

ABSTRACT

We present a fast equivalent layer method for processing large-scale magnetic data. We demonstrate that the sensitivity matrix associated with an equivalent layer of dipoles ~~assumes~~can be arranged to a Block-Toeplitz Toeplitz-Block (BTTB) structure for the case in which observations and dipoles are aligned on a horizontal and regularly-spaced grid. The product of a BTTB matrix and an arbitrary vector represents a discrete convolution and can be efficiently computed via 2D Fast Fourier Transform. In this case, the matrix-vector product uses only the elements forming the first column of the BTTB matrix, saving computational time and memory. Our convolutional equivalent layer method uses this approach to compute the matrix-vector products in the iterative conjugate gradient algorithm with the purpose of estimating the physical-property distribution over the equivalent layer ~~with~~for large data sets. Synthetic tests with a mid-size 100×50 grid of total-field anomaly data show a decrease of $\approx 10^4$ in floating-point operations and $\approx 25\times$ in computation runtime of our method compared to the classical approach of solving the least-squares normal equations via Cholesky decomposition. ~~Better~~Faster results are obtained for millions of data, showing drastic decreases in RAM usage and runtime, allowing to perform magnetic data processing of large data sets on regular desktop computers. Our results also show that, compared to the classical Fourier approach, the magnetic data processing with our method requires similar computation time, but produces significantly smaller border effects without using any padding scheme and also is ~~much~~ more robust to deal with data on irregularly spaced points or on undulating observation surfaces. A test with 1,310,000 irregularly spaced field data over the Carajás Province, Brazil, confirms the ~~sueess~~efficiency of our method by taking ≈ 385.56 seconds to estimate the physical-property distribution over the equivalent layer and ≈ 2.64 seconds to compute the upward continuation.

INTRODUCTION

Large-scale data processing with tens of thousands of data ~~is~~ is a reality in all areas of geophysics including the geophysical potential fields. The potential-field data processing includes convolution integrals which can be solved either in the space or Fourier domains. The earliest techniques of potential-field data processing were developed in the space domain. For example, Peters (1949) accomplished, in the space domain, the second and fourth derivatives of magnetic data and the upward- and downward-continuations of magnetic data by deriving coefficients that are used in a graphical convolution with the magnetic data.

~~However~~However, the techniques for processing potential-field data in space-domain were soon substituted by the Fourier-domain techniques. Dean (1958) pointed out that the operations of second ~~derivative~~order derivatives, analytic continuation, smoothing, the removing of residuals or regionals, and others for processing potential-field data are similar to the electric filter circuits in Fourier domain. Dean (1958) was the first to develop the ~~theory of linear filter~~theory of linear filters in Fourier domain for gravity and magnetic processing and to present filters in Fourier domain (Dean, 1958, see Table I, p 113) for some theoretical geophysical operations (e.g., derivatives and upward and downward continuations). Gunn (1975) presented a comprehensive analysis of processing potential-field data in Fourier domain.

An approach for processing potential-field data in space domain is the equivalent-layer technique. The equations deductions of the equivalent layer as a solution of the Laplace's equation in the region above the source was first presented by Kellogg (1929) and detailed explanations can also be found in Blakely (1996). Although the equivalent-layer technique has been known since the 1960s in geophysical literature (Danes, 1961; Bott, 1967; Dampney,

1969), its use has become feasible only recently because of the advances in computational power. In magnetic data procesing, some authors explored this technique for calculating the first and second vertical derivatives of the fields (Emilia, 1973), reduction to the pole (Silva, 1986; Oliveira Jr. et al., 2013; Li et al., 2014), upward/downward continuations (Hansen and Miyazaki, 1984; Li and Oldenburg, 2010) and total magnetic induction vector components calculation (Sun et al., 2019).

Together with the rise in computational processing power, some works tried new implementations to increase the efficiency of the equivalent layer. In Leão and Silva (1989) the authors used a shifting window over the layer, -increasing to increase the number of linear systems to be solved, but to reduced the size of each linear system is reduced at the same time. Another approach for a fast equivalent layer was proposed by Li and Oldenburg (2010) who transformed the full sensitivity matrix into a sparse one by using the compression of the coefficient matrix using fast wavelet transforms based on orthonormal, compactly supported wavelets. Oliveira Jr. et al. (2013) divided the equivalent layer into a grid of fixed source windows. Instead of directly calculating the the physical-property distribution of a finite set of equivalent sources (e.g., dipoles, point of masses) arranged in the entire equivalent layer, Oliveira Jr. et al. (2013) estimated the coefficients of a bivariate polynomial function describing the physical-property distribution within each equivalent-source window. The estimated polynomial-coefficients are transformed into the physical-property distribution and thus any standard linear transformation of the data can be performed. Grounded on excess mass constraint, Siqueira et al. (2017) proposed an iterative method for processing large gravimetric data using the equivalent layer without requiring the solution of a linear system. In Siqueira et al. (2017), the initial mass distribution over the equivalent layer is proportional to observed gravity data and it is updated at each iteration by adding mass

corrections that are proportional to the residuals of observed and estimated data.

One of the greatest obstacles to the use of the equivalent-layer technique for processing potential-field data is the solution of the associated linear system. A wide variety of applications in mathematics and engineering that fall into Toeplitz systems propelled the development of a large variety of methods for solving them. Direct methods were conceived by Levinson (1946) and by Trench (1964). Currently, the iterative method of conjugate gradient is used in most cases, in Chan and Jin (2007) the authors presented an introduction on the topic for 1D data structures of Toeplitz matrices and also for 2D data structures, which they called block-Toeplitz Toeplitz-block matrices. In both cases, the solving strategy is to ~~embed~~ embed the Toeplitz/BTTB matrix into a Circulant/Block-Circulant Circulant-Block matrix, calculate its eigenvalues by a 1D or 2D fast Fourier transform of its first column, respectively and carry the matrix-vector product between kernel and parameters at each iteration of the conjugate gradient method in a very fast manner.

In potential field methods, the properties of Toeplitz system have been used for downward continuation (Zhang et al., 2016) and for 3D gravity-data inversion using a 2D multilayer model (Zhang and Wong, 2015). More recently, Hogue et al. (2020) provided an overview on modeling the gravity and magnetic kernels using the BTTB structures and Renaud et al. (2020) used BTTB the structures for inversion of both gravity and magnetic data to recover sparse subsurface structures. Takahashi et al. (2020) combined the fast equivalent source technique presented by Siqueira et al. (2017) with the concept of symmetric block-Toeplitz Toeplitz-block (BTTB) matrices to introduce the convolutional equivalent layer for gravimetric data technique. Takahashi et al. (2020) showed that the BTTB structure appears when the sensitivity matrix of the linear system, required to solve the gravimetric equivalent layer, is calculated on a regular spaced grid of dataset with constant height and

each equivalent source is exactly beneath each observed data point. This work showed an decrease in the order of 10^4 in floating-point operations needed to estimate the equivalent sources; thus, the Takahashi et al. (2020) method was able to efficiently process very large gravity data sets. Moreover, Takahashi et al. (2020) method yielded neither significant boundary effects nor noise amplification.

In this work, the convolutional equivalent layer using the block-Toeplitz Toeplitz-block idea, presented in Takahashi et al. (2020), will be used to solve the linear system required to estimate the physical property that produces a magnetic field on regular grids. Here, we ~~achive~~achieve very fast solutions using a conjugate gradient algorithm combined with the fast Fourier transform. We present a novel method of exploring the symmetric structures of the second derivatives of the inverse of the distance contained in the magnetic kernel, to keep the memory RAM usage to the minimal by using only one equivalent source to carry the calculations of the forward problem. We also show tests of the magnetic convolutional equivalent layer when irregular grids are used. The convergence of the conjugate gradient maintains in an acceptable level even using irregular grids. Our results show the good performance of our method in producing fast and robust solutions for processing large amounts of magnetic data using the equivalent layer technique.

METHODOLOGY

Classical equivalent layer for magnetic data

Let \mathbf{d}^o be the $N \times 1$ observed data vector $(x_i, y_i, z_i), i = 1, \dots, N$, whose i th element is the total-field anomaly d_i^o produced by arbitrarily magnetized sources ~~at the position~~ (x_i, y_i, z_i) , $i = 1, \dots, N$, ~~of arranged in~~ a right-handed Cartesian coordinate system with x -, y - and z -

axis pointing to north, east and down, respectively. We consider that the total-field anomaly data d_i^o represent the discrete values of a harmonic function. Besides, we consider that the main geomagnetic field direction ~~at the study area is constant at the limited study area and~~ can be defined by the unit vector

$$\hat{\mathbf{F}} = \begin{bmatrix} F_x \\ F_y \\ F_z \end{bmatrix} = \begin{bmatrix} \cos(I_0) \cos(D_0) \\ \cos(I_0) \sin(D_0) \\ \sin(I_0) \end{bmatrix}, \quad (1)$$

with constant inclination I_0 and declination D_0 . In this case, d_i^o can be approximated by the predicted total-field anomaly

$$\Delta T_i = \sum_{j=1}^M p_j a_{ij}, \quad (2)$$

which describes the magnetic induction exerted ~~at~~ at the observation point ~~(x_i, y_i, z_i)~~, by a discrete layer of M dipoles (equivalent sources) defined on the horizontal plane $z = z_c$, where p_j is the magnetic moment intensity (in A m^2) of the j th dipole, that has unit volume and is located at the point (x_j, y_j, z_c) . In equation 2, a_{ij} is the harmonic function

$$a_{ij} = c_m \frac{\mu_0}{4\pi} \hat{\mathbf{F}}^\top \mathbf{H}_{ij} \hat{\mathbf{u}}, \quad (3)$$

the unit vector

$$\hat{\mathbf{u}} = \begin{bmatrix} u_x \\ u_y \\ u_z \end{bmatrix} = \begin{bmatrix} \cos(I) \cos(D) \\ \cos(I) \sin(D) \\ \sin(I) \end{bmatrix}, \quad (4)$$

defines the magnetization direction of all dipoles, with constant inclination I and declination D , $\mu_0 = 4\pi 10^{-7} \text{ H/m}$ is the magnetic constant, $c_m = 10^9$ is a factor that transforms the

magnetic induction from Tesla (T) to nanotesla (nT) and \mathbf{H}_{ij} is a 3×3 matrix

$$\mathbf{H}_{ij} = \begin{bmatrix} h_{ij}^{xx} & h_{ij}^{xy} & h_{ij}^{xz} \\ h_{ij}^{xy} & h_{ij}^{yy} & h_{ij}^{yz} \\ h_{ij}^{xz} & h_{ij}^{yz} & h_{ij}^{zz} \end{bmatrix}, \quad (5)$$

where

$$h_{ij}^{\alpha\beta} = \begin{cases} \frac{3(\alpha_i - \alpha_j)^2}{r_{ij}^5} - \frac{1}{r_{ij}^3}, & \alpha = \beta \\ \frac{3(\alpha_i - \alpha_j)(\beta_i - \beta_j)}{r_{ij}^5}, & \alpha \neq \beta \end{cases}, \quad \alpha, \beta = x, y, z, \quad (6)$$

are the second derivatives of the inverse distance function

$$\frac{1}{r_{ij}} = \frac{1}{\sqrt{(x_i - x_j)^2 + (y_i - y_j)^2 + (z_i - z_j)^2}} \quad (7)$$

with respect to the coordinates of the observation point (x_i, y_i, z_i) .

Equation 2 can be rewritten in matrix notation as follows:

$$\mathbf{d}(\mathbf{p}) = \mathbf{A}\mathbf{p}, \quad (8)$$

where $\mathbf{d}(\mathbf{p})$ is the $N \times 1$ predicted data vector with i th element defined as the predicted total-field anomaly ΔT_i (equation 2), \mathbf{p} is the $M \times 1$ parameter vector whose j th element is the magnetic moment intensity p_j of the j th dipole and \mathbf{A} is the $N \times M$ sensitivity matrix with element ij defined by the harmonic function a_{ij} (equation 3). In the classical equivalent-layer technique, the common approach for estimating the parameter vector \mathbf{p} from the observed total-field anomaly data \mathbf{d}^o is solving the least-squares normal equations

$$\mathbf{A}^\top \mathbf{A} \mathbf{p} = \mathbf{A}^\top \mathbf{d}^o. \quad (9)$$

Equation 9 is usually solved by first computing the Cholesky factor \mathbf{G} of [the positive-definite](#) matrix $\mathbf{A}^\top \mathbf{A}$ and then using it to solve the linear systems (Golub and Loan, 2013, p. 262):

$$\begin{aligned} \mathbf{G}\mathbf{w} &= \mathbf{A}^\top \mathbf{d}^o \\ \mathbf{G}^\top \tilde{\mathbf{p}} &= \mathbf{w} \end{aligned}, \quad (10)$$

where \mathbf{w} is a ~~dummy~~^{temporarily} variable. This approach to ~~estimate~~^{estimating} the parameter vector will be referenced throughout this work as the *classical method*. The computational cost associated with the classical method can be very high when dealing with large datasets. In the following subsections, we will show how to explore the structure of the sensitivity matrix \mathbf{A} and efficiently solve the least-squares normal equations (equation 9).

Matrix \mathbf{A} in terms of matrix components $\mathbf{A}_{\alpha\beta}$

To access the structure of the sensitivity matrix \mathbf{A} (equation 8), let us first rewrite its elements a_{ij} (equation 3) in the following way:

$$a_{ij} = a_{ij}^{xx} + a_{ij}^{xy} + a_{ij}^{xz} + a_{ij}^{yy} + a_{ij}^{yz} + a_{ij}^{zz}, \quad (11)$$

where

$$a_{ij}^{\alpha\beta} = \begin{cases} c_m \frac{\mu_0}{4\pi} (F_\alpha u_\beta) h_{ij}^{\alpha\beta} & , \quad \alpha = \beta \\ c_m \frac{\mu_0}{4\pi} (F_\alpha u_\beta + F_\beta u_\alpha) h_{ij}^{\alpha\beta} & , \quad \alpha, \beta = x, y, z, \\ & , \quad \alpha \neq \beta \end{cases} \quad (12)$$

are defined by the elements of $\hat{\mathbf{F}}$ (equation 1), $\hat{\mathbf{u}}$ (equation 4) and \mathbf{H}_{ij} (equations 5 and 6).

Then, we can rewrite the sensitivity matrix \mathbf{A} (equation 8) according to:

$$\mathbf{A} = \mathbf{A}_{xx} + \mathbf{A}_{xy} + \mathbf{A}_{xz} + \mathbf{A}_{yy} + \mathbf{A}_{yz} + \mathbf{A}_{zz}, \quad (13)$$

where $\mathbf{A}_{\alpha\beta}$ are $N \times M$ matrices with elements ij defined by $a_{ij}^{\alpha\beta}$ (equation 12).

Now we can define the structure of \mathbf{A} in terms of its components $\mathbf{A}_{\alpha\beta}$ (equation 13). To do this, we consider the particular case in which the observed total-field anomaly is located on an $N_x \times N_y$ regular grid of points spaced by Δ_x and Δ_y along the x - and y -directions, respectively, on a constant vertical coordinate z_0 . We also consider that the equivalent layer

is ~~formed by one dipole~~ arranged such that one dipole is located right below each observation point, at ~~the~~a constant coordinate z_c . In this case, the number of equivalent sources M is equal to the number of data N and, consequently, matrices \mathbf{A} and $\mathbf{A}_{\alpha\beta}$ become square ($N \times N$). Besides, the horizontal coordinates x_i and y_i of the observation points can be defined by

$$x_i = x_1 + [k(i) - 1] \Delta_x \quad (14)$$

and

$$y_i = y_1 + [l(i) - 1] \Delta_y, \quad (15)$$

where x_1 and y_1 are the ~~lower limits start grid coordinates~~ for x_i and y_i , respectively, and $k(i)$ and $l(i)$ are integer functions defined according to the orientation of the data grid (Figure 1). For *x-oriented grids*, the integer functions are given by

$$k(i) = i - \left\lceil \frac{i}{N_x} \right\rceil N_x + N_x \quad (16)$$

and

$$l(i) = \left\lceil \frac{i}{N_x} \right\rceil. \quad (17)$$

For *y-oriented grids*, the integer functions are given by

$$k(i) = \left\lceil \frac{i}{N_y} \right\rceil \quad (18)$$

and

$$l(i) = i - \left\lceil \frac{i}{N_y} \right\rceil N_y + N_y. \quad (19)$$

In equations 16–19, $\lceil \cdot \rceil$ denotes the ceiling function (e.g., Graham et al., 1994, p. 67-68). Equations 14–19 can also be used to define the coordinates x_j and y_j of the equivalent sources, but with index j instead of i .

By using equations 14–19 to define the coordinates x_i and y_i of the observation points and x_j and y_j of the equivalent sources, we can rewrite the elements $h_{ij}^{\alpha\beta}$ (equation 6) of matrix \mathbf{H}_{ij} (equation 5) as follows:

$$h_{ij}^{xx} = \frac{3(\Delta k_{ij} \Delta_x)^2}{r_{ij}^5} - \frac{1}{r_{ij}^3}, \quad (20)$$

$$h_{ij}^{yy} = \frac{3(\Delta l_{ij} \Delta_y)^2}{r_{ij}^5} - \frac{1}{r_{ij}^3}, \quad (21)$$

$$h_{ij}^{zz} = \frac{3\Delta_z^2}{r_{ij}^5} - \frac{1}{r_{ij}^3}, \quad (22)$$

$$h_{ij}^{xy} = \frac{3(\Delta k_{ij} \Delta_x)(\Delta l_{ij} \Delta_y)}{r_{ij}^5}, \quad (23)$$

$$h_{ij}^{xz} = \frac{3(\Delta k_{ij} \Delta_x) \Delta_z}{r_{ij}^5} \quad (24)$$

and

$$h_{ij}^{yz} = \frac{3(\Delta l_{ij} \Delta_y) \Delta_z}{r_{ij}^5}, \quad (25)$$

where $\Delta_z = z_c - z_0$,

$$\Delta k_{ij} = \frac{x_i - x_j}{\Delta_x} = k(i) - k(j), \quad (26)$$

$$\Delta l_{ij} = \frac{y_i - y_j}{\Delta_y} = l(i) - l(j) \quad (27)$$

and

$$\frac{1}{r_{ij}} = \frac{1}{\sqrt{(\Delta k_{ij} \Delta_x)^2 + (\Delta l_{ij} \Delta_y)^2 + \Delta_z^2}}. \quad (28)$$

Note that the integer functions $k(i)$, $k(j)$, $l(i)$ and $l(j)$ (equations 16–19) defining Δk_{ij} (equation 26), Δl_{ij} (equation 27) and $\frac{1}{r_{ij}}$ (equation 28) assume different forms depending on the defined grid orientation. Despite of that, it can be shown that

$$\Delta k_{ij} = -\Delta k_{ji}, \quad (29)$$

$$\Delta l_{ij} = -\Delta l_{ji} \quad (30)$$

and

$$\frac{1}{r_{ij}} = \frac{1}{r_{ji}} \quad (31)$$

for any grid orientation.

General structure of matrices $\mathbf{A}_{\alpha\beta}$

By using equations 20–28 to compute $a_{ij}^{\alpha\beta}$ (equation 12), we can show that matrices $\mathbf{A}_{\alpha\beta}$ (equation 13) assume well-defined structures that can be conveniently represented with *block indices* q and p (Takahashi et al., 2020). These indices are defined by the integer functions Δk_{ij} and Δl_{ij} (equations 26 and 27) in terms of the indices i of the observation points (x_i, y_i, z_0) and j of the equivalent sources (x_j, y_j, z_c) . For *x-oriented grids* (Figure 1), $Q = N_y$, $P = N_x$ and the block indices q and p are given by:

$$q \equiv q(i, j) = \Delta l_{ij} \quad (32)$$

and

$$p \equiv p(i, j) = \Delta k_{ij}, \quad (33)$$

where Δk_{ij} and Δl_{ij} (equations 26 and 27) are defined by integer functions $k(i)$, $k(j)$, $l(i)$ and $l(j)$ given by equations 16 and 17. For *y-oriented grids* (Figure 1), $Q = N_x$, $P = N_y$ and the block indices q and p are given by:

$$q \equiv q(i, j) = \Delta k_{ij} \quad (34)$$

and

$$p \equiv p(i, j) = \Delta l_{ij}, \quad (35)$$

where Δk_{ij} and Δl_{ij} (equations 26 and 27) are defined by integer functions $k(i)$, $k(j)$, $l(i)$ and $l(j)$ given by equations 18 and 19. Equations 32–35 show that q varies from $-Q + 1$ to

$Q - 1$ and p from $-P + 1$ to $P - 1$, regardless of the grid orientation. They Notice these equations differ from those presented by Takahashi et al. (2020) due to the absence of the module by not being represented as the absolute values.

Let us consider the small regular grid of $N_x = 3$ and $N_y = 2$ points shown by Figure 1. This grid may represent observation points (x_i, y_i, z_0) with constant vertical coordinate z_0 or equivalent sources (x_j, y_j, z_c) with constant vertical coordinate $z_c > z_0$. In both cases, the horizontal coordinates are defined by equations 14 and 15. Given an index i , associated with an observation point, and an index j , associated with an equivalent source, we can compute Δk_{ij} (equation 26), Δl_{ij} (equation 27) and $\frac{1}{r_{ij}}$ (equation 28). The matrices $\Delta \mathbf{K}$ and $\Delta \mathbf{L}$ having elements ij defined by Δk_{ij} and Δl_{ij} , respectively, assume different forms, depending on the grid orientation. For x -oriented grids (Figure 1), they are given by:

$$\Delta \mathbf{K} = \begin{bmatrix} 0 & -1 & -2 & 0 & -1 & -2 \\ 1 & 0 & -1 & 1 & 0 & -1 \\ 2 & 1 & 0 & 2 & 1 & 0 \\ 0 & -1 & -2 & 0 & -1 & -2 \\ 1 & 0 & -1 & 1 & 0 & -1 \\ 2 & 1 & 0 & 2 & 1 & 0 \end{bmatrix} \quad (36)$$

and

$$\Delta \mathbf{L} = \begin{bmatrix} 0 & 0 & 0 & -1 & -1 & -1 \\ 0 & 0 & 0 & -1 & -1 & -1 \\ 0 & 0 & 0 & -1 & -1 & -1 \\ 1 & 1 & 1 & 0 & 0 & 0 \\ 1 & 1 & 1 & 0 & 0 & 0 \\ 1 & 1 & 1 & 0 & 0 & 0 \end{bmatrix}. \quad (37)$$

For y -oriented grids (Figure 1), they are given by:

$$\Delta\mathbf{K} = \begin{bmatrix} 0 & 0 & -1 & -1 & -2 & -2 \\ 0 & 0 & -1 & -1 & -2 & -2 \\ 1 & 1 & 0 & 0 & -1 & -1 \\ 1 & 1 & 0 & 0 & -1 & -1 \\ 2 & 2 & 1 & 1 & 0 & 0 \\ 2 & 2 & 1 & 1 & 0 & 0 \end{bmatrix} \quad (38)$$

and

$$\Delta\mathbf{L} = \begin{bmatrix} 0 & -1 & 0 & -1 & 0 & -1 \\ 1 & 0 & 1 & 0 & 1 & 0 \\ 0 & -1 & 0 & -1 & 0 & -1 \\ 1 & 0 & 1 & 0 & 1 & 0 \\ 0 & -1 & 0 & -1 & 0 & -1 \\ 1 & 0 & 1 & 0 & 1 & 0 \end{bmatrix}. \quad (39)$$

For example, consider the matrix coordinates x_5, y_5 for the observation point and x_3, y_3 for the source in a x -oriented grid, these are defined by the matrix index $i = 5$ and $j = 3$. Using equations 16 and 17 the grid indices $k(i) = 2$, $l(i) = 2$ and $k(j) = 3$, $l(j) = 1$ can be calculated. Equations 26 and 27 define $\Delta k_{ij} = -1$ and $\Delta l_{ij} = 1$, which represents the value in the fifth row and third column of matrices $\Delta\mathbf{K}$ and $\Delta\mathbf{L}$ (equations 36 and 37), respectively. If we use Using the matrix coordinates x_4 (y_4) for the observation point and x_2 (y_2) for the source, lead us to the same values for Δk_{ij} and Δl_{ij} (fourth row and second column of equations 36 and 37). These examples (equations 36–39) show that different combinations of indices i and j result in integer functions Δk_{ij} and Δl_{ij} (equations 26 and 27) having the same numerical value. In these cases, not only the numerical values of the corresponding elements $a_{ij}^{\alpha\beta}$ (equation 12), but also their associated block indices q and

p (equations 32–35) are the same. The contrary is also true: if the elements $a_{ij}^{\alpha\beta}$ having have different associated block indices q and p , they also have different numerical values. Because of that, using the alternative notation $a_{qp}^{\alpha\beta}$ to define the elements $a_{ij}^{\alpha\beta}$ in terms of its associated block indices q and p is a good approach to investigate-investigating the structure of a given matrix component $\mathbf{A}_{\alpha\beta}$ (equation 13). This approach allows identifying elements $a_{ij}^{\alpha\beta}$ having the same numerical value only by inspecting their associated block indices.

Note that for x -oriented grids, matrices $\Delta\mathbf{K}$ (equation 36) and $\Delta\mathbf{L}$ (equation 37) define the block indices p (equation 33) and q (equation 32), respectively. In this case, they are composed of $Q \times Q$ blocks with $P \times P$ elements each, where $Q = N_y$ and $P = N_x$. For y -oriented grids, matrices $\Delta\mathbf{K}$ (equation 38) and $\Delta\mathbf{L}$ (equation 39) define the block indices q (equation 34) and p (equation 35), respectively. In this case, they are also composed of $Q \times Q$ blocks with $P \times P$ elements each, but now $Q = N_x$ and $P = N_y$. The examples shown by equations 36–39 also illustrate that, regardless of grid orientation, (i) the block index q is constant inside each block; (ii) blocks disposed along the same block diagonal are equal to each other; (iii) the block index p is constant on each diagonal of a given block; (iv) elements of a given block located on the same diagonal are also equal do each other. The results obtained with the small grid shown in Figure 1 can be easily generalized for larger grids. Based on the well-defined structure of block indices, we can define matrices $\mathbf{A}_{\alpha\beta}$ in a general form

$$\mathbf{A}_{\alpha\beta} = \begin{bmatrix} \mathbf{A}_{\alpha\beta}^0 & \mathbf{A}_{\alpha\beta}^{-1} & \cdots & \mathbf{A}_{\alpha\beta}^{-Q+1} \\ \mathbf{A}_{\alpha\beta}^1 & \ddots & \ddots & \vdots \\ \vdots & \ddots & \ddots & \mathbf{A}_{\alpha\beta}^{-1} \\ \mathbf{A}_{\alpha\beta}^{Q-1} & \cdots & \mathbf{A}_{\alpha\beta}^1 & \mathbf{A}_{\alpha\beta}^0 \end{bmatrix}_{N \times N}, \quad (40)$$

with blocks $\mathbf{A}_{\alpha\beta}^q$, $q = -Q + 1, \dots, Q - 1$, given by

$$\mathbf{A}_{\alpha\beta}^q = \begin{bmatrix} a_{q0}^{\alpha\beta} & a_{q(-1)}^{\alpha\beta} & \cdots & a_{q(-P+1)}^{\alpha\beta} \\ a_{q1}^{\alpha\beta} & \ddots & \ddots & \vdots \\ \vdots & \ddots & \ddots & a_{q(-1)}^{\alpha\beta} \\ a_{q(P-1)}^{\alpha\beta} & \cdots & a_{q1}^{\alpha\beta} & a_{q0}^{\alpha\beta} \end{bmatrix}_{P \times P}, \quad (41)$$

formed by elements $a_{qp}^{\alpha\beta}$, $p = -P + 1, \dots, P - 1$. This well-defined structure (equations 40 and 41) of matrix components $\mathbf{A}_{\alpha\beta}$ (equation 13) is called Block-Toeplitz Toeplitz-Block (BTTB) (e.g., Chan and Jin, 2007, p. 67).

Detailed structure of matrices \mathbf{A}_{xx} , \mathbf{A}_{yy} and \mathbf{A}_{zz}

Equations 40 and 41 define the general BTTB structure of all matrix components $\mathbf{A}_{\alpha\beta}$, but there are some differences between them. Let us consider the matrix component \mathbf{A}_{xx} , with elements a_{ij}^{xx} (equation 12) defined by the second derivative calculated in h_{ij}^{xx} (equation 20). It can be easily verified from equations 29 and 31 that $h_{ij}^{xx} = h_{ji}^{xx}$. As a consequence, $a_{ij}^{xx} = a_{ji}^{xx}$, which means that

$$\mathbf{A}_{xx} = (\mathbf{A}_{xx})^\top \quad (42)$$

for any grid orientation. Now, let us investigate the elements a_{qp}^{xx} forming the blocks \mathbf{A}_{xx}^q . For x -oriented grids (Figure 1), the block indices q and p are defined by equations 32 and 33 and a_{qp}^{xx} can be rewritten as follows:

$$a_{qp}^{xx} = c_m \frac{\mu_0}{4\pi} (F_x u_x) \frac{3(p\Delta_x)^2}{r_{qp}^5} - \frac{1}{r_{qp}^3}, \quad (43)$$

where

$$\frac{1}{r_{qp}} = \frac{1}{\sqrt{(p\Delta_x)^2 + (q\Delta_y)^2 + \Delta_z^2}}. \quad (44)$$

For y -oriented grids (Figure 1), the block indices q and p are defined by equations 34 and 35 and a_{qp}^{xx} can be rewritten as follows:

$$a_{qp}^{xx} = c_m \frac{\mu_0}{4\pi} (F_x u_x) \frac{3(q \Delta_x)^2}{r_{qp}^5} - \frac{1}{r_{qp}^3}, \quad (45)$$

where

$$\frac{1}{r_{qp}} = \frac{1}{\sqrt{(q \Delta_x)^2 + (p \Delta_y)^2 + \Delta_z^2}}. \quad (46)$$

From equations 43–46, we can easily verify that

$$\mathbf{A}_{xx}^q = \mathbf{A}_{xx}^{(-q)} \quad (47)$$

and

$$\mathbf{A}_{xx}^q = (\mathbf{A}_{xx}^q)^\top. \quad (48)$$

Note that these symmetries are valid for any grid orientation. From this results we conclude the matrix component \mathbf{A}_{xx} is *symmetric-Block-Toeplitz symmetric-Toeplitz-Block* for any grid orientation. The same reasoning can be used to show that matrices \mathbf{A}_{yy} and \mathbf{A}_{zz} also have this symmetric structure.

Detailed structure of matrix \mathbf{A}_{xy}

Let \mathbf{A}_{xy} be a matrix component with elements a_{ij}^{xy} (equation 12) defined by the second derivative calculated in h_{ij}^{xy} (equation 23). It can be easily verified from equations 29–31 that $h_{ij}^{xy} = h_{ji}^{xy}$. As a consequence, $a_{ij}^{xy} = a_{ji}^{xy}$, which means that

$$\mathbf{A}_{xy} = (\mathbf{A}_{xy})^\top \quad (49)$$

for any grid orientation. For x -oriented grids (Figure 1), the block indices q and p are defined by equations 32 and 33 and a_{qp}^{xy} can be rewritten as follows:

$$a_{qp}^{xy} = c_m \frac{\mu_0}{4\pi} (F_x u_y + F_y u_x) \frac{3(p \Delta_x)(q \Delta_y)}{r_{qp}^5}, \quad (50)$$

with $\frac{1}{r_{qp}}$ defined by equation 44. For y -oriented grids (Figure 1), the block indices q and p are defined by equations 34 and 35 and a_{qp}^{xy} can be rewritten as follows:

$$a_{qp}^{xy} = c_m \frac{\mu_0}{4\pi} (F_x u_y + F_y u_x) \frac{3(q\Delta_x)(p\Delta_y)}{r_{qp}^5}, \quad (51)$$

with $\frac{1}{r_{qp}}$ defined by equation 46. From equations 44, 46, 50 and 51, we can show that

$$\mathbf{A}_{xy}^q = -\mathbf{A}_{xy}^{(-q)} \quad (52)$$

and

$$\mathbf{A}_{xy}^q = -(\mathbf{A}_{xy}^q)^\top. \quad (53)$$

Note that these symmetries are valid for any grid orientation. From this results we conclude the matrix component \mathbf{A}_{xy} is *skew symmetric-Block-Toeplitz skew symmetric-Toeplitz-Block* for any grid orientation.

Detailed structure of matrices \mathbf{A}_{xz} and \mathbf{A}_{yz}

Let \mathbf{A}_{xz} be a matrix component with elements a_{ij}^{xz} (equation 12) defined by the second derivative calculated in h_{ij}^{xz} (equation 24). It can be easily verified from equations 29–31 that $h_{ij}^{xz} = -h_{ji}^{xz}$. As a consequence, $a_{ij}^{xz} = -a_{ji}^{xz}$, which means that

$$\mathbf{A}_{xz} = -(\mathbf{A}_{xz})^\top \quad (54)$$

for any grid orientation. For x -oriented grids (Figure 1), the block indices q and p are defined by equations 32 and 33 and a_{qp}^{xz} can be rewritten as follows:

$$a_{qp}^{xz} = c_m \frac{\mu_0}{4\pi} (F_x u_z + F_z u_x) \frac{3(p\Delta_x)\Delta_z}{r_{qp}^5}, \quad (55)$$

with $\frac{1}{r_{qp}}$ defined by equation 44. In this case, we can see that

$$\mathbf{A}_{xz}^q = \mathbf{A}_{xz}^{(-q)} \quad (56)$$

and

$$\mathbf{A}_{xz}^q = -(\mathbf{A}_{xz}^q)^\top . \quad (57)$$

This structure is called *symmetric-Block-Toeplitz skew symmetric-Toeplitz-Block* and is valid only for x -oriented grids. For y -oriented grids (Figure 1), the block indices q and p are defined by equations 34 and 35 and a_{qp}^{xz} can be rewritten as follows:

$$a_{qp}^{xz} = c_m \frac{\mu_0}{4\pi} (F_x u_z + F_z u_x) \frac{3(q\Delta_x)\Delta_z}{r_{qp}^5} , \quad (58)$$

with $\frac{1}{r_{qp}}$ defined by equation 46. Now, we conclude that

$$\mathbf{A}_{xz}^q = -\mathbf{A}_{xz}^{(-q)} \quad (59)$$

and

$$\mathbf{A}_{xz}^q = (\mathbf{A}_{xz}^q)^\top . \quad (60)$$

This structure is called *skew symmetric-Block-Toeplitz symmetric-Toeplitz-Block* and is valid only for y -oriented grids.

The same reasoning can be followed to show that

$$\mathbf{A}_{yz} = -(\mathbf{A}_{yz})^\top \quad (61)$$

for any grid orientation. Besides, we can also show that

$$\mathbf{A}_{yz}^q = -\mathbf{A}_{yz}^{(-q)} \quad (62)$$

and

$$\mathbf{A}_{yz}^q = (\mathbf{A}_{yz}^q)^\top \quad (63)$$

for x -oriented grids (*skew symmetric-Block-Toeplitz symmetric-Toeplitz-Block*), while

$$\mathbf{A}_{yz}^q = \mathbf{A}_{yz}^{(-q)} \quad (64)$$

and

$$\mathbf{A}_{yz}^q = -(\mathbf{A}_{yz}^q)^\top \quad (65)$$

for y -oriented grids (*symmetric-Block-Toeplitz skew symmetric-Toeplitz-Block*).

Convolutional equivalent layer

The computational cost associated with the classical method to estimate the parameter vector \mathbf{p} by solving the linear system 9 can be very high or even prohibitive when dealing with large data sets. In these cases, a well-known alternative is solving the normal equations (equation 9) iteratively by using the *standard Conjugate Gradient Least Squares (CGLS) method*:

Algorithm 1 Standard CGLS pseudocode (Aster et al., 2019, p. 166).

Input: \mathbf{A} and \mathbf{d}^o .

Output: Estimated parameter vector $\tilde{\mathbf{p}}$.

Set $it = 0$, $\tilde{\mathbf{p}}_{(it)} = \mathbf{0}$, $\mathbf{c}_{(it-1)} = \mathbf{0}$, $\beta_{(it)} = 0$, $\mathbf{s}_{(it)} = \mathbf{d}^o$ and $\mathbf{r}_{(it)} = \mathbf{A}^\top \mathbf{s}_{(it)}$.

$$1 - \text{If } it > 0, \beta_{(it)} = \frac{\|\mathbf{r}_{(it)}\|_2^2}{\|\mathbf{r}_{(it-1)}\|_2^2}$$

$$2 - \mathbf{c}_{(it)} = \mathbf{r}_{(it)} + \beta_{(it)} \mathbf{c}_{(it-1)}$$

$$3 - \alpha_{(it)} = \frac{\|\mathbf{r}_{(it)}\|_2^2}{\|\mathbf{A} \mathbf{c}_{(it)}\|_2^2}$$

$$4 - \tilde{\mathbf{p}}_{(it+1)} = \tilde{\mathbf{p}}_{(it)} + \alpha_{(it)} \mathbf{c}_{(it)}$$

$$5 - \mathbf{s}_{(it+1)} = \mathbf{s}_{(it)} - \alpha_{(it)} \mathbf{A} \mathbf{c}_{(it)}$$

$$6 - \mathbf{r}_{(it+1)} = \mathbf{A}^\top \mathbf{s}_{(it+1)}$$

$$7 - it = it + 1$$

$$8 - \text{Repeat previous steps until convergence } \underline{(\delta)}.$$

Setting a convergence **criteria** criterion δ (Algorithm 1) based on the minimum tolerance

of the residuals is a good option to carry out this algorithm efficiently and still obtaining very good results. Another possibility is to set an invariance to the Euclidean norm of residuals between iterations, which would increase algorithm runtime, but with smaller residuals. We chose the ~~first latter~~ option, as we ~~achieve satisfactory~~ ~~could achieve better~~ results.

Note that the standard CGLS solution (Algorithm 1) requires neither inverse matrix nor matrix-matrix product. Instead, it only requires: one matrix-vector product out of the loop and two matrix-vector products per iteration (in steps 3 and 6). These products can be efficiently computed by using the 2D FFT, as a discrete convolution (see Appendix A). Takahashi et al. (2020) used this approach to develop an efficient algorithm for gravity data processing. This modified approach in which the standard CGLS method is modified to efficiently compute the matrix-vector products will be referenced throughout this work as the *convolutional equivalent layer method*.

Computational performance

In this ~~seetions~~ section we compare the efficiency of the classical (equation 10), standard CGLS (Algorithm 1) and the convolutional equivalent layer method (Algorithm 1 with matrix-vector products computed according to Appendix A). To do this, we compute the total number of *flops* associated to them (Golub and Loan, 2013, p. 12).

For the classical method, we have $\frac{1}{2}N^3$ flops to compute the lower triangle of $\mathbf{A}^\top \mathbf{A}$; $\frac{1}{3}N^3$ flops to compute the Cholesky factor \mathbf{G} of $\mathbf{A}^\top \mathbf{A}$ (Golub and Loan, 2013, p. 164); $2N^2$ flops to compute the matrix-vector product $\mathbf{A}^\top \mathbf{d}^o$; and $2N^2$ flops to solve the triangular systems given by equation 10 (Golub and Loan, 2013, p. 106). The resultant flop count for

the classical method is

$$f_{classical} = \frac{5}{6}N^3 + 4N^2 . \quad (66)$$

For the standard CGLS method (Algorithm 1) we have $2N^2$ to compute the matrix-vector product $\mathbf{A}^\top \mathbf{s}_{(it)}$ out of the loop; $4N$ in step 1; $2N$ in step 2; $2N^2 + 2N$ in step 3; $2N$ in step 4; $2N$ in step 5; and $2N^2$ in step 6. The resultant flop count is given by:

$$f_{cglsls} = 2N^2 + it(4N^2 + 12N) . \quad (67)$$

To compute the flops count of our method, we need only to replace the flops associated with matrix-vector products in the standard CGLS method by those associated with 2D convolution defined in Appendix A, which consists of $\kappa 4N \log_2(4N)$ flops to compute the 2D FFT for each matrix $\mathbf{L}_{\alpha\beta}$ (equation A-14); $\kappa 4N \log_2(4N)$ flops to compute $\mathbf{F}_{2Q} \mathbf{V} \mathbf{F}_{2P}$ via 2D FFT; $24N$ flops to compute the Hadamard product $\mathbf{L} \circ (\mathbf{F}_{2Q} \mathbf{V} \mathbf{F}_{2P})$; and $\kappa 4N \log_2(4N)$ flops to compute the IDFT ([inverse discrete Fourier transform](#)) in equation A-16. We use $\kappa = 5$ for the *radix-2* algorithm (Van Loan, 1992, p. 15). By replacing these flops into Algorithm 1, we obtain the complete number of flops

$$f_{conv} = \kappa 16N \log_2(4N) + 24N + it(\kappa 16N \log_2(4N) + 60N) . \quad (68)$$

Figure 2 shows a comparison between $f_{classical}$ (equation 66), f_{cglsls} (equation 67) and f_{conv} (equation 68) for different numbers of observation points up to 1,000,000. As we can see, the total flops count associated with our method is $\textcolor{red}{10^7}$ $\textcolor{blue}{7}$ orders of magnitude smaller than that associated with the classical method and $\textcolor{red}{10^3}$ $\textcolor{blue}{3}$ orders of magnitude smaller than that associated with the standard CGLS method [by](#) [when](#) using a maximum number of iterations $N^{it} = 50$.

Figure 3 shows the time necessary to build matrix \mathbf{A} (equation 13) and solve the linear

system for N varying up to 10,000. With $N = 10,000$, the classical method takes more than sixty-three seconds, the standard CGLS more than twelve seconds, while our method takes only half a second. The CPU used for this test was a Intel Core i7-7700HQ@2.8GHz.

Table 1 shows a comparison between the RAM memory storage associated with each method. ~~The In the~~ classical and standard CGLS methods ~~have to store~~, the whole matrix \mathbf{A} (equation 13) ~~has to be stored~~. For example, a dataset with $N = 10,000$ observation points has an associated sensitivity matrix \mathbf{A} formed by $N^2 = 100,000,000$ elements and takes approximately 763 Megabytes of memory (8 bytes per element). Using the same number of observation points $N = 10,000$, our method requires only 1.831 Megabytes to store the first columns of the BCCB matrices $\mathbf{C}_{\alpha\beta}$ (equation A-6) and 0.6104 Megabytes to store the complex matrix \mathbf{L} (equation A-18) (16 bytes per element). For a bigger dataset with $N = 1,000,000$, the amount of necessary RAM goes to 7,629,395, 183.096 and 61.035 Megabytes, respectively.

APPLICATION TO SYNTHETIC DATA

Our convolutional equivalent layer method requires a regular data grid located on a horizontal and flat observation surface. Here, we evaluate the performance of our method by applying it to simulated airborne magnetic surveys formed by i) a regular data grid on a flat surface; ii) irregular data grids on a flat surface; and iii) regular data grid on undulating surfaces. Note that the simulated surveys in (ii) and (iii) violate the premises of our method.

Simulated airborne surveys

The ~~upper and middle~~ first and second rows in Figure 4 show, respectively, the simulated flight patterns and noise-corrupted total-field anomalies of the airborne magnetic surveys used in our tests. The ~~lower~~ third row in Figure 4 shows the true upward-continued total-field anomalies at $z = -1, 300$ m. The fourth row in Figure 4 shows the true reduced to pole total-field anomalies. All magnetic data (~~middle~~ second and lower rows in Figure 4) are produced by the same three synthetic bodies: two prisms and one sphere with constant total-magnetization vector having inclination, declination and intensity of 0° 35.26, 45° , and 2.8284 3.4641 A/m, respectively. The simulated main geomagnetic field has inclination and declination of 10° ~~and~~37 $^\circ$ 35.26 and 45° , respectively.

Figure 4a shows the simulated airborne survey on a regular grid of 100×50 observation points (totaling $N = 5, 000$ observation points), with a grid spacing of $\Delta x = 101.01$ m and $\Delta y = 163.265$ m along the x - and y -axis, respectively. The noise-corrupted total-field anomaly (~~middle panel~~ second panel of Figure 4a) is calculated at $z = -900$ m, with a pseudorandom Gaussian noise added having null mean and standard deviation of 0.2961 nT.

Figures 4b and 4c show the simulated surveys on irregular grids obtained by perturbing the horizontal coordinates of the regular grid (upper panel in Figure 4a). For the survey shown in Figure 4b, the x and y coordinates are perturbed with sequences of pseudorandom Gaussian noises having null mean and standard deviations equal to 20% of the corresponding grid spacing, which results in absolute values of 20.2 m and 32.6 m, along the x - and y -directions, respectively. For the survey shown in Figure 4c, the standard deviations are equal to 30% of the corresponding grid spacing, which results in absolute values of 30.3

m and 49.0 m along the x - and y -directions, respectively. Their noise-corrupted total-field anomalies (~~middle second~~ panels in Figures 4b and 4c) are calculated on their corresponding irregular grids, on a flat observation surface at $z = -900$ m, with ~~a~~-pseudorandom Gaussian noise added having null mean and standard deviation of 0.2961 nT.

Figures 4d and 4e show the simulated surveys on the same regular grid as shown in Figure 4a (upper panel). The difference is that ~~they are~~observation points are located no longer on a flat, but on undulating surfaces. For the survey shown in Figure 4d, the z coordinates of the undulating surface are defined by a sequence of pseudorandom Gaussian noise having mean -900 m and standard deviation equal to 5% of 900 m, which corresponds to 45 m. For the survey shown in Figure 4e, the standard deviation is equal to 10% of 900 m, which corresponds to 90 m. The noise-corrupted total-field anomalies of these simulated surveys (~~middle second~~ panels in Figures 4d and 4e) are calculated on their corresponding undulating surfaces (upper panels in Figures 4d and 4e), on the same regular grid shown in Figure 4a, with ~~a~~-pseudorandom Gaussian noise added having null mean and standard deviation of 0.2961 nT.

Tests with a regular data grid on a flat surface

Figure 5 show the difference between the simulated (~~middle second~~ row in Figure 4) and predicted data (not shown) obtained by using the classical (the upper row) and our method (the ~~middle second~~ row). From now on, we designate this difference as data residuals. The lower row in Figure 5 shows the convergence curve of our method.

The data residuals using the classical method (equation 10) are shown in the upper panel of Figure 5a, with mean ~~0.3627~~0.4118 nT and standard deviation ~~0.2724~~0.3780 nT.

This process took 17.10 seconds. Using our method, the data residuals (the middle panel in Figure 5a) have mean 0.5223 – 0.9972 nT and standard deviation 0.4323 – 1.3904 nT. In this case, however, the processing time was only 0.18 – 0.25 seconds. As expected, the Euclidean norm of the data residuals produced by our method (lower panel in Figure 5a) decreases ~~and the convergence criterion is~~. ~~The convergence criterion was~~ satisfied close to iteration 50. ~~This result shows that, in practice, our method converges way before N iterations, where N is the number of observations. Setting the convergence to N iterations, besides being unnecessary, it also demands a larger computer processing time, even in this synthetic test with a small number of $N = 5,000$ observations.~~

Tests with irregular data grids on a flat surface

Figure 5b shows the results obtained with the irregular data grid perturbed by using 20% of the regular grid spacing. In this Figure we can see that the data residuals using the classical method (upper panel) yield a good data fit with mean 0.3630 – 0.4084 nT and standard deviation 0.2731 – 0.3862 nT. Using our method, the data residuals (middle panel in Figure 5b) also produced an acceptable data fitting with mean of 0.7147 – 1.3125 nT and standard deviation of 0.5622 – 1.7187 nT. The Euclidean norm of the data residuals obtained by our method (lower panel in Figure 5b) decreases, as expected, and converges to a constant value close to iteration 50.

Figure 5b–c shows the results obtained with the irregular data grid perturbed by using 30% of the regular grid spacing. ~~This figure shows that the~~ ~~The~~ data residuals obtained by the classical method (upper panel in Figure 5c) produced an acceptable data ~~fitting~~^{fit}, having mean 0.3634 – 0.4070 nT and standard deviation 0.2735 – 0.3899 nT. Using our method,

the data residuals (middle panel in Figure 5c) with mean $\underline{0.9186}$ – $\underline{1.5129}$ nT and standard deviation $\underline{0.7354}$ – $\underline{1.8526}$ nT also produced a good data fitting. The convergence of our method (lower panel in Figure 5c) shows that, similarly to the previous results, the Euclidean norm of the residuals decreases; ~~however it starts increasing without achieving an invariance.~~ ~~In this case, we have to stop the algorithm at iteration 40~~ converging to a constant value ~~close to iteration 50~~. Note that this good result was obtained by using a very perturbed data grid (upper panel in Figure 4c).

Tests with regular data grid an undulating surfaces

Figure 5d shows the results obtained with data on the undulating surface varying 5% of $z = 900$ m. In this case, the data residuals either using the classical method (upper panel in Figure 5d) or our method (middle panel in Figure 5d) reveal acceptable data fittings. Using the classical method, data residuals have mean $\underline{0.3712}$ – $\underline{0.4316}$ nT and standard deviation $\underline{0.2870}$ – $\underline{0.4762}$ nT. Using our method, they have mean $\underline{0.9542}$ – $\underline{2.1069}$ nT and standard deviation $\underline{0.8943}$ – $\underline{2.5023}$ nT. Likewise, the Euclidean norm of the data residuals produced by our method (lower panel in Figure 5d) decreases up to iteration 50 and reaches ~~an invariance~~–~~the convergence criterion~~ in the subsequent iterations (~~mean residulas are less than 0.00015 between iterations~~).

Figure 5e shows the results obtained with data on the undulating surface varying 10% of $z = 900$ m. By using the classical approach, the data residuals (upper panel in Figure 5e) yielded a good data fitting, with mean $\underline{0.3865}$ – $\underline{0.4818}$ nT and standard deviation $\underline{0.3216}$ – $\underline{0.6565}$ nT. By using our method, the data residuals (middle panel in Figure 5e) yielded a worse data fitting with mean $\underline{1.6109}$ – $\underline{3.4981}$ nT and standard deviation $\underline{1.6231}$ – $\underline{3.8153}$ nT.

The convergence curve (lower panel in Figure 5e) reveals the inadequacy of our method in dealing with observations on rugged surfaces, as the Euclidean norm of the data residuals ~~decreases slower than do not decrease as much as~~ in previous tests. We stress that, in this test, the undulating surface (upper panel in Figure 4e) varies in a broad range ~~of flight values,~~ from $z = -570$ m to about $z = -1,230$ m. Thus, this simulated airborne magnetic survey greatly violates the requirement of a flat observation surface demanded by our method.

Although our method is formulated to deal with magnetic observations measured on a horizontally regular grid, on a flat surface, the results obtained with synthetic data show that our method is robust in dealing either with irregular grids in the horizontal directions or with uneven surfaces. However, the robustness of our method has limitations. ~~The performance limitation of our method depends on the degree of departure of the (i) x - and y -coordinates from those of the closest regular grid and (ii) the z -coordinates from a constant value. High departures~~ ~~High discrepancies~~ in the x -, y , and z -coordinates lead to unacceptable data fittings (large data residuals), as shown the middle panels in Figures 5c and 5e.

Magnetic data processing

We performed the upward continuations of the synthetic total-field anomalies (~~middle second~~ row in Figure 4) by using the classical method, our convolutional equivalent layer method, and the classical approach in the Fourier domain, which consists in computing the Fourier transform of the total-field anomaly (e.g., Blakely, 1996, p. 317).

Figure 6 shows the ~~continuation residuals defined as the~~ differences between the true

upward-continued total-field anomalies (~~lower~~third row in Figure 4) and the predicted upward-continued total-field anomalies (not shown). We conveniently denote these differences as continuation residuals.

~~Figure 6 shows that the~~ The continuation residuals obtained by using the classical method (upper row) and our method (middle row) are similar to each other in most of the tests. ~~One exception is~~ The exceptions are the synthetic test with data over irregular grid (Figures 4c and 5c) and over an undulating surface (Figures 4e and 5e), which greatly violates the requirement of regular grids or a flat observation surface, demanded by our method. Note that the maximum absolute value of the continuation residuals produced by using our method (middle panel in Figure 6e) are ~~≈ 2.5~~ ≈ 2 times greater than those produced by the classical method (upper panel in Figure 6e). ~~Besides, they are generally concentrated at the boundaries of the study area.~~

In contrast, the continuation residuals obtained by using the classical Fourier approach (lower row in Figure 6) are, in most of the tests, approximately ~~6–2~~ ≈ 2 times greater than those produced by the classical method (upper row in Figure 6) and ~~4–1.5~~ ≈ 2 times greater than those produced by our method (middle row in Figure 6). Note that, ~~similarly~~ similar to our method, the maximum absolute values of the continuation residuals obtained by using the classical Fourier approach are located at the boundaries of the simulated area. However, the values are significantly higher.

~~We~~ Figure 7 shows the differences between the true reduced to pole total-field anomalies (fourth row in Figure 4) and the predicted reduced to pole total-field anomalies (not shown). The true reduced to pole total-field anomalies are generated by using only induced magnetization, with $I_0 = 90^\circ$ and $D_0 = 0^\circ$. Figure 7 shows that the reduced to pole residuals

obtained by using the classical method (upper row) and our method (middle row) have differences when high irregular grids or non flat surfaces are used (Figures 7c and 7e). The absolute values of the reduced to pole residuals are almost ≈ 2 times greater than those of classical method when the 10% standard deviation was used (upper and middle panels in Figure 7e, respectively). As in the continuation test, they are generally concentrated at the boundaries of the study area.

The reduced to pole residuals obtained by using the classical Fourier approach (lower row in Figure 7) are approximately 3.5 times greater than those produced by the classical method (upper row in Figure 7) and 3 times greater than those produced by our method (middle row in Figure 6).

Important to note that the reduction to pole, either using the equivalent layer or the Fourier approach, has the requirement of a previously knowledge of the sources magnetization directions (equation 4) to obtain a correct source parameter estimative, otherwise, only non-phase dependent processing can be used (upward continuation for example).

We also call attention to the following aspects: In applying the classical method, our method, or the classical Fourier approach, we do not expand the data by using a padding scheme. The data residuals (upper and middle rows in Figure 5) ~~and the continuation residuals, the continuation~~ (Figure 6) ~~and~~ reduction to pole residuals (Figure 7) are shown without removing ~~the~~ edge effects. The computational time required by our method is much lower than that required by the classical method and has the same order of magnitude of that required by the classical Fourier approach. However, the classical Fourier approach shows upward-continued ~~and reduced to pole~~ data with strong border effects if no padding scheme is applied to expand the data.

APPLICATION TO FIELD DATA

We applied the convolutional equivalent layer method to the aeromagnetic data of Carajás, northern Brazil. The survey is composed of 131 flight lines along north-south direction with line spacing of $\Delta y = 3,000$ m. Data were measured with average spacing $\Delta x = 7.65$ m along lines, with an average distance to the ground of 900 m. The total number of observation points is $N = 6,081,345$. Figure ??-8a shows the observed total-field anomaly data over the study area.

We compare the results obtained with an interpolated regular grid of $10,000 \times 131$ points, by using the nearest neighbor algorithm, and a decimated irregular grid, also with $10,000 \times 131$ points, totaling. In both cases the total $N = 1,310,000$ observation points in both cases are in the original undulating surface of the flight lines. The decimated grid was generated by choosing the nearest observation points in comparison of the regular grid presented in the interpolation. The mean and standard deviation of this irregular decimated from the regular interpolated are 6.8386 m and 107.7343 m in the x -direction and 30.8799 m and 28.3849 m in the y -direction, respectively. Both application were made with an Intel core i7 7700HQ@2.8GHz processor in single-processing and single-threading modes. Figures ??-a and ??-b show, respectively, the data obtained by interpolation and decimation. With 1,310,000 observation points, it would be necessary 12.49 Terabytes of RAM to store the full sensitivity matrix with the classical method. In this case, our method uses only 59.97 Megabytes, allowing regular desktop computers to be able to process this amount of data.

As the study area is very large, the main magnetic field varies with position. For this application, we set the main field direction as that of a mid location (latitude -6.5° and longitude -50.75°) where the declination is -19.86° and the inclination is -7.4391° . Both

values were calculated using the magnetic field calculator from NOAA at 1st January, 2014 (epoch of the survey). We set the equivalent layer at 300 meters above the ground (600 m below the data). By applying Figure 8b shows the residuals obtained after using our method to the interpolated regular grid (Figure ??a), we obtain the predicted data shown in Figure ??a and data residuals (Figure ??b) fit the interpolated data with mean 0.0762 nT and the standard deviation 0.4886 nT, revealing an acceptable data fitting. Our method took ≈ 390.80 seconds to converge at about 200 iterations.

By applying Figure 8c shows the residuals obtained after using our method to the decimated irregular grid (Figure ??b), we obtain the predicted data shown in Figure ??a and data residuals (Figure ??b) fit the decimated data with mean 0.0717 nT and standard deviation 0.3144 nT with a better fit than the produced by the interpolated data. In this case, our method took ≈ 385.56 seconds to converge at about 200 iterations (Figure ??8d). The convergence curve reveals a good convergence rate obtained with the decimated irregular grid. This result shows the robustness of our method in processing irregular grid. Notice that we used 200 iterations in our method of the interpolated regular grid and the mean residual still decreasing up to 2000. This happens because the invariance convergence criterion was met as the mean residuals are very small and each iteration decreases less than 0.00001.

With 1,310,000 observation points, it would be necessary 12.49 Terabytes of RAM to store the full sensitivity matrix with the classical method. In this case, our method uses only 59.97 Megabytes, allowing regular desktop computers to be able to process this amount of data.

Finally, Figure ??9a shows the upward-continued magnetic data to a horizontal plane

located at an altitude of 5, 000 m using the estimated equivalent layer obtained by applying our method to the decimated irregular grid(Figure ??b). This process took ≈ 2.64 seconds, showing good results without visible errors or border effects. Figure 9b shows the upward-continued magnetic data to a horizontal plane located at an altitude of 5, 000 m using the classical Fourier filtering method to the decimated irregular grid. This process took ≈ 0.5 seconds. Altough, no border effect could be observed, this result showed unexpected high values instead of the expected attenuation of the anomalies. We stress that we did not use a padding scheme to expand the data.

CONCLUSIONS

We have proposed a fast equivalent-layer technique for processing magnetic data called convolutional equivalent layer method. We have demonstrated that the sensitivity matrix associated with planar equivalent layers of dipoles has a BTTB structure for the particular case in which the dipoles are aligned with the horizontal and regular grid of magnetic data. The product of such matrices and arbitrary vectors represents a 2D discrete convolution that can be efficiently computed via 2D Fast Fourier Transform by using only the elements forming the first column of the matrix. By using this property, we have developed a fast and memory efficient iterative method for estimating the physical-property distribution on the equivalent layer.

Comparisons between the estimated physical-property distribution obtained with our method and the classical approach that solves the least-squares normal equations via Cholesky decomposition show similar results. The differences in total number of floating-point operations (flops), memory usage and computation time, however, are noticeable. For a mid-size grid of 100×50 points, the total number of flops is about four orders of magnitude smaller

than that required by the classical method. Besides, our method uses less than 1% of the RAM and takes about 3% of the computation time associated with the classical method in this case. Significantly better results can be obtained with larger data sets.

Tests with synthetic data show that the computational time required by our method has the same order of magnitude of that required by the classical approach in the Fourier domain to perform magnetic data processing. However, the classical Fourier approach shows considerable larger border effects if no previous padding scheme is used to expand the data. Besides, although both methods require the magnetic data be on a planar and regular grid, tests with synthetic data show the robustness of our method to deal with data either on irregular grids or on undulating observation surfaces.

While the classical equivalent-layer method would require 12.49 Terabytes of RAM to store the full sensitivity matrix associated with the irregular grid of 1,310,000 observation points over the Carajás Province, northern Brazil, our method requires only 59.97 Megabytes. When performed on a standard laptop computer with an Intel Core i7 7700HQ@2.8GHz processor in single-processing and single-threading modes, the total times spent by our method to estimate the physical-property distribution over the equivalent layer and to compute the upward continuation of the 1,310,000 magnetic observations over the Carajás province was approximately 385.56 seconds and 2.64 seconds.

Further investigation could usefully explore different preconditioning strategies to improve the convergence rate of our method. Besides, considerably more work will need to be done to generalize our convolutional equivalent layer method for dealing with irregularly spaced data sets on undulating observation surfaces.

FIGURES

TABLES

APPENDIX A

BTTB MATRIX-VECTOR PRODUCT

This appendix follows a similar approach to that presented by Takahashi et al. (2020) to efficiently compute the product of the sensitivity matrix \mathbf{A} (equations 13) and a generic vector \mathbf{b} . Let this product be represented by

$$\mathbf{t} = \mathbf{A} \mathbf{b} , \quad (\text{A-1})$$

where

$$\mathbf{t} = \mathbf{t}_{xx} + \mathbf{t}_{xy} + \mathbf{t}_{xz} + \mathbf{t}_{yy} + \mathbf{t}_{yz} + \mathbf{t}_{zz} \quad (\text{A-2})$$

and

$$\mathbf{t}_{\alpha\beta} = \mathbf{A}_{\alpha\beta} \mathbf{b} . \quad (\text{A-3})$$

Let us also consider that vectors

$$\mathbf{t}_{\alpha\beta} = \begin{bmatrix} \mathbf{t}_{\alpha\beta}^0 \\ \vdots \\ \mathbf{t}_{\alpha\beta}^{Q-1} \end{bmatrix}_{N \times 1} \quad (\text{A-4})$$

and

$$\mathbf{b} = \begin{bmatrix} \mathbf{b}^0 \\ \vdots \\ \mathbf{b}^{Q-1} \end{bmatrix}_{N \times 1} \quad (\text{A-5})$$

are composed of $P \times 1$ vectors $\mathbf{t}_{\alpha\beta}^q$ and \mathbf{b}^q , respectively, where q is the block index (equations 32 and 34). From equation A-3, we obtain an auxiliary matrix-vector product given by

$$\mathbf{w}_{\alpha\beta} = \mathbf{C}_{\alpha\beta} \mathbf{v} , \quad (\text{A-6})$$

where $\mathbf{C}_{\alpha\beta}$ is a $4N \times 4N$ block circulant matrix with circulant blocks (BCCB) (e.g., Davis, 1979, p. 184),

$$\mathbf{w}_{\alpha\beta} = \begin{bmatrix} \mathbf{w}_{\alpha\beta}^0 \\ \vdots \\ \mathbf{w}_{\alpha\beta}^{Q-1} \\ \mathbf{0}_{2N \times 1} \end{bmatrix}_{4N \times 1}, \quad (\text{A-7})$$

$$\mathbf{w}_{\alpha\beta}^q = \begin{bmatrix} \mathbf{t}_{\alpha\beta}^q \\ \mathbf{0}_{P \times 1} \end{bmatrix}_{2P \times 1}, \quad (\text{A-8})$$

$$\mathbf{v} = \begin{bmatrix} \mathbf{v}^0 \\ \vdots \\ \mathbf{v}^{Q-1} \\ \mathbf{0}_{2N \times 1} \end{bmatrix}_{4N \times 1} \quad (\text{A-9})$$

and

$$\mathbf{v}^q = \begin{bmatrix} \mathbf{b}^q \\ \mathbf{0}_{P \times 1} \end{bmatrix}_{2P \times 1}, \quad (\text{A-10})$$

with $\mathbf{0}_{2N \times 1}$ and $\mathbf{0}_{P \times 1}$ being vectors of zeros. The key point here is that the auxiliary matrix-vector product (equation A-6) represents a 2D discrete convolution and can be efficiently computed by using the 2D Fast Fourier Transform (2D FFT).

The BCCB matrix $\mathbf{C}_{\alpha\beta}$ (equation A-6) is formed by $2Q \times 2Q$ blocks, where each block $\mathbf{C}_{\alpha\beta}^q$ is a $2P \times 2P$ circulant matrix. The entire BCCB matrix $\mathbf{C}_{\alpha\beta}$ is defined by properly

downshifting its first block column

$$[\mathbf{C}_{\alpha\beta}]_{(0)} = \begin{bmatrix} \mathbf{C}_{\alpha\beta}^0 \\ \vdots \\ \mathbf{C}_{\alpha\beta}^{Q-1} \\ \mathbf{0}_{2P \times 2P} \\ \mathbf{C}_{\alpha\beta}^{-Q+1} \\ \vdots \\ \mathbf{C}_{\alpha\beta}^{-1} \end{bmatrix}_{4N \times 2P}, \quad (\text{A-11})$$

where $\mathbf{0}_{2P \times 2P}$ is a matrix of zeros. Similarly, each block $\mathbf{C}_{\alpha\beta}^q$, $q = -Q + 1, \dots, Q - 1$, is obtained by properly downshifting its first column

$$\mathbf{c}_{\alpha\beta}^q = \begin{bmatrix} a_{q0}^{\alpha\beta} \\ \vdots \\ a_{q(P-1)}^{\alpha\beta} \\ 0 \\ a_{q(-P+1)}^{\alpha\beta} \\ \vdots \\ a_{q(-1)}^{\alpha\beta} \end{bmatrix}_{2P \times 1}, \quad (\text{A-12})$$

where $a_{qp}^{\alpha\beta}$, $p = -P + 1, \dots, P - 1$, are the elements of matrix component $\mathbf{A}_{\alpha\beta}$ described in terms of block indices q and p (equations 32–35). The BCCB matrix $\mathbf{C}_{\alpha\beta}$ is diagonalized by $\mathbf{F}_{2Q} \otimes \mathbf{F}_{2P}$, where “ \otimes ” denotes the Kronecker product (e.g., Horn and Johnson, 1991, p. 242) and \mathbf{F}_{2Q} and \mathbf{F}_{2P} are the $2Q \times 2Q$ and $2P \times 2P$ unitary DFT matrices (Davis, 1979, p. 31). Due to this property, the auxiliary matrix-vector product (equation A-6) can be computed as follows (Takahashi et al., 2020):

$$\mathbf{F}_{2Q}^* [\mathbf{L}_{\alpha\beta} \circ (\mathbf{F}_{2Q} \mathbf{V} \mathbf{F}_{2P})] \mathbf{F}_{2P}^* = \mathbf{W}_{\alpha\beta}, \quad (\text{A-13})$$

where “ \circ ” denotes the Hadamard (element-wise) product (e.g., Horn and Johnson, 1991, p. 298), “ $*$ ” denotes the complex conjugate, $\mathbf{W}_{\alpha\beta}$ and \mathbf{V} are $2Q \times 2P$ matrices obtained by rearranging, respectively, vectors $\mathbf{w}_{\alpha\beta}$ (equation A-7) and \mathbf{v} (equation A-9) along their rows and $\mathbf{L}_{\alpha\beta}$ is a $2Q \times 2P$ matrix given by

$$\mathbf{L}_{\alpha\beta} = \sqrt{4QP} \mathbf{F}_{2Q} \mathbf{G}_{\alpha\beta} \mathbf{F}_{2P}, \quad (\text{A-14})$$

with

$$\mathbf{G}_{\alpha\beta} = \begin{bmatrix} (\mathbf{c}_{\alpha\beta}^0)^{\top} \\ \vdots \\ (\mathbf{c}_{\alpha\beta}^{Q-1})^{\top} \\ \mathbf{0}_{1 \times 2P} \\ (\mathbf{c}_{\alpha\beta}^{-Q+1})^{\top} \\ \vdots \\ (\mathbf{c}_{\alpha\beta}^{-1})^{\top} \end{bmatrix}_{2Q \times 2P}, \quad (\text{A-15})$$

defined by the first columns $\mathbf{c}_{\alpha\beta}^q$ (equation A-12), $q = -Q + 1, \dots, Q - 1$, of all circulant blocks $\mathbf{C}_{\alpha\beta}^q$ (equation A-11). Hence, the whole BCCB matrix $\mathbf{C}_{\alpha\beta}$ does not have to be formed, but only its first column. Besides, the symmetries defined by equations 42–65 imply that all elements of $\mathbf{G}_{\alpha\beta}$ can be obtained by using only the first column of $\mathbf{A}_{\alpha\beta}$. Consequently, the whole matrices $\mathbf{A}_{\alpha\beta}$ do not have to be formed as well, but only their first columns.

It is important noting that the left side of equation A-13 represents the 2D Inverse Discrete Fourier Transform (2D IDFT) of the term in brackets. This term, in turn, represents the Hadamard product of $\mathbf{L}_{\alpha\beta}$ (equation A-14) and the 2D Discrete Fourier Transform (2D DFT) of \mathbf{V} . Similarly, equation A-14 shows that $\mathbf{L}_{\alpha\beta}$ is obtained by computing the 2D DFT of matrix $\mathbf{G}_{\alpha\beta}$ (equation A-15). Hence, equations A-13 and A-14 can be efficiently

computed by using the 2D FFT. After that, the elements of vector $\mathbf{t}_{\alpha\beta}$ (equation A-3) can be retrieved from the first quadrant of matrix $\mathbf{W}_{\alpha\beta}$ (equation A-13). By combining the results obtained for all components $\alpha\beta$, $\alpha, \beta = x, y, z$, we can show that

$$\mathbf{F}_{2Q}^* [\mathbf{L} \circ (\mathbf{F}_{2Q} \mathbf{V} \mathbf{F}_{2P})] \mathbf{F}_{2P}^* = \mathbf{W}, \quad (\text{A-16})$$

where

$$\mathbf{W} = \mathbf{W}_{xx} + \mathbf{W}_{xy} + \mathbf{W}_{xz} + \mathbf{W}_{yy} + \mathbf{W}_{yz} + \mathbf{W}_{zz} \quad (\text{A-17})$$

and

$$\mathbf{L} = \mathbf{L}_{xx} + \mathbf{L}_{xy} + \mathbf{L}_{xz} + \mathbf{L}_{yy} + \mathbf{L}_{yz} + \mathbf{L}_{zz}, \quad (\text{A-18})$$

with $\mathbf{L}_{\alpha\beta}$ defined by equation A-14. Then, the elements of \mathbf{t} (equation A-1) are obtained from the first quadrant of \mathbf{W} (equations A-16 and A-17).

Finally, it can be shown that the product

$$\mathbf{t} = \mathbf{A}^\top \mathbf{b} \quad (\text{A-19})$$

can be computed by using equation A-16. The difference is that, in this case, matrices $\mathbf{G}_{\alpha\beta}$ (equation A-15) are defined by using the new vectors

$$\mathbf{c}_{\alpha\beta}^q = \begin{bmatrix} a_{q0}^{\alpha\beta} \\ \vdots \\ a_{q(-P+1)}^{\alpha\beta} \\ 0 \\ a_{q(P-1)}^{\alpha\beta} \\ \vdots \\ a_{q1}^{\alpha\beta} \end{bmatrix}_{2P \times 1}. \quad (\text{A-20})$$

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1 This table shows the RAM memory usage (in Megabytes) for storing the whole $N \times N$ matrix \mathbf{A} (equation 13), the first columns of the BCCB matrices $\mathbf{C}_{\alpha\beta}$ (equation A-6) (both need 8 bytes per element) and the matrix \mathbf{L} (equation A-18) (16 bytes per element).

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1 Schematic representation of an $N_x \times N_y$ regular grid of points (black dots) with $N_x = 3$ and $N_y = 2$, where each point has an associated index. This index may represent i or j , that are associated with observation points (x_i, y_i, z_0) and equivalent sources (x_j, y_j, z_c) . Left panel shows an example of x -oriented grid, with indices varying along x -axis, while right panel shows an example of y -oriented grid, with indices varying along y -axis.

2 Number of flops associated with classical method (equation 66), the standard CGLS method (equation 67) and our method (equation 68, all of them with $N^{it} = 50$). The number of observation points N varies from 5,000 to 1,000,000.

3 Comparison between the runtime of the equivalent-layer technique using the classical method, standard CGLS method and our method. The values for the standard CGLS and our method use $N^{it} = 50$ iterations.

4 Synthetic tests: the simulated airborne magnetic surveys - The first row shows the grids of observation points and the undulating observation surfaces that simulate the airborne magnetic surveys. The second row shows the noise-corrupted total-field anomalies produced by the synthetic sources and calculated on the simulated airborne magnetic survey shown in the first row. The third row shows the noise-free total-field anomalies produced by the synthetic sources at $z = -1,300$ m (the true upward-continued total-field anomalies). The fourth row shows the noise-free total-field anomalies produced by the synthetic sources at inclination $I_0 = 90^\circ$ (the true reduced to pole total-field anomalies). The results shown in these last three rows were obtained by using the simulated airborne magnetic surveys as follows: (a) A regular grid of 100×50 observation points in the x - and y -directions and a flat observation surface at $z = -900$ m. An irregular grid with uncertainties of (b) 20% and (c) 30% in the x - and y -coordinates and a flat observation surface at $z = -900$ m . A

regular grid of 100×50 observation points in the $x-$ and $y-$ coordinates and an undulating observation surface with uncertainties of (d) 5% and (e) 10%. The black lines represent the horizontal projection of the sources .

5 Synthetic tests: the data residuals and convergence - The first row shows the data residuals using the classical method. The second and third rows show, respectively, the data residuals and the convergence curves using the convolutional equivalent layer (our method). The results shown in these three rows were obtained by using the simulated airborne magnetic surveys shown in Figure 4, i.e.: (a) A regular grid of 100×50 observation points in the $x-$ and $y-$ directions and a flat observation surface at $z = -900$ m. An irregular grid with uncertainties of (b) 20% and (c) 30% in the $x-$ and $y-$ coordinates and a flat observation surface at $z = -900$ m . A regular grid of 100×50 observation points in the $x-$ and $y-$ coordinates and an undulating observation surface with uncertainties of (d) 5% and (e) 10%. The black lines represent the horizontal projection of the sources .

6 Synthetic tests: the data residuals of the upward-continued total-field anomalies (third row in Figure 4). The data residuals of the upward-continued total-field anomalies are defined as the difference between the noise-free total-field anomaly produced by the synthetic sources at $z = -1,300$ m (third row in Figure 4) and the predicted total-field anomaly at $z = -1,300$ m obtained by using three methods: the classical method (first row); the convolutional equivalent layer (second row); and the classic approach in the Fourier domain (third row). The results shown in these three rows were obtained by using the simulated airborne magnetic surveys shown in Figure 4, i.e.: (a) A regular grid of 100×50 observation points in the $x-$ and $y-$ directions and a flat observation surface at $z = -900$ m. An irregular grid with uncertainties of (b) 20% and (c) 30% in the $x-$ and $y-$ coordinates and a flat observation surface at $z = -900$ m . A regular grid of 100×50 observation points in

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7 Synthetic tests: the data residuals of the reduced to pole total-field anomalies (fourth row in Figure 4). The data residuals of the reduced to pole total-field anomalies are defined as the difference between the noise-free total-field anomaly produced by the synthetic sources at the pole (fourth row in Figure 4) and the predicted total-field anomaly obtained by using three methods: the classical method (first row); the convolutional equivalent layer (second row); and the classic approach in the Fourier domain (third row). The results shown in these three rows were obtained by using the simulated airborne magnetic surveys shown in Figure 4, i.e.: (a) A regular grid of 100×50 observation points in the x - and y -directions and a flat observation surface at $z = -900$ m. An irregular grid with uncertainties of (b) 20% and (c) 30% in the x - and y -coordinates and a flat observation surface at $z = -900$ m . A regular grid of 100×50 observation points in the x - and y -coordinates and an undulating observation surface with uncertainties of (d) 5% and (e) 10%. The black lines represent the horizontal projection of the sources .

8 (a) Observed total-field anomaly over the Carajás Province, northen Brazil. The aeromagnetic survey was flown in 131 north-south flight lines at an average altitude of 900 m, totaling $N = 6,081,345$ observation points. (b) Data residuals, defined as the difference between the regular interpolated grid data (not shown) and the predicted data (not shown), with mean of 0.0762 nT and standard deviation of 0.4886 nT. (c) Data residuals, defined as the difference between the irregular decimated grid data (not shown) and the predicted data (not shown), with mean of 0.0717 nT and standard deviation of 0.3144 nT. (d) Convergence curve using our method to the decimated irregular grid of the real data of Carajás Province, Brazil.

9 Upward continuations of real data of Carajás Province, Brazil at altitude of 5,000
m by using: (a) the convolutional equivalent layer (our method) and (b) the classical Fourier
method.

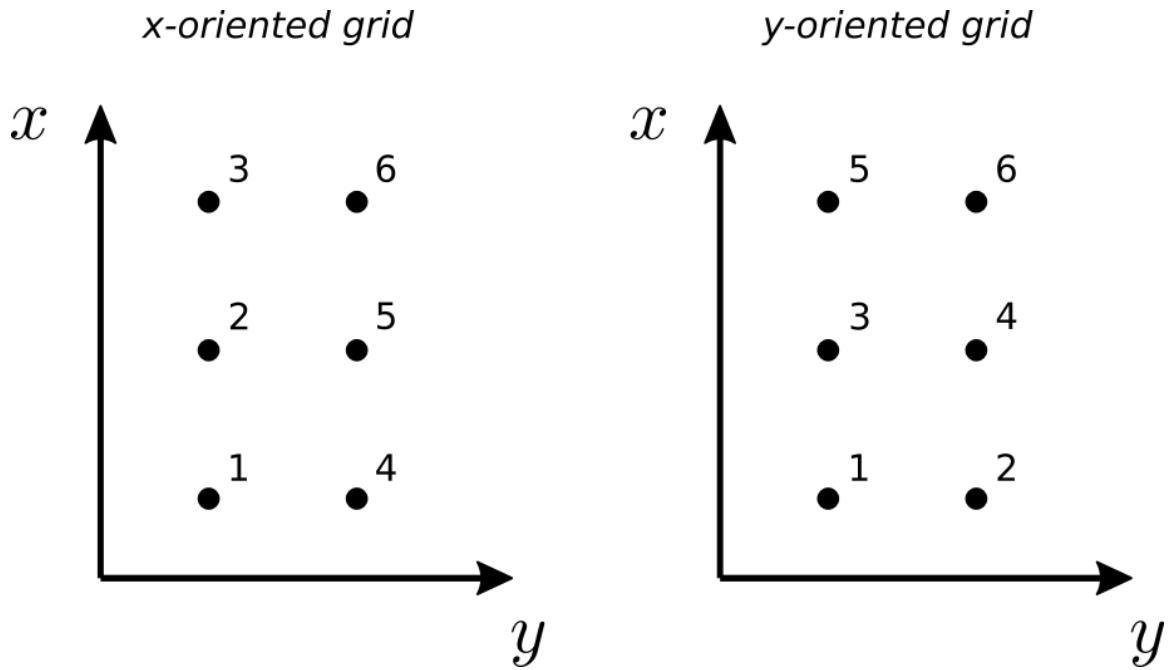


Figure 1: Schematic representation of an $N_x \times N_y$ regular grid of points (black dots) with $N_x = 3$ and $N_y = 2$, where each point has an associated index. This index may represent i or j , that are associated with observation points (x_i, y_i, z_0) and equivalent sources (x_j, y_j, z_c) . Left panel shows an example of x -oriented grid, with indices varying along x -axis, while right panel shows an example of y -oriented grid, with indices varying along y -axis.

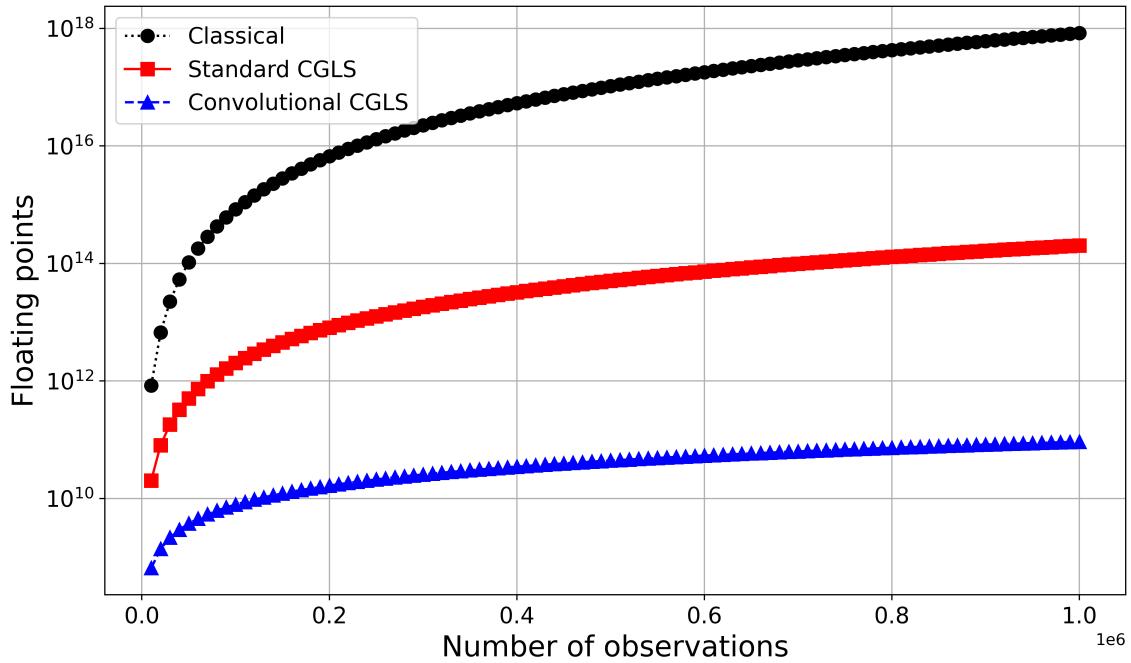


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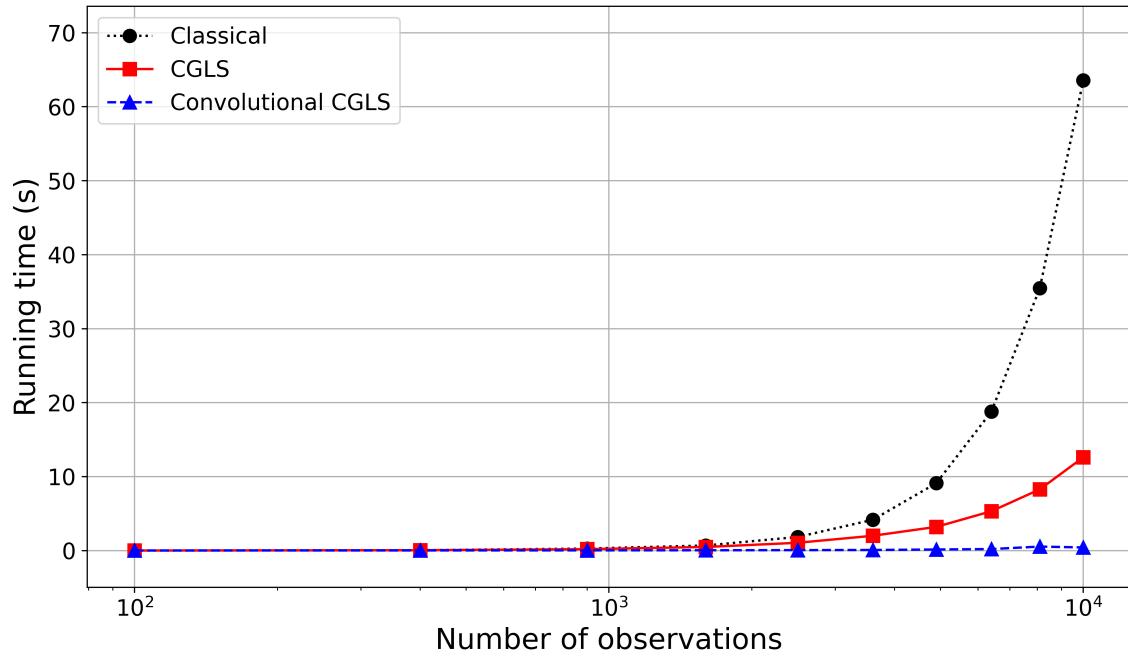


Figure 3: Comparison between the runtime of the equivalent-layer technique using the classical method, standard CGLS method and our method. The values for the standard CGLS and our method use $N^{it} = 50$ iterations.

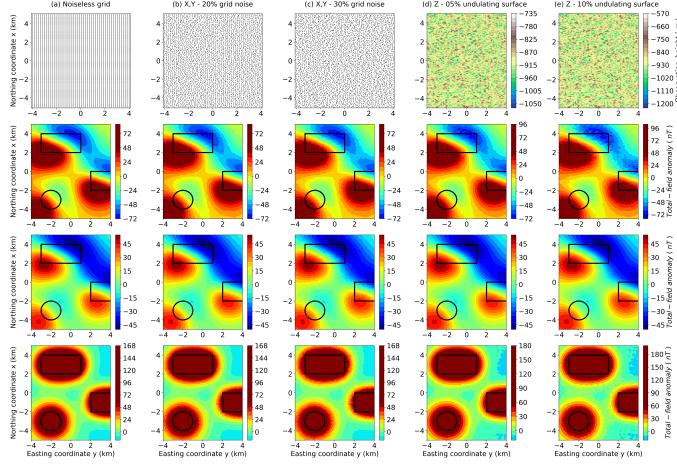


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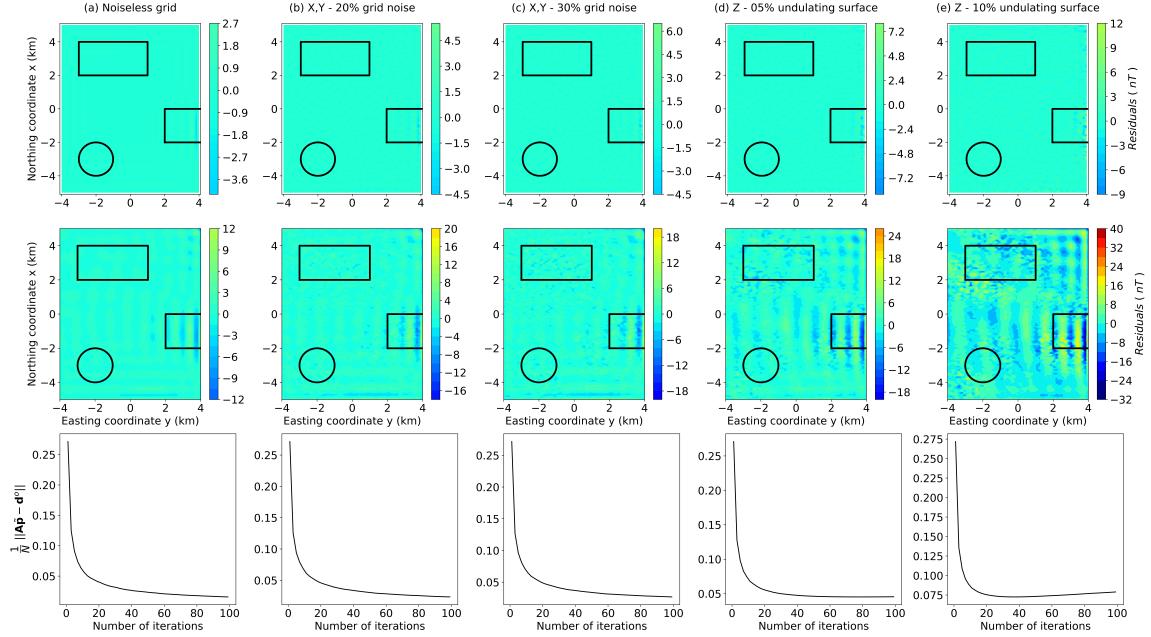


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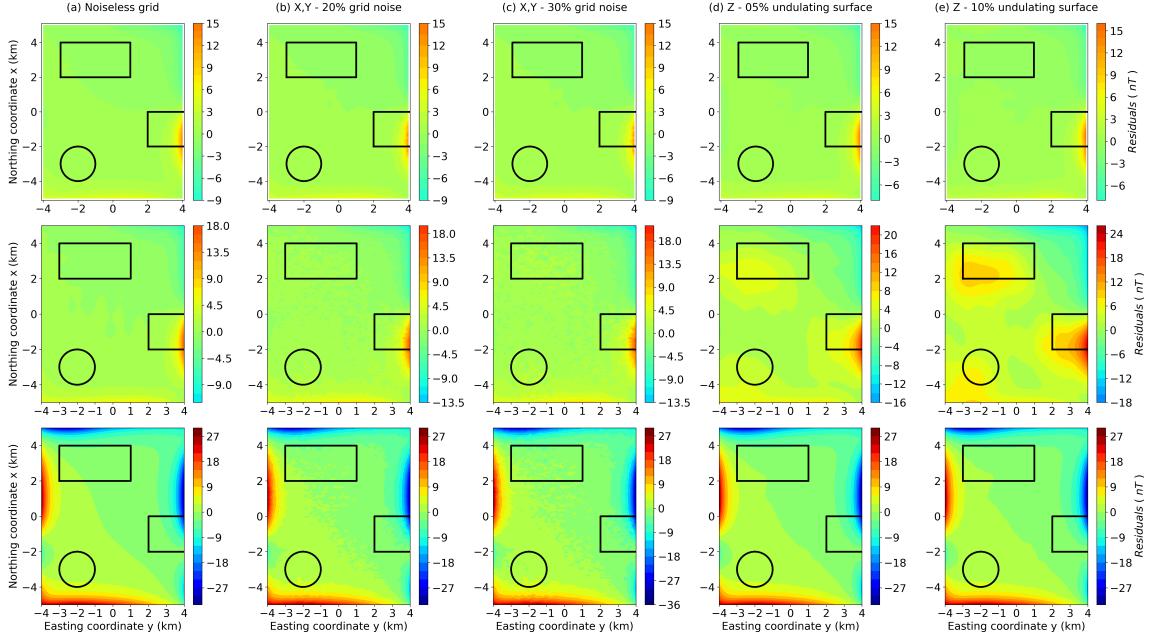


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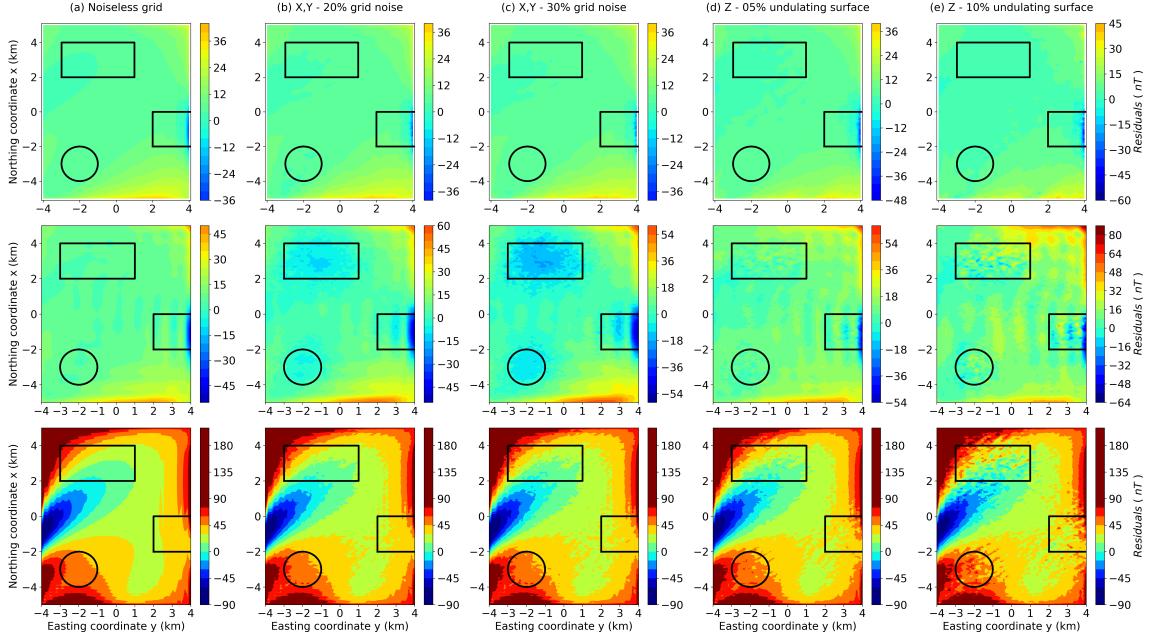


Figure 7: Synthetic tests: the data residuals of the reduced to pole total-field anomalies (fourth row in Figure 4). The data residuals of the reduced to pole total-field anomalies are defined as the difference between the noise-free total-field anomaly produced by the synthetic sources at the pole (fourth row in Figure 4) and the predicted total-field anomaly obtained by using three methods: the classical method (first row); the convolutional equivalent layer (second row); and the classic approach in the Fourier domain (third row). The results shown in these three rows were obtained by using the simulated airborne magnetic surveys shown in Figure 4, i.e.: (a) A regular grid of 100×50 observation points in the x - and y -directions and a flat observation surface at $z = -900$ m. An irregular grid with uncertainties of (b) 20% and (c) 30% in the x - and y -coordinates and a flat observation surface at $z = -900$ m. A regular grid of 100×50 observation points in the x - and y -coordinates and an undulating observation surface with uncertainties of (d) 5% and (e) 10%. The black lines represent the horizontal projection of the sources .

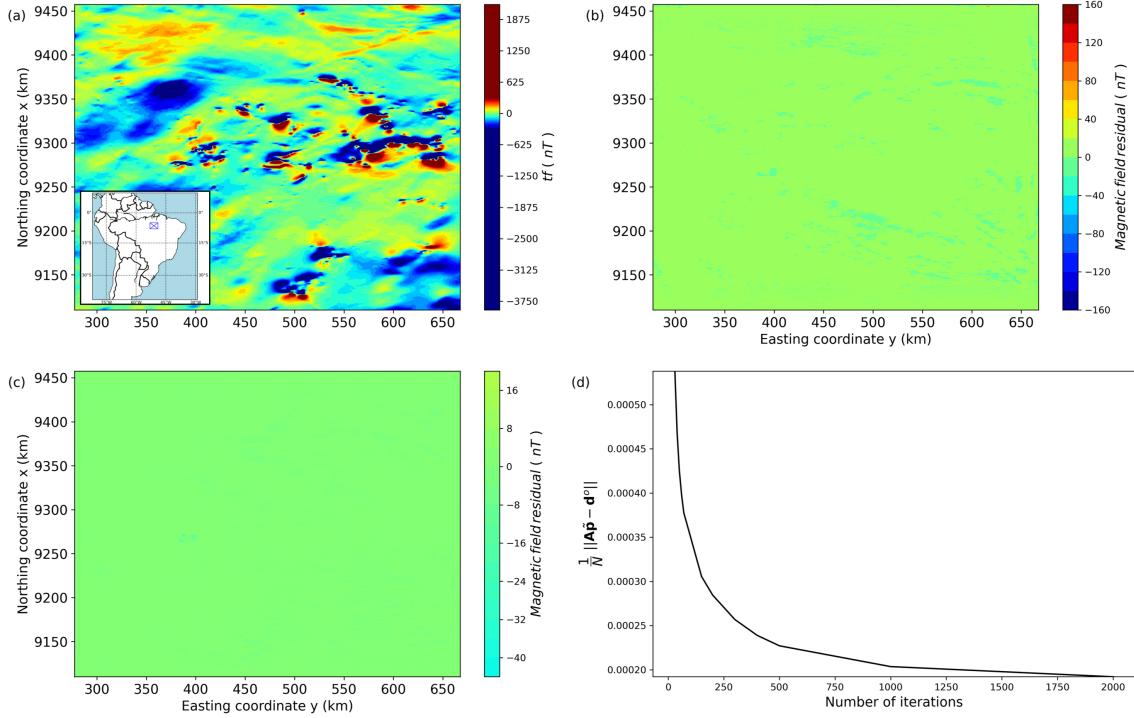


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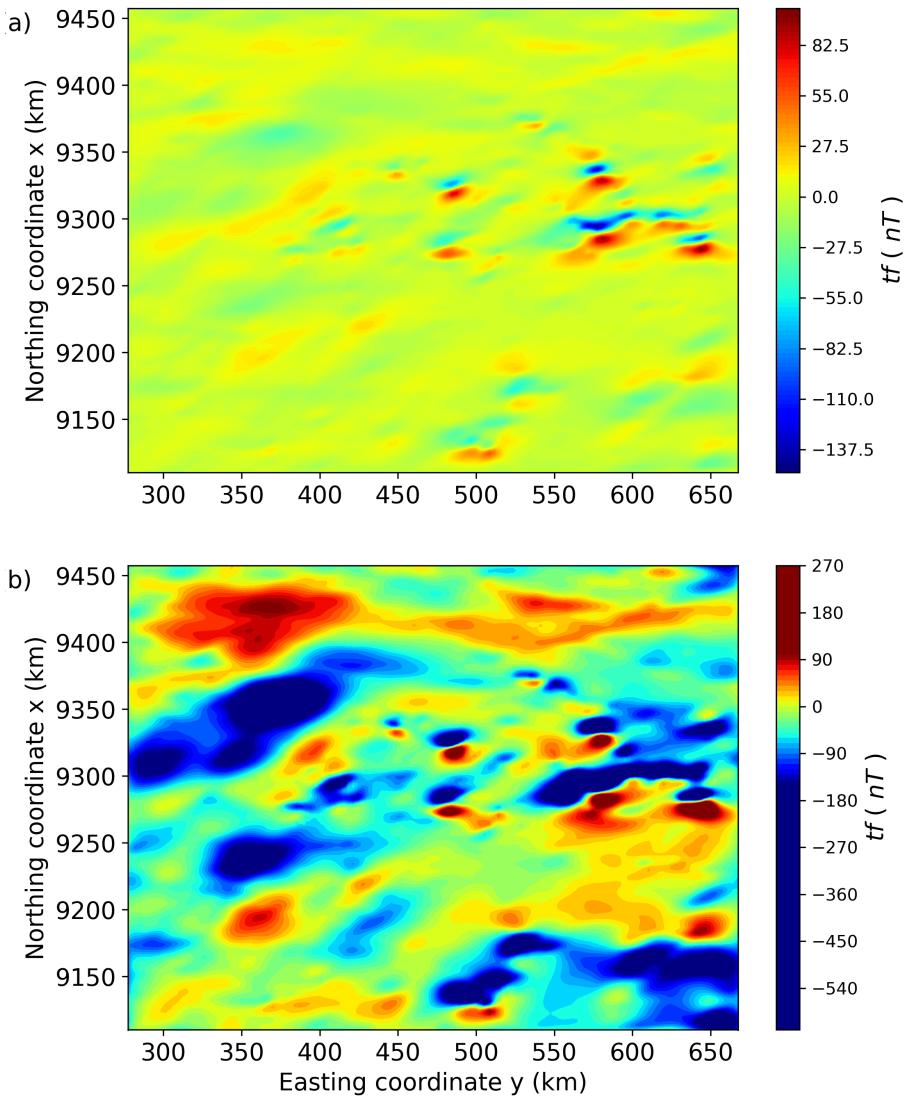


Figure 9: Upward continuations of real data of Carajás Province, Brazil at altitude of 5,000 m by using: (a) the convolutional equivalent layer (our method) and (b) the classical Fourier method.

N	A	First columns of matrices $\mathbf{C}_{\alpha\beta}$	L
100	0.0763	0.0183	0.00610
400	1.22	0.0744	0.0248
2,500	48	0.458	0.1528
10,000	763	1.831	0.6104
40,000	12,207	7.32	2.4416
250,000	476,837	45.768	15.3
500,000	1,907,349	91.56	30.518
1,000,000	7,629,395	183.096	61.035

Table 1: This table shows the RAM memory usage (in Megabytes) for storing the whole $N \times N$ matrix **A** (equation 13), the first columns of the BCCB matrices $\mathbf{C}_{\alpha\beta}$ (equation A-6) (both need 8 bytes per element) and the matrix **L** (equation A-18) (16 bytes per element).