An efficient regularization method for a large scale ill-posed geothermal problem

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Abstract

The inverse geothermal problem consists of estimating the temperature distribution below the earth’s surface using measurements on the surface. The problem is important since temperature governs a variety of geologic processes, including the generation of magmas and the deformation style of rocks. Since the thermal properties of rocks depend strongly on temperature the problem is non-linear.

The problem is formulated as an ill-posed operator equation, where the righthand side is the heat-flux at the surface level. Since the problem is ill-posed regularization is needed. In this study we demonstrate that Tikhonov regularization can be implemented efficiently for solving the operator equation. The algorithm is based on having a code for solving a well-posed problem related to the above mentioned operator. The algorithm is designed in such a way that it can deal with both 2D and 3D calculations.

Numerical results, for 2D domains, show that the algorithm works well and the inverse problem can be solved accurately with a realistic noise level in the surface data.
1 Introduction

Temperature controls many physical properties of rocks and governs a variety of geologic processes, including the generation of magmas and the deformation style of rocks [1, 3, 20, 26], see also [9, 10, 23]. Thus, knowledge of the thermal structure of the lithosphere is essential for understanding geophysical observations and dynamic processes in the deep interior. However, temperature cannot be directly measured beyond a few kilometers depth, in deep mines or drill holes [12, 21]. To obtain the information about the temperature distribution in the lithosphere, the surface observations, e.g. surface heat flux, are extrapolated to the lithosphere base using a heat conduction model [5, 8]. Since the thermal conductivity of rocks are temperature and pressure dependent [11] the problem is non-linear. For the one-dimensional case an analytical solution can be found [7, 14] but for the two or three-dimensional case numerical modeling is required. Several previous studies investigated this type of problems using the least squares approach, see [32, 29], and also [19, 9], but these authors did not strictly treat the nonlinear aspect resulting from the temperature-dependent thermal properties of rocks.

The inverse geothermal problem can be formulated as a boundary value problem for the stationary heat equation in a domain, but where the temperature and heat-flux data are only known on a part of the boundary. This is often referred to as the Cauchy problem for the heat equation [2, 15], see also [4] and [5]. The problem is ill-posed and a regularization technique is needed to stabilize the computations, see [35] for a good overview of regularization techniques used in geophysics.

We reformulate the problem as a non-linear operator equation; where the operator maps the heat-flux at the lower boundary to the heat-flux at the surface level. The operator can thus be evaluated by solving a well-posed boundary value problem for the temperature distribution inside the domain under consideration. The inverse problem is solved by using Tikhonov regularization. In our implementation we solve the normal equations, that correspond to the minimization of the Tikhonov functional, using the conjugate gradient method. The bulk of the computational work in the algorithm consists of solving several linear systems of equations originating from a finite difference discretization of the underlying well-posed boundary value problem. We also show how to treat the non-linearity of the problem. An important feature of the algorithm is that it can be generalized to the case surface data is available on a 2D area and the temperature distribution for a 3D region below the surface is desired.

The outline of this paper is as follows: In Section 2 we introduce a well-posed boundary value problem for the heat-equation that allows us to reformulate the inverse geothermal problem as a non-linear operator equation. We also give the the details of our thermal model. In Section 3 we start by investigating
the inverse geothermal problem for the case of a constant thermal conductivity and heat production. For this case we can use the singular value decomposition and demonstrate the ill-posedness of the problem. Also we apply Tikhonov regularization and give numerical results. The non-linear case is treated in Section 4. We show how to implement Tikhonov regularization efficiently by using the conjugate gradient method. Finally, we give numerical results in Section 4.2 and present conclusions in Section 5.

2 The Direct Problem

For geothermal calculations the goal is to estimate the temperature distribution in the lithosphere \cite{9}. The model is two dimensional and confined to the domain \((x, z) \in \Omega = (0, L_x) \times (0, L_z)\), where \(x\) is the lateral coordinate, and \(z\) is the depth.

The boundary of the domain \(\Omega\) consists of the surface where measurements are available, the base where we specify the heat flux, and the sides where we assume a zero flux. Thus the thermal model is the following: The temperature \(T(x, z) \in C^2(\Omega) \cap C^1(\bar{\Omega})\) satisfies,

\[
\begin{cases}
(k T_x)_x + (k T_z)_z + A_p = 0, & 0 < x < L_x, 0 < z < L_z, \\
T_x(0, z) = T_x(L_x, z) = 0, & 0 < z < L_z, \\
T(x, 0) = T_0(x), & 0 < x < L_x, \\
k T_z(x, L_z) = Q_m(x), & 0 < x < L_x,
\end{cases}
\]  

(2.1)

where \(k := k(x, z, T)\) is the thermal conductivity and \(A_p := A_p(x, z, T)\), or more commonly \(A_p := A_p(x, z)\), is the heat production. The above problem is well-posed \cite{13} and provided that we have access to the heat-flow \(Q_m\) and the surface temperature \(T_0\) we can compute the temperature distribution \(T(x, z)\) in the domain \(\Omega\). This allows us to define a non-linear operator as follows: In (2.1) we consider the surface temperature \(T_0\) to be fixed. If, in addition, we specify the heat-flux \(Q_m\) then we have a well-posed problem that we can solve for the temperature distribution \(T(x, z)\) in the domain. Having obtained \(T(x, z)\) we can compute the surface heat-flux \(Q_0\). This means that we can introduce a mapping from \(Q_m\) to \(Q_0\) that we denote by

\[
Q_0 = k T_z(x, 0) = \mathcal{K}(k, A_p, T_0) Q_m,
\]  

(2.2)

In order to simplify the notation in the paper we also introduce a function

\[
[T, Q_0] = \mathcal{F}(k, A_p, T_0, Q_m), \quad Q_0 = k T_z|_{z=0}.
\]  

(2.3)

Note that since the model is non-linear an iterative procedure is required for evaluating the operator \(\mathcal{K}\) or the function \(\mathcal{F}\). We refer to evaluating the operator
$K$, for a given heat-flux $Q_m$, i.e. solving (2.1), as the direct problem. In contrast solving operator equation (2.2), for a given surface heat-flux $Q_0$, is referred to as the inverse problem, see Section 3.

2.1 Numerical Solution of the Direct Problem

If the thermal conductivity and the heat production are independent of the temperature, i.e. $\kappa := \kappa(x, z)$ and $A_p := A_p(x, z)$, then solving (2.1) is a standard problem.

By introducing a computational grid $(x_i, z_j)$, $i = 1, \ldots, N$ and $j = 1, \ldots, M$, on the domain $\Omega$, we represent the function $T(x, z)$ by its values at the grid points, i.e. the unknowns are $T = (T_{i,j}) = (T(x_i, z_j))$. Similarly the boundary data $T_0$ and $Q_m$ are represented by vectors, e.g. $Q_m = (Q_m(x_i))$. Thus, having a code for solving (2.1) effectively means that we have have a functional relation

$$[T, Q_0] = FDM(\kappa, A_p, T_0, Q_m),$$

(2.4)

where $T$, $\kappa$, and $A_p$ are matrices, and $T_0$, $Q_0$ and $Q_m$ are vectors. The above function represents a discrete approximation of the function $F$, see (2.3), and can be used to approximate the operator $K$, introduced in (2.2).

In our work we use finite differences [31] to discretize the differential equation. The details of our code are described briefly. At interior points we discretize the governing equation using a symmetric difference approximation,

$$(\kappa T_x)(x_i, z_j) = \frac{1}{\Delta x^2} \left( \kappa_{i-\frac{1}{2},j} T_{i-1,j} - (\kappa_{i-\frac{1}{2},j} + \frac{\kappa_i + \frac{1}{2},j}{\kappa_i + \frac{1}{2},j}) T_{i,j} + \kappa_{i+\frac{1}{2},j} T_{i+1,j} \right),$$

(2.5)

where the values of $\kappa(x, z)$ at half-index points are evaluated using linear interpolation. Neumann-type boundary conditions, e.g. $\kappa T_z(x, L_z) = Q_m(x)$ or $T_z(0, z) = 0$, are implemented using one-sided finite differences. The finite difference approximation leads to a linear system of equations $Au = b$, where $u$ is a vector that contains the temperature values at the grid points. Having solved for $u$ we can compute the surface heat flux using a one sided difference quotient.

We remark that the linear system $Au = b$ is large and sparse but can be solved effectively using either direct or iterative methods. Also, the matrix $A$ only depends on the thermal conductivity $\kappa$. The righthand side $b$ contains the boundary data, i.e. $T_0$ and $Q_m$, and also the heat production values $A_p$ at interior points.
Table 1: The parameters used in our model of the thermal conductivity $\kappa$ and heat production $A_p$ in the different lithosphere layers, see [6, 11, 14].

<table>
<thead>
<tr>
<th>Layer</th>
<th>$\kappa_0$ (W/m°C)</th>
<th>$c$ (1/o°C)</th>
<th>$A_p$ (W/m$^3$)</th>
</tr>
</thead>
<tbody>
<tr>
<td>Sediment</td>
<td>3.0</td>
<td>0</td>
<td>$1.5 \cdot 10^{-6}$</td>
</tr>
<tr>
<td>Upper Crust</td>
<td>3.0</td>
<td>$1.5 \cdot 10^{-3}$</td>
<td>$2.0 \cdot 10^{-6}$</td>
</tr>
<tr>
<td>Lower Crust</td>
<td>2.6</td>
<td>$1.0 \cdot 10^{-4}$</td>
<td>$0.45 \cdot 10^{-6}$</td>
</tr>
<tr>
<td>Mantle</td>
<td>2.5</td>
<td>$-2.5 \cdot 10^{-4}$</td>
<td>$0.02 \cdot 10^{-6}$</td>
</tr>
</tbody>
</table>

2.2 A Numerical Example with Artificial Data

In order to illustrate the geothermal problem we give numerical results for case where we picked a realistic heat-flux $Q_m$ at the base of the model and solved the problem (2.1) using finite differences. Note that this is an artificially generated test problem with a known solution. For the actual application the heat-flux at the base of the model, i.e. $Q_m$, would be unknown. Having a test problem with a known heat flux $Q_m$ is convenient for testing purposes.

In our previous paper [9] we used actual measured data at the surface to estimate $Q_m$ for the eastern Tibetan margin. In this work we use the same model for the thermal conductivity and the heat production. The thermal conductivity, at crustal conditions, is assumed to vary inversely with temperature [7, 8], i.e.

$$\kappa = \frac{\kappa_0}{1 + cT},$$  

(2.6)

where $\kappa_0$ is the thermal conductivity at surface conditions and $c$ is an experimentally determined constant. At low temperatures the thermal conductivity generally decreases with increasing temperature, i.e. $c > 0$, while in the range $300 - 500$°C the conductivity varies very little with temperature. At temperatures above $500$°C the conductivity instead increases with temperature, i.e. $c < 0$ [30]. Further we assume that the heat production, i.e. $A_p$ in (2.1), is constant in each lithosphere layer. For discussions of more detailed heat production models see [27, 18]. The specific values used in our simulations are listed in Table 1, see also Figure 1. Note that the heat production is not temperature dependent.

For the numerical simulations, with the above thermal model, we picked a computational grid $(x_i, z_j)$, with $N = 300$ equidistant grid points in the $x$–direction, and $M = 250$ grid points in the $z$–direction. The physical size of the computational domain was where $L_x = 560$ km and $L_z = 80$ km. At the surface we used a constant temperature $T_0 = 10$°C. The heat-flux $Q_m$ is seen in Figure 1.

Starting from the initial guess $T^{(0)}(x, z) = 0$ we used our finite difference code
to solve (2.1) iteratively,

\[ [T^{(k+1)}, Q^{(k+1)}_0] = FDM(\kappa(T^{(k)}), A_p(T^{(k)}), T_0, Q_m); \]

until the update \( \|T^{(k+1)} - T^{(k)}\|_F \) is sufficiently small. At each step we solve a sparse linear system of size 75000 × 75000, with 371412 non-zero elements.

The convergence is rather fast, see Figure 1. The final temperature distribution \( T(x, z) \), after convergence, is also displayed. We emphasize that the choice of the initial approximation \( T^{(0)}(x, z) = 0 \) is not very good and thus we start with a large error. Still only about five iterations are needed to reach a sufficiently small error.

It is interesting to note that the temperature in the mantle layer varies in the interval 500 < \( T < 1150^\circ\text{C} \). This means that the thermal conductivity changes significantly, \( 5.1 < \kappa < 6.3\text{W/m}\text{oC} \) [34], within the mantle layer. Thus a constant value for the thermal conductivity in the whole mantle layer would not give an accurate solution.

3 The Linear Inverse Problem and Regularization

In the previous section we presented the thermal model that is used for our test simulations. Provided that we have access to the heat-flow \( Q_m \) at the base of the model and the surface temperature \( T_0 \) we can solve (2.1) and obtain the temperature distribution \( T(x, z) \). In practice we cannot measure the heat-flux \( Q_m \). Instead we measure both the temperature \( T_0 \) and the heat-flux \( Q_0 \) at surface level; i.e. we have Cauchy data \([T_0, Q_0]\) and want to compute the corresponding heat-flux \( Q_m \). Thus we consider (2.2) as a non-linear operator equation,

\[ KQ_m = Q_0, \quad K := K(\kappa, A_p, T_0), \]

(3.7)

that can be solved for the desired heat-flux \( Q_m \).

In our model the thermal conductivity and heat production are temperature dependent. Even though we are interested in solving the non-linear equation (3.7) we begin by studying the case where the coefficients \( \kappa \) and \( A_p \) are independent of the temperature. We assume that an approximate temperature distribution \( T^{(0)} \) is available and use the notation

\[ \kappa := \kappa(x, z) := \kappa(x, z, T^{(0)}), \quad \text{and}, \quad A_p := A_p(x, z) := A_p(x, z, T^{(0)}), \]

in the remainder of this section.
In the case where the thermal conductivity $\kappa$ and heat production $A_p$ are both independent of the temperature we can transform the operator (2.2) into a linear operator. The procedure is as follows: First, find a function $V(x, z)$ that satisfies (2.1) with the heat-flux $T_z(L_z, x) = 0$ at the base of the model, i.e.

$$[V, Q_1] = F(\kappa, A_p, T_0, 0).$$

Second, set the surface temperature and the heat production to zero, so that $W$ is defined by

$$[W, Q_2] = F(\kappa, 0, 0, Q_m).$$

Then $T = V + W$ and we have $Q_0 = Q_1 + Q_2$. We conclude that

$$\tilde{\kappa}Q_m = Q_2 = Q_0 - Q_1, \quad \tilde{\kappa} = \kappa(\kappa, 0, 0),$$

and that $\tilde{\kappa}$ is a linear operator.
3.1 Analysis of the Linear Problem

In this section we study the linear operator \( \tilde{K} \) analytically. In order to be able to carry out the analysis we simplify the model and assume that the thermal conductivity \( \kappa \) is a constant.

Taking \( A_p = 0, T_0 = 0 \) and \( \kappa \) a constant, the problem (2.1) transforms into

\[
\begin{cases}
T_{xx} + T_{zz} = 0, & 0 < x < L_x, \quad 0 < z < L_z, \\
T_x(0, z) = T_x(L_x, z) = 0, & 0 < z < L_z, \\
T(x, 0) = 0, & 0 < x < L_x, \\
\kappa T_z(x, L_z) = Q_m(x), & 0 < x < L_x.
\end{cases}
\]

(3.11)

By separation of variables, we can write the solution as

\[
T(x, z) = \sum_{n=0}^{\infty} a_n \cos\left(\frac{n\pi x}{L_x}\right) \left(-e^{-\frac{n\pi}{L_x}z} + e^{\frac{n\pi}{L_x}z}\right).
\]

(3.12)

The condition \( \kappa T_z(x, L_z) = Q_m(x) \) can also be written as

\[
\kappa T_z(x, L_z) = \sum_{n=1}^{\infty} \kappa a_n \cos\left(\frac{n\pi x}{L_x}\right) \left(\frac{n\pi}{L_x} e^{-\frac{n\pi}{L_x}L_z} + \frac{n\pi}{L_x} e^{\frac{n\pi}{L_x}L_z}\right) = Q_m(x),
\]

If we then write \( Q_m(x) \) as a Fourier cosine series,

\[
Q_m(x) = \sum_{n=1}^{\infty} A_n \cos\left(\frac{n\pi x}{L_x}\right), \quad \text{where} \quad A_n = \frac{2}{L_x} \int_0^{L_x} Q_m(x) \cos\left(\frac{n\pi x}{L_x}\right) dx,
\]

(3.13)

and compare the two expressions we find that the constants \( a_n \) can be written as

\[
a_n = \frac{2\kappa^{-1}}{n\pi(e^{-\frac{n\pi}{L_x}L_z} + e^{\frac{n\pi}{L_x}L_z})} \int_0^{L_x} Q_m(x) \cos\left(\frac{n\pi x}{L_x}\right) dx.
\]

(3.14)

At the earth’s surface the heat flux \( \kappa T_z(x, 0) = Q_0 \) is then given by

\[
\kappa T_z(x, 0) = \sum_{n=1}^{\infty} \kappa \frac{2n\pi a_n}{L_x} \cos\left(\frac{n\pi x}{L_x}\right) = Q_0(x),
\]

substituting for \( a_n \) and reorganizing the expression gives the result

\[
\int_0^{L_x} \sum_{n=0}^{\infty} \frac{4}{L_x(e^{-\frac{n\pi}{L_x}L_z} + e^{\frac{n\pi}{L_x}L_z})} \cos\left(\frac{n\pi x'}{L_x}\right) \cos\left(\frac{n\pi x}{L_x}\right) Q_m(x') dx' = Q_0(x).
\]

The above solution formula can be interpreted as follows: If proper function spaces are introduced the operator \( \tilde{K} \) is compact and its singular system [13] is
given by
\[
\{\sigma_n, u_n, v_n\} = \left\{ \frac{1}{e^{-n\pi \frac{L_z}{L_x}} + e^{n\pi \frac{L_z}{L_x}}}, \frac{2}{\sqrt{L_x}} \cos\left(\frac{n\pi L_x}{L_z}\right), \frac{2}{\sqrt{L_x}} \cos\left(\frac{n\pi L_x}{L_z}\right) \right\}. \quad (3.15)
\]

The degree of ill-posedness for an operator equation, e.g. (3.10), is defined in terms of its singular values [13, p. 40]. If the singular values decay as \(O(n^{-\alpha})\), for some \(\alpha > 0\), then the problem is mildly ill-posed. If on the other hand the singular values decay exponentially, i.e. as \(O(e^{-n\alpha})\), then the operator equation is severely ill-posed. In our case the singular values decays exponentially, which indicates that the operator equation \(\tilde{K}Q_m = Q_0\) is severely ill-posed.

### 3.2 Numerical Solution of the Linear Inverse Problem

After having introduced a computational grid \((x_i, z_j)\) the functions \(T, \kappa, \) and \(A_p\) are represented by matrices, and \(Q_0, T_0, \) and \(Q_m\) are vectors. The linear operator \(\tilde{K}\), see (3.10), is approximated by a matrix-relation
\[
\tilde{K}Q_m := FDM(\kappa, 0, 0, Q_m). \quad (3.16)
\]
The product of the matrix \(\tilde{K}\) and a vector \(Q_m\) can be computed by solving a well-posed boundary value problem using our code. Let \(\{e_i\}\) be the standard basis on \(\mathbb{R}^N\). Since the above relation is linear we can compute the matrix \(\tilde{K}\) explicitly by,
\[
\tilde{K}(i, i) = \tilde{K}e_i = FDM(\kappa, 0, 0, e_i), \quad i = 1, \ldots, N. \quad (3.17)
\]
In the remainder of this section we are concerned with solving the ill-conditioned linear system of equations,
\[
\tilde{K}Q_m = \tilde{Q}_0, \quad (3.18)
\]
where the vector \(\tilde{Q}_0\) is computed as, see (3.10),
\[
[V, Q_1] = FDM(\kappa, A_p, T_0, 0), \quad \text{and} \quad \tilde{Q}_0 = Q_0 - Q_1. \quad (3.19)
\]
The linear system (3.18) is ill-conditioned and regularization is needed[13, 16]. We use Tikhonov regularization[25, 33]; i.e. the regularized solution is computed by minimizing the Tikhonov functional,
\[
\|\tilde{K}Q_m^\lambda - \tilde{Q}_0\|_2^2 + \lambda^2 \|Q_m^\lambda\|_2^2, \quad (3.20)
\]
where \(\lambda > 0\) is a regularization parameter. If the matrix \(\tilde{K}\) is explicitly known then the solution can be computed efficiently using the singular value decomposition [16].
The above techniques can be summarized as follows: For the case of temperature
independent coefficients the inverse problem \((3.7)\) can be solved using Tikhonov
regularization. We obtain a functional relation,

\[ Q^\lambda_m := LT(\kappa, A_p, T_0, Q_0, \lambda). \] (3.21)

where, as previously, \(\kappa\) and \(A_p\) are matrices, and \(Q^\lambda_0, T_0,\) and \(Q_0\) are vectors,
and \(\lambda\) is a scalar.

The regularization parameter \(\lambda\) controls the weight given to the regularization
term, \(\lambda^2 \|Q^\lambda_m\|^2\), relative to the residual \(\|KQ^\lambda_m - \tilde{Q}_0\|^2\). If only a little regular-
ization is used, i.e. if \(\lambda\) is small, then the noise can be magnified during the
computations and destroy the solution. If more regularization is used, i.e. a
larger \(\lambda\), then the noise is effectively filtered out but the solution may be too
smooth and the residual might be large.

The appropriate choice of the parameter \(\lambda\) represents a compromise between
accuracy, i.e. keeping the residual small, and stability, i.e. preventing the
noise from being magnified too much. There are several standard methods for
selecting a good regularization parameter for a concrete problem. Two of the
most widely used are the Discrepancy principle and the L-curve.

The Discrepancy principle, attributed to Morozov[24], can be used if the noise
level in the data is available. The regularization parameter \(\lambda\) is then defined as

\[ \lambda := \sup\{\lambda > 0 \mid \|KQ^\lambda_m - \tilde{Q}_0\| \leq \tau \delta\}, \] (3.22)

where \(\delta\) is the noise level, and \(\tau \geq 1\) is a suitable constant. This means that the
method finds a solution where the residual is of the same order of magnitude
as the noise level in the data. The discrepancy principle is known to work
well under very general assumptions regarding the ill-posed problem its applied
to[13].

The L-curve, originally introduced in [22], see also [17], which is a plot of the
residual norm \(\|KQ^\lambda_m - \tilde{Q}_0\|^2\) versus \(\|Q^\lambda_m\|^2\) for different values of \(\lambda\), is a con-
vienient graphical tool for analysis of ill-posed problems. The L-curve displays
the compromise between the minimization of the two quantities. For discrete
ill-posed problems, the L-curve, when plotted in a log–log scale, often has a char-
acteristic L-shaped appearance with a distinct corner separating the vertical and
horizontal parts of the curve. Less regularization corresponds to the uppermost
part of the curve and more regularization corresponds to the rightmost of the
curve. The optimal regularization parameter, that ensures a solution with a
good balance between the noise magnification and the approximation error is
the value \(\lambda\) that corresponds to the corner[16].
Figure 2: The singular values \( \{\sigma_k\} \) for the computed matrix \( \tilde{K} \). The singular values decay exponentially, illustrating that the operator equation \( \tilde{K}Q_m = \tilde{Q}_0 \) is severely ill-posed. Only about 60 of the values \( \sigma_k \) are above the machine precision.

### 3.3 A Numerical Example

In order to illustrate the ill-posedness of the operator \( \tilde{K} \), and also the effects of applying Tikhonov regularization we solve a numerically generated test problem and present the results. The test problem is the one presented in Section 2.2, i.e. the grid \((x_i, z_j)\) consists of \( N = 300 \) points in the \( x \)-direction and \( M = 250 \) points in the \( z \)-direction.

By selecting a vector \( Q_m \), representing the exact heat flux at the base of the model, and solving the direct problem repeatedly until convergence, we obtain Cauchy data \([T_0, Q_0]\), at the surface level, matrices \( \kappa \) and \( A_p \) representing the thermal conductivity and heat production on the grid.

First, using the technique described in Section 3, we compute the matrix \( \tilde{K} \), of size \( 300 \times 300 \), representing the linear operator. Since the matrix is relatively small we can compute its singular value decomposition. The results are displayed in Figure 2. The results clearly demonstrate that the problem is severely ill-posed.

Secondly we add artificially generated noise to the Cauchy data \([T_0, Q_0]\). For this test we use normally distributed noise of variance \( \epsilon_1 = 5 \cdot 10^{-3} \, \text{W/m} \, \text{C} \) and \( \epsilon_2 = 5 \cdot 10^{-5} \, \text{W/m} \, \text{m}^2 \, \text{C} \) to obtain the simulated measurements \([T_0^*1, Q_0^*2]\). The simulated data vectors are illustrated in Figure 3. The noise is of a realistic order of magnitude. In order to find a good value of the regularization parameter we use the \( L \)-curve, and minimize the Tikhonov functional for a range of parameter values \( \lambda \). The method predicts that \( \lambda = 6.3 \cdot 10^{-3} \) is an appropriate value for the
regularization parameter. Even though the data contains significant amounts of noise the reconstruction of the heat flux at the base of the model is very successful.

In order to illustrate the ill-posedness of the problem we also solve the equation using less regularization. The solution $Q^\lambda_m$ computed using $\lambda = 7 \cdot 10^{-4}$ is displayed in Figure 4. Clearly the solution is highly oscillating as a result of noise magnification during the computations. Also we present a solution computed using $\lambda = 9 \cdot 10^{-2}$. In this case too much regularization results in a too smooth solution and features from the exact solution are not reconstructed accurately.
Figure 4: We display the Tikhonov solutions $Q_m^\lambda$, using $\lambda = 7 \cdot 10^{-4}$ (left, blue curve) and $\lambda = 9 \cdot 10^{-2}$ (right, blue curve). In the first case the not enough regularization has been applied and in the second case we have too much regularization. The exact solution $Q_m$ (black dashed curve) is also presented.

4 The Non-linear Problem

We recall that our goal is to solve the non-linear case, where both the thermal conductivity $\kappa$ and the heat production $A_p$ depend on the temperature. Our idea is to solve a sequence of linear problems using Tikhonov regularization. Let $T^{(0)}$ be an approximate temperature distribution. The algorithm is as follows:

First solve the linearized problem using Tikhonov regularization to obtain

$$Q^{(k+1), \lambda}_m := LT(\kappa(T^{(k)}), A_p(T^{(k)}), T_0, Q_0, \lambda).$$  \hfill (4.23)

The new approximate heat-flux $Q^{(k+1), \lambda}_m$, at the base of the model, is then used to compute the next temperature distribution

$$T^{(k+1)} := FDM(\kappa(T^{(k)}), A_p(T^{(k)}), T_0, Q^{(k+1), \lambda}_m).$$  \hfill (4.24)

The process is repeated until the update $\|T^{(k+1)} - T^{(k)}\|_F$ is sufficiently small.

The algorithm works well, provided that a sufficiently large regularization parameter $\lambda$ is used, but is computationally demanding. At each step we need to form a linear system of equations $Au = b$, see Section 2.1. For relatively small grid sizes, i.e. 2D calculations, it is feasible to solve the linear system using Gaussian elimination. If the matrix $\tilde{K}$, see (3.17), is desired then we make use of the observation that the matrix $A$ only depends on the thermal conductivity $\kappa(T^{(k)})$. Thus we can compute the $LU$ decomposition of the sparse matrix $A$ and reuse it when evaluating $\tilde{K}e_i$, for all the basis vectors $e_i$. After having computed the matrix $\tilde{K}$ explicitly we can easily find the regularized solution of the inverse problem by solving the least squares problem (3.20).
4.1 Modified Normal Equations and an Iterative Method

For large grid sizes, e.g. 3D calculations, it is necessary to avoid forming $\tilde{K}$, see (3.17), explicitly. Instead we note that the minimization problem (3.20) is equivalent to solving the modified normal equations,

$$(\tilde{K}^T \tilde{K} + \lambda^2 I)Q^\lambda_m = \tilde{K}^T \tilde{Q}_0.$$ (4.25)

Instead of forming the matrix $\tilde{K}$ explicitly we solve the normal equations using the conjugate gradient method[28]. The matrix-vector product $\tilde{K}Q_m$ can be evaluated by solving the direct problem using finite differences, i.e.

$$\tilde{K}Q_m := FDM(\kappa, 0, 0, Q_m).$$ (4.26)

The core of the finite difference code is a linear system of equations $Au = b$, where $u_k = T(x_i, z_j)$, $k = (i-1)n+j$. The right hand side $b$ is zero except for a block that contains the vector $Q_m$. In matrix form this can be written as

$$b = W_2 Q_m, \quad W_2 \in \mathbb{R}^{NM \times N},$$

where $W_2$ is a matrix that places the values in $Q_m$ at the correct locations in the vector $b$. After having computed the temperatures $T_{ij}$ at the grid points a finite difference formula is used to obtain the heat-flux at the surface level. This can also be represented in matrix form as

$$\tilde{Q}_0 = W_1 u, \quad W_1 \in \mathbb{R}^{N \times NM},$$

where $W_1$ is a matrix that computes a finite difference approximation of the surface heat flux given the temperature values stored in the vector $u$. The matrix $\tilde{K}$ can be written in factorized form as,

$$\tilde{Q}_0 = W_1 u = W_1 A^{-1} b = W_1 A^{-1} W_2 Q_m = \tilde{K}Q_m,$$ (4.27)

and its transpose can be written as,

$$\tilde{K}^T = W_2^T (A^T)^{-1} W_1^T.$$ (4.28)

In order to evaluate a matrix-vector product $\tilde{K}^T Q$ we need to solve a system of equations involving the matrix $A^T$. Our finite difference code is written so it can compute matrix-vector products involving both $\tilde{K}$ and $\tilde{K}^T$. In order to evaluate the product,

$$(\tilde{K}^T \tilde{K} + \lambda^2 I)Q_m$$

we need to form the matrix $A$ once and we need to solve two linear systems of equations (one with $A$ and one with $A^T$). Note that if the LU decomposition of $A$ is computed then the relation $A^T P^T = U^T L^T$ can be used to solve the system involving $A^T$ efficiently. If the matrix $A$ is too large, e.g. the 3D case, then instead we use an iterative method, e.g. CG or GMRES[28].
Thus in the implementation of (4.23) the minimization problem (3.20) is solved
by finding the solution $Q^{(k+1),\lambda}$ of the linear system of equations,
\[
((\tilde{K}^{(k)})^T\tilde{K}^{(k)} + \lambda^2 I)Q^{(k+1),\lambda} = (\tilde{R}^{(k)})^T\tilde{Q}_0,
\]
where $\tilde{K}^{(k)}$ is the matrix defined by the relation (4.26) above; with $\kappa := \kappa(T^{(k)})$.
The previous solution $Q^{m,\lambda}_0$ can be used as an initial guess to start the CG
iterations.

4.2 Numerical Tests for the Non-Linear Problem

Here we present the results from the numerical testing of the methods. For all
tests we use the numerically constructed example from Section 2.2. The grid
$(x_i, z_j)$ consists of $N = 300$ points in the $x$–direction and $M = 250$ points in
the $z$–direction. For the tests a vector $Q_m$, representing the exact heat-flux at
the base of the model, was selected and the direct problem solved repeatedly
until convergence. In all cases the direct problems were solved with sufficient
accuracy that the generated Cauchy data $[T_0, Q_0]$ can be considered exact.

To the generated Cauchy data we add normally distributed noise of variance $\epsilon_1$
and $\epsilon_2$ to obtain the simulated measurements $[T_0^{\epsilon_1}, Q_0^{\epsilon_2}]$. The vectors $T_0$ and
$Q_0$ are of different orders of magnitude so it is important to use different noise
levels for the temperature and heat-flux vectors respectively. For our tests we
used the noise levels $\epsilon_1 = 5 \times 10^{-3} \, ^\circ C$ and $\epsilon_2 = 5 \times 10^{-5} \, W/m^2\circ C$. The simulated
data vectors are illustrated in Figure 5. The noise levels are realistic. The same
data vectors and noise are used for all tests presented in the text below.

The purpose of the tests is to illustrate how the algorithm for solving the non-

![Figure 5: The simulated Cauchy data at the surface level. We display the heat-flux $Q_0$ (left) and the temperature $T_0$. In both cases we display both the exact function (blue,dashed curve) and the noisy function (black, solid curve).]
Iteration number: $k$

Step size: $||T(k) - T(k-1)||_F$

Horizontal Coordinate: $x$ [km]

Heat-flux: $Q_m$

Figure 6: Results for non-linear Tikhonov regularization. We display the heat-flux $Q_m$ at the base of the model (right, dashed blue curve) the approximation $Q_m^{(10),\lambda}$ (right, black curve), and the convergence history $\|T^{(k),\lambda} - T^{(k-1),\lambda}\|_F$, measured in the Frobenius norm (left).

For the first test we use the simplest implementation of Tikhonov regularization, i.e. the steps (4.24) and 4.24), works in practice. For the linear problem, i.e. the steps (4.24) and 4.24), works in practice. For the Tikhonov regularization to work well a good regularization parameter is needed. We suggest that the $L$-curve is used and in our tests we use the value $\lambda = 6.3 \cdot 10^{-3}$ that was obtained by the $L$-curve for the constant coefficient case, see Section 3.2.

The results are presented in Figure 6. The convergence is rapid and only 3 iterations is enough to reach a level where the first decimal digits of the temperatures $T^{(k)}$ stops changing. The large initial error is due to the poor starting guess $T^{(0)} = 0$.

We remark that this is a very inefficient approach. At each step of the algorithm we need to solve two different types of finite difference equations. First, the functional relation (3.19) is implemented by first forming the matrix in the finite difference equations and computing its $LU$ decomposition. Second, we need to form the finite difference equations defined by the relation (3.16), compute its $LU$ decomposition and solve $2N$ triangular linear systems to obtain the matrix $\tilde{K}^{(k)}$. Minimizing the small Tikhonov functional, e.g. (3.20), is cheap by comparison and the cost can be neglected.

In our second test we solve the same problem again using the iterative implemen-
Figure 7: Results obtained using the iterative implementation of Tikhonov regularization. We display the convergence history $\|T^{(k)} - T^{(k-1)}\|_F$, measured in the Frobenius norm (left) and also the number of CG iterations needed to reach the stopping criteria at each step of the algorithm (right). The stopping criteria for the CG iterations were set at either $10^{-9}$ (blue curves) and $10^{-5}$ (red curves).

The computational work for the second test is as follows. For each step of the algorithm we need to form two linear systems corresponding to (3.19) and (3.16). The first system is solved once and the second needs to be solved twice during each step in the CG algorithm. The amount of work required is significantly less than for the first test above.

In order to further reduce the amount of computational work we stop the CG iterations when the relative residual is less than $10^{-5}$. This means the linear systems are solved less accurately. The results are also displayed in Figure 7. Here it is enough to perform 5 iterations to reach convergence. A total 10 systems of finite difference equations needs to be formed, and 58 linear system solves are needed. The obtained solution $Q^{(5)}_m$, after 5 steps of the algorithm, is displayed in Figure 8.
5 Concluding Remarks

In this paper we have presented the inverse geothermal problem. The application is important since many geophysical process are influenced by the temperature. The physical properties of rocks depend strongly on the temperature and the problem is non-linear.

We have demonstrated that the problem is severely ill-posed and regularization is needed. One of the most popular regularization techniques is Tikhonovs method. The method requires that we rewrite the geothermal problem as an operator equation, and minimize a related least squares problem. The tests in the paper demonstrates that this approach works well. Though the results may be sensitive with respect to the thermal model used the influence from random measurement errors is relatively small and the computations are numerically stable.

Our implementation is based on having a finite difference solver for a related, well-posed, direct problem. We show how to use software for the direct problem when implementing Tikhonov regularization for the inverse problem. Our algorithm is iterative and the conjugate gradient method is used for solving the normal equations that appear due to the Tihonov regularization. The method is efficient and solving the inverse problem requires only about an order of magnitude additional work compared to solving only the direct problem. It is worth noting that instead starting from a finite element solver for the direct problem would work equally well.
In our work we consider only 2D models however the algorithm can be expected
to work well for 3D calculations as well. As long as the corresponding direct
problem can be solved we can also solve the inverse problem. Calculations on a
realistic full 3D model is something we intend to do in the future.

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