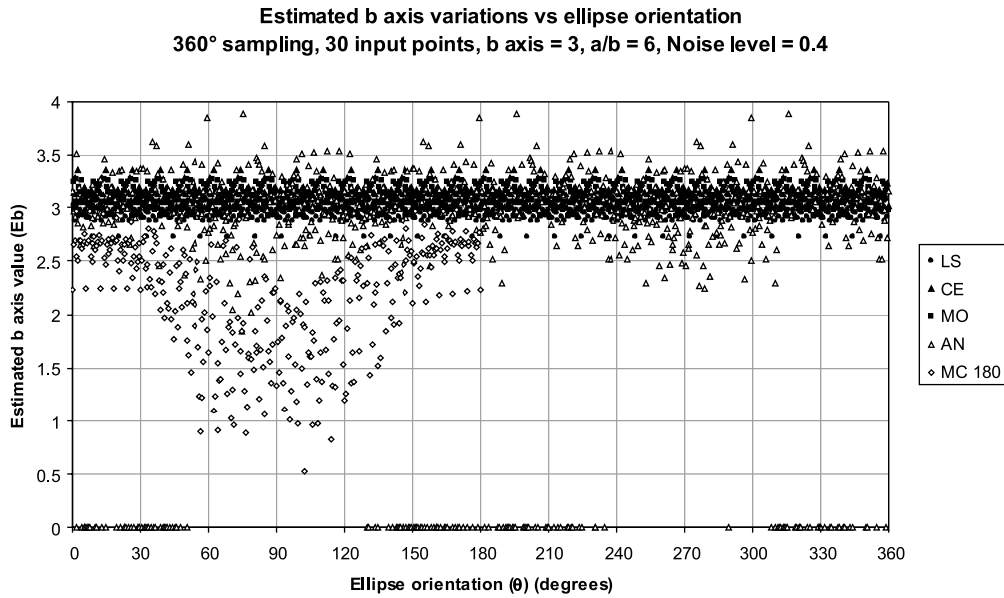


Fig. 13. E_b vs. θ , $a/b=6$. See text for details.

samples that are close to the b semi axis magnitude (see previous section and Fig. 8). The analytical and Mohr circle algorithms are the least accurate and least precise. The analytical algorithm displays many non- numerical results (plotting as zeros on Fig. 13) for the $a/b=6$ case and the Mohr circle method shows consistent underestimates of E_b in both plots.

4.3.3. Major semi axis orientation

Figs. 14 and 15 show $E\theta$ vs. θ for $a/b=2$ and $a/b=6$, respectively. The constrained, least squares and mean object methods show good accuracy and precision on both plots. The accuracy is greater in Fig. 15 because the ellipse axial

ratio is larger so the addition of noise has a reduced effect on $E\theta$. As with the E_a and E_b plots, the analytical and Mohr circle algorithms are the least accurate. In this case the analytical algorithm does not function very well with $a/b=6$ as there are many short intervals where $E\theta$ is offset from the $E\theta=\theta$ line by $70\text{--}90^\circ$ and the data has slightly gentler gradients than the other datasets. These intervals correlate with the null data in Figs. 11 and 13.

5. Discussion and conclusions

In noisy datasets, all the methods tested show some

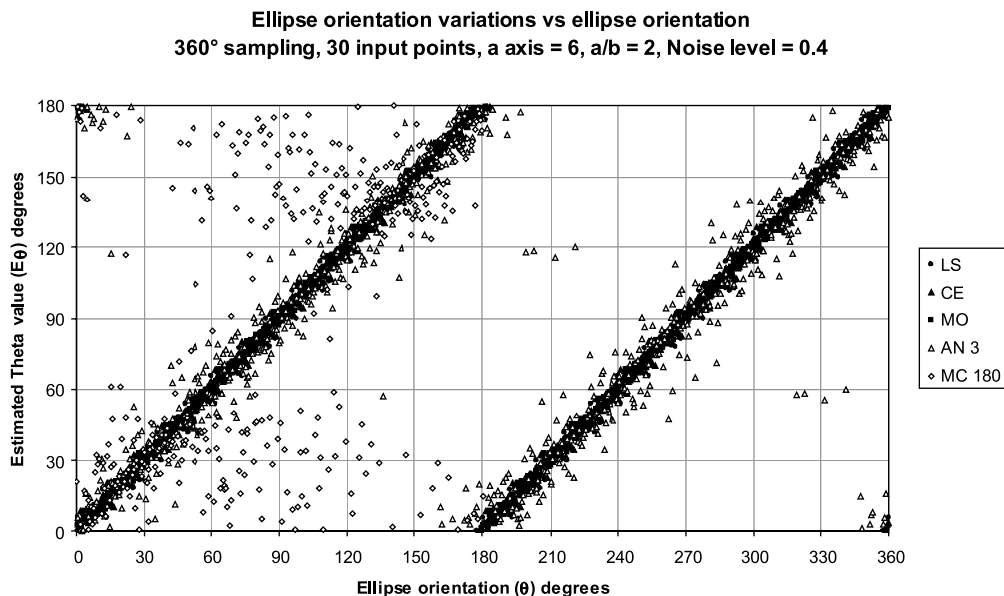


Fig. 14. E_θ vs. θ , $a/b=2$. See text for details.

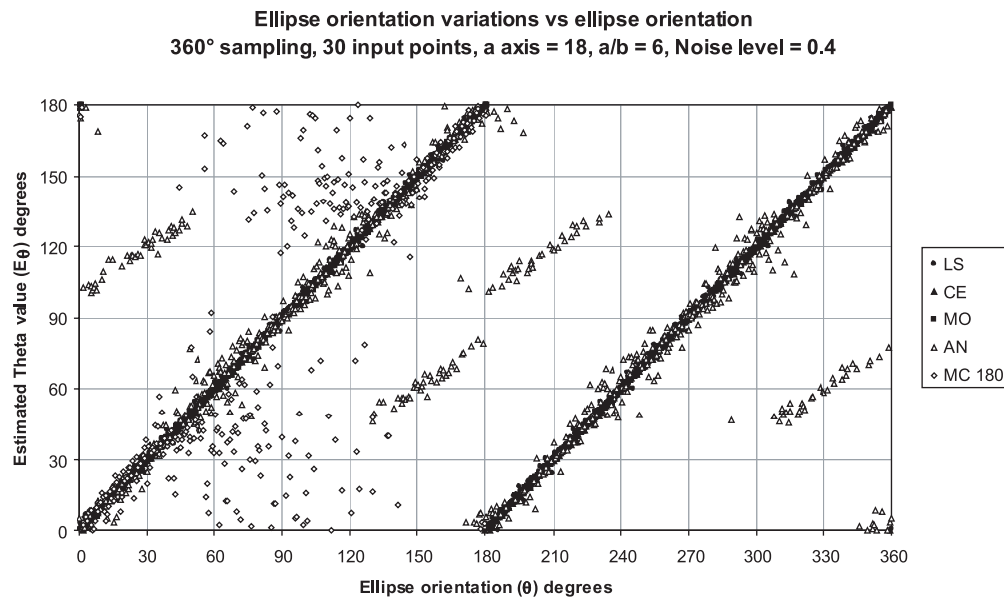


Fig. 15. $E\theta$ vs. θ , $alb=6$. See text for details.

common characteristics. For a given noise level, as the a/b values increase the Ea values become less accurate and tend toward underestimates, whereas the Eb and $E\theta$ values increase in accuracy up to a/b values of 4. This explanation proposed here is that it is due to the strong bias of ellipse radii magnitudes to the short semi axis magnitude with sample directions at constant angular spacings, as shown in Fig. 8. The results reported in this paper can be summarised in the context of specific boundary conditions, such that depending on the characteristics of a given situation (number of data, knowledge of noise levels in data, priority of semi axis magnitude vs. orientation in results, etc.), the best performing method can be identified.

If the dataset is small, for example 3–5 strain measurements from a single exposed surface, the least squares algorithm is the best choice for strain or curvature data. The analytical or Mohr circle methods could also be used but may be less accurate if the data is noisy or large axial ratios (≥ 6) are present. The presence of noise in the data may create many erroneous or non-numerical results in the analytical algorithm. If a visual estimate of the maximum strain axis can be utilised to select the sample directions then the analytical method will be more accurate. However, this defeats the object of using a method where a priori knowledge of the ellipse parameters is not required.

If a larger dataset is available ($n \geq 5$) for example nearest grain centres in a Fry analysis, or azimuth-dependant curvature from a gridded mapped surface then the mean object ellipse or least squares algorithms are the best choices, respectively. The mean object ellipse algorithm is also the best choice for strain data if an estimate of ellipse axial ratio is required. Where orientation is of importance then the mean object ellipse and least

squares algorithm are good choices although the constrained ellipse approach for strain data is also good. Overall, there is no situation where the analytical or Mohr circle approaches are the best performers although at a/b values less than 6 and noise levels less than 0.4 the Mohr circle method does not perform significantly worse than the other algorithms and it has the advantage of being useful for both strain and curvature datasets. Therefore, it may be suitable for strain or curvature analysis in low to moderately deformed areas. The Mohr circle method may actually be better suited to smaller datasets to avoid some of the problems with the pole construction with larger datasets.

These results should enable methodologies to be optimised in a range of geological applications, from strain analysis of small datasets to automated curvature analysis of millions of data points. Although we have attempted to review a broad range of existing methods, there are undoubtedly additional algorithms and concepts that might further refine approaches to the problem of ellipse fitting to geological data.

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