BOUNDARY INTEGRAL FORMULAE FOR THE RECONSTRUCTION OF ELECTRIC AND ELECTROMAGNETIC INHOMOGENEITIES OF SMALL VOLUME

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Abstract. In this paper we discuss the approximate reconstruction of inhomogeneities of small volume. The data used for the reconstruction consist of boundary integrals of the (observed) electromagnetic fields. The numerical algorithms discussed are based on highly accurate asymptotic formulae for the electromagnetic fields in the presence of small volume inhomogeneities.

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1. INTRODUCTION

The problem of determining interior information about a medium from boundary field measurements is one that is not in general well posed. If, however, in advance we have additional structural information about the medium, then we may be able to determine specific features with “higher resolution”. As an example consider a bounded domain consisting of a homogeneously conducting background medium with volumetrically small inhomogeneities of conductivity different from the background. The geometry of each of the inhomogeneities may take the form of

$$\epsilon B + z,$$

where B is some fixed bounded domain containing the origin, or,

$$\{x + \eta h(x)n(x) : x \in \sigma_0, \eta \in (-\epsilon, \epsilon)\},$$

where σ_0 is a simple curve, $n(x)$ is a unit vectorfield orthogonal to σ_0 , and $h(x)$ is some smooth positive function. That is, we assume the inhomogeneities are either small domains of diameter $\approx \epsilon$, or they are thin strips of thickness $\approx \epsilon$ (this latter case only pertains to two dimensions). When in the shape of strips, the inhomogeneities are required to have bounded, nonzero conductivity, in order that the limiting ($\epsilon \rightarrow 0$) voltage potential coincide with that of the pure background medium.

In both of the settings described above, asymptotic formulae have been derived for the voltage potential as $\epsilon \rightarrow 0$. These formulae express the ϵ -dependent voltage potential in terms of the background potential, a certain

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Neumann-function, and the relevant polarization tensors [5, 8, 11]. In this paper we will use these formulae to develop an efficient algorithm to determine the location and/or shape of the inhomogeneities.

To be quite precise, let us assume we are given some finite set of small inhomogeneities ω_ϵ

$$\omega_\epsilon = \bigcup_{j=1}^m \omega_\epsilon^j$$

where, either all the inhomogeneities are of the form

$$\omega_\epsilon^j = \epsilon B_j + z_j,$$

for some smooth, bounded, star-shaped domains, B_j , or they are all of the form

$$\omega_\epsilon^j = \{x + \eta h^{(j)}(x)n(x) : x \in \sigma_0^j, \eta \in (-\epsilon, \epsilon)\},$$

for some simple curves, σ_0^j . Here $n(x)$ is a unit normal vectorfield to σ_0^j , and $h^{(j)}$ is some positive smooth function. The potential u_ϵ satisfies

$$\begin{aligned} \nabla \cdot \gamma_\epsilon \nabla u_\epsilon &= 0 & \text{in } \Omega \\ \gamma_\epsilon \frac{\partial u_\epsilon}{\partial \nu} &= g & \text{on } \partial\Omega \end{aligned} \tag{1.1}$$

with conductivity

$$\gamma_\epsilon = \begin{cases} \gamma_0 & x \in \Omega \setminus \overline{\omega_\epsilon}, \\ \gamma_j & x \in \omega_\epsilon^j. \end{cases} \tag{1.2}$$

We know that as ϵ approaches zero, the potential u_ϵ converges to the background potential, u , which satisfies

$$\begin{aligned} \Delta u &= 0 & \text{in } \Omega, \\ \gamma_0 \frac{\partial u}{\partial \nu} &= g & \text{on } \partial\Omega. \end{aligned} \tag{1.3}$$

In the case of m star-shaped imperfections of type $\epsilon B_j + z_j$ (of either finite or extreme conductivity) that are *well separated* from each other and from the boundary, it has been shown in [8, 11] that the following asymptotic formula holds:

$$u_\epsilon(y) = u(y) + \epsilon^n \sum_{j=1}^m \left(\frac{\gamma_0}{\gamma_j} - 1 \right) \nabla_x N(z_j, y) \cdot M^{(j)} \nabla u(z_j) + o(\epsilon^n). \tag{1.4}$$

$N(x, y)$ is the standard Neumann function

$$\begin{aligned} \Delta_x N(x, y) &= -\delta_y & \text{in } \Omega \\ \frac{\partial N}{\partial \nu_x}(x, y) &= -\frac{1}{|\partial\Omega|} & \text{on } \partial\Omega, \end{aligned} \tag{1.5}$$

and each $M^{(j)}$ is a $n \times n$, symmetric, positive definite matrix (called the polarization tensor). $M^{(j)}$ is given by

$$M_{kl}^{(j)} = |B_j| \delta_{kl} - \int_{\partial B_j} y_k \frac{\partial \phi_l^+}{\partial \nu} d\sigma_y, \tag{1.6}$$

where, for $0 < c_j = \frac{\gamma_0}{\gamma_j} < \infty$ and $1 \leq l \leq n$, $\phi_l(y)$ is the unique solution to

$$\begin{aligned} \Delta \phi_l &= 0 & \text{in } B_j, \\ \Delta \phi_l &= 0 & \text{in } \mathbb{R}^n \setminus \overline{B_j}, \\ c_j \frac{\partial \phi_l^+}{\partial \nu} - \frac{\partial \phi_l^-}{\partial \nu} &= (c_j - 1)\nu_l & \text{on } \partial B_j, \end{aligned} \tag{1.7}$$

with ϕ_l continuous across ∂B_j , and

$$\lim_{|y| \rightarrow \infty} \phi_l(y) = 0.$$

Note that the polarization tensor depends on the conductivity aspect ratio and the size and the shape of the inhomogeneity. Also note that the above asymptotic formula for u_ϵ holds in $\mathbb{R}^n$ for any $n \geq 2$ (later we shall specialize to two dimensions). It is sometimes convenient to emphasize the dependence of $M^{(j)}$ on the conductivity aspect ratio, $c_j = \gamma_0/\gamma_j$, in which case we write $M^{(j)}(c_j)$. The tensor $M^{(j)}$, introduced here, is the same as that introduced in [8], and is exactly c_j^{-1} times that introduced in [16].

There are very similar formulae for the (rescaled) polarization tensor in the case $\gamma_j = 0$ or $\gamma_j = \infty$, see [8] and [11]. The remainder, $o(\epsilon^n)$ in (1.4) signifies a term which goes to zero faster than ϵ^n uniformly in y , for y bounded away from the inhomogeneities. (It has been shown rigorously that this term is $O(\epsilon^{n+1/2})$, but numerical computations seem to suggest it goes to zero faster than this when the B_j are smooth.)

Remark 1. If (for fixed inhomogeneity B_j) we define $\varphi_l = -\phi_l + y_l$, then φ_l satisfies

$$\begin{aligned} \nabla_y \cdot \gamma(y) \nabla_y \varphi_l &= 0 & \text{in } \mathbb{R}^n, \\ \varphi_l - y_l &\rightarrow 0 & \text{as } |y| \rightarrow \infty. \end{aligned}$$

Here $\gamma(y)$ is given by

$$\gamma(y) = \begin{cases} \gamma_0 & y \in \mathbb{R}^n \setminus \overline{B_j}, \\ \gamma_j & x \in B_j. \end{cases} \tag{1.8}$$

A simple calculation now shows that

$$M_{kl}^{(j)} = c_j^{-1} \int_{B_j} \frac{\partial}{\partial y_k} \varphi_l(y) \, dy. \tag{1.9}$$

For $n = 2$ there is a very special relationship between the polarization tensor corresponding to aspect ratio c and that corresponding to aspect ratio c^{-1} . Based on the formula just above, and convex duality, we get that

$$M^{(j)}(c_j^{-1}) = c_j R M^{(j)}(c_j) R^T,$$

where R denotes the matrix corresponding to a counterclockwise rotation by $\pi/2$, *i.e.*, the matrix given by $R_{11} = R_{22} = 0$, $R_{12} = -R_{21} = -1$.

When the m conductivity imperfections ω_ϵ^j all take the form of the thin strips described above, and when these strips are *well separated* from each other and from the boundary, a similar asymptotic formula has been formally derived in [5]:

$$u_\epsilon(y) = u(y) - \epsilon \sum_{j=1}^m \int_{\sigma_0^j} \nabla_x N(x, y) \cdot A^{(j)}(x) \nabla u(x) h^{(j)}(x) ds_x + o(\epsilon). \tag{1.10}$$

The matrix $A^{(j)}(x)$ is similar to the polarization tensor $M^{(j)}$. If $\tau(x)$ and $n(x)$ are the unit tangential and normal fields to σ_0^j , then $A^{(j)}(x)$ is the symmetric matrix with eigenvectors $\tau(x)$ and $n(x)$, with respective eigenvalues $2(c_j^{-1} - 1)$ and $2(1 - c_j)$. Although this formula has not been proven rigorously, the computational evidence in [5] attests to its accuracy. We also note that this formula only pertains to two dimensions.

Our first goal is to apply these asymptotic formulae for the purpose of identifying the location and certain properties of the shapes of the inhomogeneities. In contrast to a (straightforward) boundary least squares fit to “the measured data”, we develop a method based on appropriate averaging, using particular “background” solutions as weights. The order ϵ^n term in the representation formula (1.4) is physically that of a voltage potential corresponding to m appropriately polarized dipoles located at the points z_j , $1 \leq j \leq m$. Similarly the order ϵ term in (1.10) is an “average” of voltage potentials corresponding to appropriately polarized dipoles distributed along the curve σ_0 . Results that address the theoretical feasibility of recovering the location of dipoles from boundary measurements are found various places in the literature; the reader may for example consult [14] and [12]. The special formulae found in [12] Section 3 for the recovery of a single inhomogeneity are very related to our brief discussion in Section 3 of the present paper.

Our second goal is to formally show that a similar “small volume” perturbation approach may be applied to the full Maxwell equations with inhomogeneities of different electric permittivity or magnetic permeability (or both). The formal arguments we present provide excellent insight when it comes to the inner workings of this “small volume” perturbation approach – an insight which may hopefully be used to adapt the approach for use in other contexts (for instance in the context of spectral data).

2. INTEGRATION AGAINST SPECIAL TEST FUNCTIONS

The general approach we take to reconstruct the locations and certain properties of the shapes of the inhomogeneities is to form the integral of “the measured boundary data” against harmonic test functions. In this section we use the formulae (1.4) and (1.10) to derive alternate expressions for these boundary integrals. These expressions are rather explicit in terms of the inhomogeneities and their locations. Suppose u_ϵ is the solution to (1.1) and w is some known harmonic function. We look for an expression for the “averaged measurement”

$$\Gamma = \int_{\partial\Omega} u_\epsilon \gamma_0 \frac{\partial w}{\partial \nu} d\sigma - \int_{\partial\Omega} \gamma_\epsilon \frac{\partial u_\epsilon}{\partial \nu} w d\sigma. \quad (2.11)$$

Integrals like Γ have already been used in several contexts to determine “interior” information from boundary data, most notably to determine the location of plane cracks, see for instance [2, 3] or [4]. If, in (2.11), we replace u_ϵ with the right hand side of (1.4), and $\gamma_\epsilon \frac{\partial u_\epsilon}{\partial \nu}$ with g , then we obtain

$$\Gamma = \int_{\partial\Omega} u \gamma_0 \frac{\partial w}{\partial \nu} d\sigma + \epsilon^n \sum_{j=1}^m \left(\frac{\gamma_0}{\gamma_j} - 1 \right) \int_{\partial\Omega} \nabla_x N(z_j, y) \cdot M^{(j)} \nabla u(z_j) \gamma_0 \frac{\partial w}{\partial \nu}(y) d\sigma_y - \int_{\partial\Omega} g w d\sigma + o(\epsilon^n).$$

The first and third terms cancel due to the harmonicity of u and w (and the fact that $\gamma_0 \frac{\partial u}{\partial \nu} = g$) and we are left with

$$\Gamma = \epsilon^n \sum_{j=1}^m \left(\frac{\gamma_0}{\gamma_j} - 1 \right) M^{(j)} \nabla u(z_j) \cdot \nabla_x \int_{\partial\Omega} N(z_j, y) \gamma_0 \frac{\partial w}{\partial \nu}(y) d\sigma_y + o(\epsilon^n).$$

By integration by parts and use of the equation for N we have

$$\begin{aligned} \int_{\partial\Omega} N(x, y) \frac{\partial w}{\partial \nu} d\sigma_y &= w(x) - \frac{1}{|\partial\Omega|} \int_{\partial\Omega} w d\sigma, \quad \text{or} \\ \nabla_x \int_{\partial\Omega} N(x, y) \gamma_0 \frac{\partial w}{\partial \nu} d\sigma_y &= \gamma_0 \nabla w(x). \end{aligned} \quad (2.12)$$

We have thus established the following theorem:

Theorem 1. *Let u_ϵ be the solution to (1.1) with inhomogeneities of type $\epsilon B_j + z_j$, with polarization tensors $M^{(j)}$. Let u be the background potential, satisfying (1.3). If w is any harmonic function, then the measurement Γ , defined by (2.11), has the expansion*

$$\Gamma = \epsilon^n \sum_{i=j}^m \gamma_0 \left(\frac{\gamma_0}{\gamma_j} - 1 \right) M^{(j)} \nabla u(z_j) \cdot \nabla w(z_j) + o(\epsilon^n).$$

The same argument can be applied to the case of strips. By replacing u_ϵ with the expansion (1.10) we obtain the approximation

$$\Gamma = -\epsilon \sum_{j=1}^m \int_{\partial\Omega} \int_{\sigma_0^j} \nabla_x N(x, y) A^{(j)}(x) \nabla u(x) h^{(j)}(x) \gamma_0 \frac{\partial w}{\partial \nu}(y) ds_x d\sigma_y + o(\epsilon),$$

which, by interchange of the order of integration, yields

$$\Gamma = -\epsilon \sum_{j=1}^m \int_{\sigma_0^j} A^{(j)}(x) \nabla u(x) \cdot \nabla_x \int_{\partial\Omega} N(x, y) \gamma_0 \frac{\partial w}{\partial \nu}(y) d\sigma_y h^{(j)}(x) ds_x + o(\epsilon).$$

Now, due to (2.12), we obtain

$$\Gamma = -\epsilon \sum_{j=1}^m \gamma_0 \int_{\sigma_0^j} \nabla w(x) \cdot A^{(j)}(x) \nabla u(x) h^{(j)}(x) ds_x + o(\epsilon).$$

We note that this last expansion has not been completely rigorously verified (due to the fact that the derivation of (1.10) is only formal).

To summarize: we have at this point at our disposal two asymptotic formulae with which to reconstruct the inhomogeneities; if the inhomogeneities are all of small diameter, then

$$\int_{\partial\Omega} u_\epsilon \gamma_0 \frac{\partial w}{\partial \nu} d\sigma - \int_{\partial\Omega} \gamma_\epsilon \frac{\partial u_\epsilon}{\partial \nu} w d\sigma \approx \epsilon^n \sum_{j=1}^m \gamma_0 \left(\frac{\gamma_0}{\gamma_j} - 1 \right) M^{(j)} \nabla u(z_j) \cdot \nabla w(z_j), \quad (2.13)$$

and if the inhomogeneities are all thin strips (in two dimensions) then

$$\int_{\partial\Omega} u_\epsilon \gamma_0 \frac{\partial w}{\partial \nu} d\sigma - \int_{\partial\Omega} \gamma_\epsilon \frac{\partial u_\epsilon}{\partial \nu} w d\sigma \approx -\epsilon \sum_{j=1}^m \gamma_0 \int_{\sigma_0^j} \nabla w(x) \cdot A^{(j)}(x) \nabla u(x) h^{(j)}(x) ds_x. \quad (2.14)$$

In both cases w is an arbitrary harmonic function.

Remark 2. Let us restrict attention to two dimensions. If we were to consider the simultaneous presence of thin inhomogeneities of thickness ϵ^2 and diametrically small inhomogeneities of diameter ϵ (*i.e.*, differently shaped inhomogeneities of same magnitude area) then a direct, additive combination of the formulae (2.13) and (2.14) gives the leading order perturbation of the “averaged data”. Perhaps more interesting, one might calculate a higher order version of (2.14) (including terms of order ϵ^2) and by combination with (2.13) now obtain an asymptotic perturbation for the “averaged data” in the simultaneous (well separated) presence of thin strips and diametrically small inhomogeneities, with thickness and diameter of same magnitude, ϵ .

Remark 3. If γ_ϵ is a small amplitude perturbation (an L^∞ -small perturbation) of the constant γ_0 , *i.e.*, if $\gamma_\epsilon(x) = \gamma_0 + \epsilon\gamma(x)$, with $\gamma \in L^\infty(\Omega)$, then it is not difficult to show that

$$\int_{\partial\Omega} u_\epsilon \gamma_0 \frac{\partial w}{\partial \nu} d\sigma - \int_{\partial\Omega} \gamma_\epsilon \frac{\partial u_\epsilon}{\partial \nu} w d\sigma = -\epsilon \int_{\Omega} \gamma(x) \nabla u_\epsilon \cdot \nabla w dx \approx -\epsilon \int_{\Omega} \gamma(x) \nabla u \cdot \nabla w dx, \quad (2.15)$$

for any harmonic function w . This approximation was for example used by Calderon [7]. The formulae (2.13) and (2.14) may be viewed as the analogues of (2.15) for special “low volume fraction perturbations”, that is for special types of L^1 -small perturbations.

3. DETECTING ONE INHOMOGENEITY

In this section we look at the simple case of only one inhomogeneity and show how formula (2.13) can be used to approximate its location and its polarization tensor. Suppose the inhomogeneity is centered at the point $p = (p_1, p_2)$, and has polarization tensor M . For simplicity we here denote the “rescaled” polarization tensor $\gamma_0 \left(\frac{\gamma_0}{\gamma_1} - 1 \right) M$ by M , and we consider only the case $n = 2$. Then, for any harmonic w

$$\int_{\partial\Omega} u_\epsilon \gamma_0 \frac{\partial w}{\partial \nu} d\sigma - \int_{\partial\Omega} \gamma_\epsilon \frac{\partial u_\epsilon}{\partial \nu} w d\sigma \approx \epsilon^2 M \nabla u(p) \cdot \nabla w(p). \quad (3.16)$$

Let us neglect the asymptotically small remainder term, and apply the currents ($g_i = \gamma_0 \nu_i$) corresponding to background potentials

$$u^{(i)} = x_i, \quad i = 1, 2.$$

First, we use the two test functions

$$w^{(1)} = x_1, \text{ and } w^{(2)} = x_2.$$

Define the “measurement” Γ_{ij} by

$$\Gamma_{ij} = \int_{\partial\Omega} u_\epsilon^{(i)} \gamma_0 \frac{\partial w^{(j)}}{\partial \nu} d\sigma - \int_{\partial\Omega} \gamma_\epsilon \frac{\partial u_\epsilon^{(i)}}{\partial \nu} w^{(j)} d\sigma.$$

From (3.16) we immediately get an approximation to the terms of the tensor $\epsilon^2 M$. In fact,

$$\epsilon^2 M_{ij} = \Gamma_{ij}.$$

Now that we have “recovered” $\epsilon^2 M$, we may use the same two currents, $\gamma_0 \nu_1$ and $\gamma_0 \nu_2$, to determine the location of the inhomogeneity. By inserting the test function

$$w^{(3)} = x_1 x_2$$

into (3.16), we obtain the linear system

$$\Gamma_{13} = \epsilon^2 (M_{11} p_2 + M_{21} p_1),$$

$$\Gamma_{23} = \epsilon^2 (M_{12} p_2 + M_{22} p_1),$$

which, since we already know $\epsilon^2 M$, is solvable for p_1 and p_2 (here we rely on the positive definiteness of M). By taking two boundary measurements we have thus directly obtained an approximation to the rescaled polarization tensor as well as the location of the inhomogeneity. The above formulas are in the same spirit as those derived in [2] for the identification of a single linear crack.

Remark 4. If there are more than one inhomogeneity but we know *a priori* that they all have the same rescaled polarization tensor, then we can easily determine the number of inhomogeneities by one boundary measurement. Simply apply boundary current $\gamma_0 \nu_1$ (so that $u = x_1$) and let $w = x_1$. Then (2.13) gives

$$\Gamma \approx \epsilon^2 m M_{11}$$

where Γ is measured. Assuming $\epsilon^2 M_{11}$ is known, we obtain m directly.

4. USING FOURIER INVERSION

In the previous section we saw an example of how the formula (2.13) can be used to directly calculate an approximation to the polarization tensor and the location of one inhomogeneity. By choosing more test functions w and possibly taking more boundary measurements, one may in a quite similar way calculate the locations and/or polarization tensors for several inhomogeneities. For a large number of inhomogeneities of different shape this becomes more difficult as the corresponding systems become highly nonlinear (and potentially, ill-posed). In this section we focus on an alternate (linear) method to determine the locations (and possibly, the shapes) of inhomogeneities of a wide range of types. This approach may require quite a number of boundary measurements, but if these are readily available, then the approach is rapid and simple to implement. This approach is based on the original idea of Calderon [7], which was, by way of the low amplitude perturbation formula (2.15), to reduce the reconstruction problem to the calculation of an inverse Fourier Transform.

We rewrite the formula (2.13)

$$\Gamma = \int_{\partial\Omega} u_\epsilon \gamma_0 \frac{\partial w}{\partial \nu} d\sigma - \int_{\partial\Omega} \gamma_\epsilon \frac{\partial u_\epsilon}{\partial \nu} w d\sigma \approx \epsilon^2 \sum_{i=1}^m M^{(i)} \nabla_x u(z_i) \cdot \nabla_x w(z_i), \quad (4.17)$$

by absorbing the constant $\gamma_0 \left(\frac{\gamma_0}{\gamma_j} - 1 \right)$ into the matrix $M^{(j)}$. For an arbitrary $\eta \in \mathbb{R}^2$, we assume we are in possession of the boundary data for the voltage potential, u_ϵ , whose corresponding background potential is given by

$$u = e^{i(\eta + i\eta^\perp) \cdot x}$$

(boundary current $= i(\eta + i\eta^\perp) \cdot \nu e^{i(\eta + i\eta^\perp) \cdot x}$) and we average this boundary data using the test function

$$w = e^{i(\eta - i\eta^\perp) \cdot x}.$$

Note that both of these functions are harmonic. Let us now view the measurement (4.17) as a function of η :

$$\Gamma(\eta) \approx \epsilon^2 \sum_{j=1}^m M^{(j)} \nabla \left(e^{i(\eta + i\eta^\perp) \cdot x} \right) \Big|_{x=z_j} \cdot \nabla \left(e^{i(\eta - i\eta^\perp) \cdot x} \right) \Big|_{x=z_j},$$

or

$$\Gamma(\eta) \approx -\epsilon^2 \sum_{j=1}^m M^{(j)} (\eta + i\eta^\perp) \cdot (\eta - i\eta^\perp) e^{2i\eta \cdot z_j}.$$

Recall that the function

$$e^{2i\eta \cdot z_j}$$

(up to a multiplicative constant) is exactly the Fourier Transform of the Dirac delta

$$\delta_{-2z_j}$$

(a point mass located at $-2z_j$). Multiplication by powers of η in Fourier space corresponds differentiation of the delta functions. The function $\Gamma(\eta)$ is therefore the Fourier Transform of a linear combination of derivatives of point masses, or

$$\check{\Gamma}(x) \approx \epsilon^2 \sum_{j=1}^m L_j \delta_{-2z_j},$$

where L_j is a second order constant coefficient, differential operator whose coefficients depend on the polarization tensor $M^{(j)}$, and $\check{\Gamma}(x)$ represent the inverse Fourier Transform of $\Gamma(\eta)$. The method of reconstruction we propose here consists in sampling the values of $\Gamma(\eta)$ at some discrete set of points and then calculating the corresponding discrete inverse Fourier Transform. After a rescaling (by $-1/2$) the support of this inverse Fourier Transform yields the location of the inhomogeneities. Once the locations are known we may calculate the polarization tensors by solving an appropriate linear system arising from (2.13).

Similarly, when we have m inhomogeneities in the form of thin strips

$$\Gamma(\eta) \approx \epsilon \sum_{j=1}^m \int_{\sigma_0^j} A^{(j)}(x)(\eta + i\eta^\perp) \cdot (\eta - i\eta^\perp) e^{i2\eta \cdot x} ds_x,$$

where γ_0 as well as $h^{(j)}(x)$ have been absorbed into $A^{(j)}(x)$. This is the Fourier Transform of a linear combination of derivatives of distributions with support on the curves $-2\sigma_0^j$, or

$$\check{\Gamma}(x) \approx \epsilon \sum_{j=1}^m L_j D_{\sigma_0^j},$$

where $\check{\Gamma}(x)$ denotes the inverse Fourier Transform of $\Gamma(\eta)$, the L_j are second order constant coefficient differential operators, and $D_{\sigma_0^j}$ are distributions supported on $-2\sigma_0^j$. In this case we also propose to calculate a discrete inverse Fourier Transform of $\Gamma(\eta)$ to approximate the locations and shapes of the curves σ_0^j .

Calderon's analysis (*cf.* [7]) showed that the inverse Fourier Transform of $\Gamma(\eta)$ yielded a good approximation to small amplitude perturbations in the conductivity (around a constant). The calculations above shows that it also yields a good approximation to (large amplitude but) small volume fraction perturbations of the conductivity. Finally we note that it is possible to construct exponential solutions, with properties similar to those above, for use in more complicated situations, for instance in connection with the Maxwell equations (see [17]).

5. NUMERICAL RESULTS

In the last section we saw that the boundary measurement

$$\Gamma(\eta) = \int_{\partial\Omega} u_\epsilon \gamma_0 \frac{\partial w}{\partial \nu} d\sigma - \int_{\partial\Omega} \gamma_\epsilon \frac{\partial u_\epsilon}{\partial \nu} w d\sigma,$$

corresponding to background potential

$$u = e^{i(\eta + i\eta^\perp) \cdot x},$$

and test function

$$w = e^{i(\eta - i\eta^\perp) \cdot x}$$

is approximately (up to rescalings) the Fourier Transform of a distribution with support at the centers of the inhomogeneities. We now consider the computational aspects of performing a discrete inverse Fourier Transform in order to identify the location of the inhomogeneities. Since the accuracy of the approximations (1.4, 1.10) (and thus the accuracy of the approximations (2.13, 2.14)) has already been carefully documented the most important computational issue that remains is the conditioning of the discrete inverse Fourier Transform, when applied to data that arise from distributions that are supported at a finite set of points or on a finite set of simple

curves. Again, reflecting the fact that the accuracy of the approximations (2.13) and (2.14) are not at issue, we shall use these two formulas to generate the data for $\Gamma(\eta)$ (as opposed to computing it from approximate solutions to the elliptic boundary value problems). We do want to emphasize that before one could contemplate to use this algorithm “in real life” there is of course one important practical issue, which we do not attempt to address, and which needs to be resolved. We are here referring to the issue of how to obtain a sufficient number of real measurements to generate the values of $\Gamma(\eta)$ needed for a satisfactory Fourier Inversion.

To arrive at some idea of the number of data (sampling) points needed for an accurate discrete Fourier inversion we remind the reader of the main assertion of the so-called Shannon’s Sampling theorem (*cf.* [9]): “a function f is completely specified (by a very explicit formula) by the sampled values $\{f(\frac{n}{2M} + c_0)\}_{n=-\infty}^{\infty}$ if and only if the support of $\hat{f}$ is contained inside a square of side $2M$ ”. For our application this suggests two things: (1) if the inhomogeneities are contained inside a square of side $2M$, then we need to sample $\Gamma(\eta)$ at a uniform, infinite, rectangular grid of meshsize $\frac{1}{2M}$ to obtain an accurate reconstruction, (2) if we only sample the points in this grid for which the absolute values of the coordinates are less than K , then the resulting discrete inverse Fourier Transform will recover the location of the inhomogeneities with a resolution of order $\delta = \frac{1}{2K}$. In summary: we need (conservatively) of the order $(2M/\delta)^2$ sampled values of $\Gamma(\eta)$ to reconstruct, with resolution δ , a collection of inhomogeneities that lie inside a square of side $2M$.

It may be useful to compare these measurement counts to those related to another method (a linear “linear sampling” method) which has recently been proposed as a tool for the reconstruction of collections of small inhomogeneities (see [6]). For the accurate reconstruction of the location of m inhomogeneities this method requires knowledge of the subspace spanned by the first $2m$ eigenvectors of the incremental Dirichlet to Neumann data operator. For moderate size m it is reasonable to assume that we are able to approximate this subspace (so as to determine the location of the inhomogeneities with a resolution of order δ) by $C_\delta m$ measurements (using a “power-like” method) where C_δ depends on δ and (the gaps in) the spectrum of the incremental Dirichlet to Neumann data operator. For a relatively small number of inhomogeneities ($m \ll (2M/\delta)^2/C_\delta$) the “linear sampling” method would therefore require fewer measurements than the Fourier Inversion based approach. For relatively large number of inhomogeneities the count should be in favor of the Fourier Inversion based approach. We note that the largest possible number of inhomogeneities that are well separated (by δ) is *a priori* bounded by $(2M/\delta)^2$. A more precise comparison of measurement counts would require a careful analysis of the actual size of C_δ and the range of validity of the bound $C_\delta m$. We also note that the present crude comparison may substantially overestimate the number of measurements needed for the Fourier Inversion based approach in the case of only a few inhomogeneities. Some of the following numerical examples clearly indicate that fewer measurements are sufficient to reconstruct the location of a very small number of inhomogeneities.

As our first example, we take the domain Ω to be the square $[-10, 10] \times [-10, 10]$ and we insert five inhomogeneities in the shape of balls, with the j -th ball positioned at the point (j, j) . In the case of balls the polarization tensor is just a multiple of the identity, $M^{(j)} = d_j I$, where d_j depends on the aspect ratio γ_0/γ_j and the relative size of the ball [8]. Here we take each d_j to be 10 and $\epsilon = 0.1$. We sample our approximation of $\Gamma(\eta)$ on the square $[-3, 3] \times [-3, 3]$ with a uniform 30×30 grid. We are thus following the recipe from above, with $M = 2.5$ and $K = 3$. We should expect a recovery of all the locations of the inhomogeneities, with a resolution $\delta = 1/6$. The discrete inverse Fourier Transform yields the grey-level (intensity) plot shown in Figure 1.

Note, though, that the reconstruction in Figure 1 corresponds to sampling $\Gamma(\eta)$ at 900 points. Next we sample the same function on $[-2.5, 2.5] \times [-2.5, 2.5]$ with a uniform 20×20 grid, thus having a total of 400 sample points. In Figure 2 we see that the five balls are still visible. If the domain has only one inhomogeneity, then we can find its location by sampling far fewer values of Γ . The reconstruction shown in Figure 3 corresponds to a single inhomogeneity at the point $(1, 1)$. We have sampled the values of Γ on the square $[-1.5, 1.5] \times [-1.5, 1.5]$ using a uniform 4×4 grid.

In order to simulate errors in the boundary measurements, as well as the error inherent in the approximations (2.13, 2.14) we add on the order of 10% of random noise to the values of Γ in the examples with five inhomogeneities. We see from Figures 4 and 5 that the reconstruction is quite stable.

FIGURE 1. Five inhomogeneities -30×30 sample points. FIGURE 2. Five inhomogeneities -20×20 sample points.

For our next example we take the case in which there are 50 equidistant balls (of same size and aspect ratio as before) lying on the diagonal between $(-5, 5)$ and $(5, 5)$. We sample the approximation to $\Gamma(\eta)$ (with 10% noise added) on the square $[-10, 10] \times [-10, 10]$ using a uniform 128×128 grid. From Figure 6 see that the placement of the 50 balls is clearly defined, but on the other hand we can not identify individual balls. We note that, due to the closeness of the balls, it is not clear how good an approximation to Γ the expression (2.13) provides. We also note that “background” noise is more visible here than in any of the previous plots: this is entirely to be expected, since the fairly high number of balls leads to much larger values for the “averaged data” and thus to much larger absolute noise. The “maximal intensity” near the location of the balls is also

FIGURE 3. One inhomogeneity -4×4 sample points. FIGURE 4. Five inhomogeneities -30×30 sample points with 10% noise.

significantly larger than before; this may be a little hard to appreciate on a grey-level plot, it is much easier to see on the corresponding surface plot, shown in Figure 7.

In our last example we take an inhomogeneity in the shape of a thin strip, *i.e.* we use the approximation to $\Gamma(\eta)$ arising from (2.14). The strip is situated about the interval $[0, 1]$ on the x -axis. We choose the background conductivity to be 1 and the conductivity on the strip to be 10, and take $\epsilon = 0.1$. We sample the approximation to $\Gamma(\eta)$ with (10% noise added) on the square $[-3, 3] \times [-3, 3]$ using a uniform 10×10 grid. The position of the strip is clearly visible in Figure 8.

The numerical experiments we have presented here correspond to but one of many possible choices for the background voltages and the test functions in the formulae (2.13) and (2.14). We are currently exploring other

FIGURE 5. Five inhomogeneities -20×20 sample points with 10% noise. FIGURE 6. Fifty inhomogeneities -128×128 sample points with 10% noise.

choices, some of which are undoubtedly more practical, as far as the required number of measurements are concerned. We are convinced that the use of approximate formulae such as (2.13) and (2.14) represents a very promising approach to the reconstruction problem, in particular when one considers reconstructions based on the use of electric as well as magnetic fields. This is the main reason why we devote the next section to the (formal) derivation of a formula analogous to (2.13) for solutions to the Maxwell equations.

FIGURE 7. Fifty inhomogeneities -128×128 sample points with 10% noise.

FIGURE 8. Strip -10×10 sample points with 10% noise.

6. IDENTIFICATION OF SMALL INHOMOGENEITIES USING ELECTROMAGNETIC FIELDS

In this section we establish the basis for the extension of our reconstruction approach to take into account all electromagnetic fields. To be more precise: we give a formal derivation of an identity similar to (2.13) for solutions to the 3-d Maxwell equations (the identity (6.32)). This formal derivation provides an extremely intuitive and accessible alternative to the rigorous, at times quite technical, derivation found in [1]. We strongly feel that the additional insight provided by this formal derivation will allow the discovery of similar identities in other contexts.

Let Ω be a bounded subset of $\mathbb{R}^3$, with smooth boundary $\partial\Omega$. Let ν denote the outward unit normal to $\partial\Omega$, and let (k_1, k_2, k_3) be an orthonormal basis of $\mathbb{R}^3$. For fixed frequency $\omega > 0$, (E_ϵ, H_ϵ) denote time-harmonic electromagnetic fields. These time-harmonic fields satisfy the following form of the Maxwell equations

$$\begin{aligned} \nabla \times E_\epsilon &= i\omega\mu_\epsilon H_\epsilon & \text{in } \Omega, \\ \nabla \times H_\epsilon &= -i\omega q_\epsilon E_\epsilon & \text{in } \Omega, \\ E_\epsilon \times \nu &= f & \text{on } \partial\Omega. \end{aligned} \quad (6.18)$$

Here f is a tangential field on the boundary $\partial\Omega$ that furthermore belongs to the Sobolev space $H_{\text{div}}^{-1/2}(\partial\Omega) = \{g \in H^{-1/2}(\partial\Omega)^3, g \cdot \nu = 0 \text{ on } \partial\Omega, \text{div}_{\partial\Omega}(g) \in H^{-1/2}(\partial\Omega)\}$. The magnetic permeability μ_ϵ and the electric permittivity q_ϵ differ from the (background) values μ_0 and q_0 inside the small inhomogeneity $B_\epsilon = \epsilon B + z$. In other words

$$\mu_\epsilon = \begin{cases} \mu_0 & \text{in } \Omega \setminus \overline{B_\epsilon}, \\ \mu_1 & \text{in } B_\epsilon, \end{cases}$$

and

$$q_\epsilon = \begin{cases} q_0 & \text{in } \Omega \setminus \overline{B_\epsilon}, \\ q_1 & \text{in } B_\epsilon. \end{cases}$$

B is a bounded smooth domain, containing the origin; μ_j and q_j are constants with $\mu_j > 0$, $\text{Re } q_j > 0$, and $\text{Im } q_j \geq 0$. For simplicity we only consider one inhomogeneity B_ϵ . The general case (with a finite number of well separated inhomogeneities) does not lead to any new difficulties. We assume that ω is not an eigenvalue for the problem (6.18), and so by assumption there exists a unique solution $(E_\epsilon, H_\epsilon) \in [H(\text{curl}, \Omega)]^2 = \{F \in L^2(\Omega)^3, \nabla \times F \in L^2(\Omega)^3\}^2$. From (6.18) we have that

$$\nabla \cdot (q_\epsilon E_\epsilon) = \nabla \cdot (\mu_\epsilon H_\epsilon) = 0 \quad \text{in } \Omega. \quad (6.19)$$

From [15] and [13] we know that the boundary map given by

$$\Lambda : E_\epsilon \times \nu|_{\partial\Omega} \mapsto H_\epsilon \times \nu|_{\partial\Omega},$$

where E_ϵ, H_ϵ satisfy (6.18) determines the functions μ_ϵ and q_ϵ , in the case when these are sufficiently smooth. The goal in the piecewise constant context, considered here, is to construct an efficient method for determining the location of B_ϵ , and certain aspects of B from only one pair $(E_\epsilon \times \nu, H_\epsilon \times \nu)$.

Let v be a smooth solution to

$$\nabla \times (\nabla \times v) - \omega^2 \mu_0 q_0 v = 0. \quad (6.20)$$

From the Maxwell equations (6.18) we obtain by integration by parts

$$\begin{aligned} \int_{\partial\Omega} (\nabla \times E_\epsilon) \times \nu \cdot v \, d\sigma - \int_{\partial\Omega} (\nabla \times v) \times \nu \cdot E_\epsilon \, d\sigma &= \omega^2 \mu_0 (q_0 - q_1) \int_{B_\epsilon} E_\epsilon \cdot v \, dx \\ &+ \left(\frac{\mu_0}{\mu_1} - 1 \right) \int_{B_\epsilon} \nabla \times E_\epsilon \cdot \nabla \times v \, dx. \end{aligned} \quad (6.21)$$

If we replace v and $\nabla \times v$ by $v(z)$ and $\nabla \times v(z)$ in B_ϵ we formally obtain

$$\begin{aligned} \int_{\partial\Omega} (\nabla \times E_\epsilon) \times \nu \cdot v \, d\sigma - \int_{\partial\Omega} (\nabla \times v) \times \nu \cdot E_\epsilon \, d\sigma &= \omega^2 \mu_0 (q_0 - q_1) \left(\int_{B_\epsilon} E_\epsilon \, dx \right) \cdot v(z) \\ &+ \left(\frac{\mu_0}{\mu_1} - 1 \right) \left(\int_{B_\epsilon} \nabla \times E_\epsilon \, dx \right) \cdot \nabla \times v(z) + o(\epsilon^3). \end{aligned} \quad (6.22)$$

We note that it is quite easy, systematically to create vector valued functions v , that satisfy (6.20). For example, let $\xi \in \mathbb{R}^3$ be such that $\xi \cdot \xi = \omega^2 \mu_0 q_0$, let k be an arbitrary vector in $\mathbb{R}^3$, and define

$$v_{\xi,k}(x) = e^{i\xi \cdot x} k + \frac{1}{\omega^2 \mu_0 q_0} \nabla \nabla \cdot (e^{i\xi \cdot x} k) = e^{i\xi \cdot x} \left(k - \frac{k \cdot \xi}{\omega^2 \mu_0 q_0} \xi \right). \quad (6.23)$$

In order to derive a formula that is similar to (2.13) we first derive an asymptotic expansion for E_ϵ inside B_ϵ , which we then insert into (6.22).

6.1. The asymptotic expansion for E_ϵ

Following a common practice in multiscale expansions we introduce the local variable $y = \frac{x-z}{\epsilon}$, where z is the “center” of B_ϵ . Let (E_0, H_0) be the solution to the following (background) Maxwell Equations

$$\begin{aligned} \nabla \times E_0 &= i\omega \mu_0 H_0 & \text{in } \Omega, \\ \nabla \times H_0 &= -i\omega q_0 E_0 & \text{in } \Omega, \\ E_0 \times \nu &= f & \text{on } \partial\Omega. \end{aligned} \quad (6.24)$$

We expect that $E_\epsilon(x)$ will differ appreciably from $E_0(x)$ for x near z , but that it will differ little from $E_0(x)$ for x far from z . Therefore, in the spirit of matched asymptotic expansions, we shall represent $E_\epsilon(x)$ by two different expansions, an inner expansion for x near z , and an outer expansion for x far from z . The outer expansion must begin with E_0 , so we write

$$E_\epsilon(x) = E_0(x) + \epsilon^{\alpha_1} E_1(x) + \epsilon^{\alpha_2} E_2(x) + \dots, \quad \text{for } |x-z| \gg O(\epsilon), \quad (6.25)$$

where $0 < \alpha_1 < \alpha_2 < \dots$, and $E_1, E_2, \dots$ are to be found. We write the inner expansion as

$$E_\epsilon(z + \epsilon y) = e_\epsilon(y) = e_0(y) + \epsilon e_1(y) + \epsilon^2 e_2(y) + \dots \quad \text{for } |y| = O(1), \quad (6.26)$$

where $e_0, e_1, \dots$ are to be found. The magnetic field H_ϵ has similar expansions. The inner and outer expansions must be asymptotically equal in some “overlap” domain within which the stretched variable y is large and $x-z$ is small. In this domain the matching condition is

$$E_0(x) + \epsilon^{\alpha_1} E_1(x) + \epsilon^{\alpha_2} E_2(x) + \dots \sim e_0(y) + \epsilon e_1(y) + \epsilon^2 e_2(y) + \dots \quad (6.27)$$

From the terms of order ϵ^0 , we obtain the first matching condition

$$e_0(y) \rightarrow E_0(z) \quad \text{as } |y| \rightarrow +\infty.$$

If we substitute (6.26) into (6.18, 6.19) and formally equate coefficients of ϵ^{-1} we get

$$\nabla_y \times e_0 = 0, \quad \nabla_y \cdot (q e_0) = 0 \quad \text{in } \mathbb{R}^3,$$

where

$$q = \begin{cases} q_0 & \text{in } \mathbb{R}^3 \setminus \overline{B}, \\ q_1 & \text{in } B. \end{cases}$$

Therefore,

$$e_0(y) = \sum_{i=1}^3 (E_0(z) \cdot k_i) \nabla_y \varphi_i(y), \quad (6.28)$$

where φ_i is the unique solution to the following problem

$$\begin{aligned} \nabla_y \cdot (q(y) \nabla_y \varphi_i) &= 0 \quad \text{in } \mathbb{R}^3, \\ \varphi_i(y) - y_i &\rightarrow 0, \quad \text{as } |y| \rightarrow +\infty. \end{aligned} \quad (6.29)$$

The asymptotic behavior of $\nabla_y \varphi_i(y)$ as $|y| \rightarrow +\infty$ (see [10]) is

$$\nabla_y \varphi_i(y) = k_i + \frac{C}{|y|^3} y + o\left(\frac{1}{|y|^2}\right),$$

where C is a constant that depends on q_0, q_1 , and B . We note that the coefficients q_j are complex, but the fact that $\operatorname{Re} q_j > 0$ is sufficient to guarantee the uniqueness and the decay mentioned above.

Exactly as for the electric field, the first order term in the asymptotic expansion of $\frac{1}{i\omega\mu_1} \nabla \times E_\epsilon = H_\epsilon$ inside B_ϵ is given by

$$h_0(y) = \sum_{i=1}^3 (H_0(z) \cdot k_i) \nabla_y \psi_i(y), \quad (6.30)$$

where ψ_i is the unique solution to the following problem

$$\begin{aligned} \nabla_y \cdot (\mu(y) \nabla_y \psi_i) &= 0 \quad \text{in } \mathbb{R}^3, \\ \psi_i(y) - y_i &\rightarrow 0, \quad \text{as } |y| \rightarrow +\infty, \end{aligned} \quad (6.31)$$

with

$$\mu = \begin{cases} \mu_0 & \text{in } \mathbb{R}^3 \setminus \overline{B}, \\ \mu_1 & \text{in } B. \end{cases}$$

Based on Remark 1.7 we immediately get that

$$\frac{q_1}{q_0} \int_B \frac{\partial}{\partial y_j} \varphi_l(y) \, dy = |B| \delta_{jl} - \int_{\partial B} y_j \frac{\partial \phi_l^+}{\partial \nu} \, d\sigma_y = M_{jl} \left(\frac{q_0}{q_1} \right),$$

where the tensor $M(\frac{q_0}{q_1})$ has been appropriately extended to complex aspect ratios. Similarly we get

$$\frac{\mu_1}{\mu_0} \int_B \frac{\partial}{\partial y_j} \psi_l(y) \, dy = M_{jl} \left(\frac{\mu_0}{\mu_1} \right).$$

6.2. The asymptotic formula for the measured data

By combining (6.22) and the (inner) asymptotic expressions (6.28, 6.30) we arrive at the following formula.

$$\begin{aligned}
& \int_{\partial\Omega} (\nabla \times E_\epsilon) \times \nu \cdot v \, d\sigma - \int_{\partial\Omega} (\nabla \times v) \times \nu \cdot \nu \times (E_\epsilon \times \nu) \, d\sigma \\
&= \epsilon^3 \omega^2 \mu_0 (q_0 - q_1) \sum_{i=1}^3 (E_0(z) \cdot k_i) \left(\int_B \nabla_y \varphi_i(y) \, dy \right) \cdot v(z) \\
&+ \left(\frac{\mu_0}{\mu_1} - 1 \right) i\omega \mu_1 \sum_{i=1}^3 (H_0(z) \cdot k_i) \left(\int_B \nabla_y \psi_i(y) \, dy \right) \cdot \nabla \times v(z) \\
&+ o(\epsilon^3) \\
&= \epsilon^3 \left[\omega^2 \mu_0 q_0 \left(\frac{q_0}{q_1} - 1 \right) M \left(\frac{q_0}{q_1} \right) E_0(z) \cdot v(z) \right. \\
&\quad \left. + i\omega \mu_0 \left(\frac{\mu_0}{\mu_1} - 1 \right) M \left(\frac{\mu_0}{\mu_1} \right) H_0(z) \cdot \nabla \times v(z) \right] + o(\epsilon^3). \tag{6.32}
\end{aligned}$$

Here (E_ϵ, H_ϵ) is the unique solution to (6.18) and (E_0, H_0) is the unique solution to the unperturbed problem (6.24). The functions φ_i and ψ_i are defined by (6.29) and (6.31). To proceed from (6.22) to (6.32) we have also used that $\nu \times (E_\epsilon \times \nu)$ is the projection of E_ϵ onto the tangent plane of $\partial\Omega$. For a rigorous derivation of this formula (and other asymptotic representations) we refer the reader to [1]. We may now insert the functions $v_{\xi,k}$ (or other solutions to (6.20) into (6.32)) to arrive at reconstruction algorithms related to those described in Sections 3–5.

Remark 5. Note that if the inhomogeneity $\epsilon B + z$ (and the domain Ω) is cylindrical (in the x_3 direction) then we have $\psi_3 = \varphi_3 = y_3$. Furthermore, the (correctly interpreted) functions $\psi_1, \psi_2, \varphi_1$ and φ_2 depend only on the variables (y_1, y_2) , and thus the polarization tensor $\epsilon^3 M(c)$ gets replaced by

$$\epsilon^2 M(c) = \epsilon^2 \begin{bmatrix} \{M_{\alpha,\beta}(c)\}_{\alpha,\beta=1}^2 & 0 \\ 0 & c^{-1}|B'| \end{bmatrix}.$$

If we now consider the TE (transverse electric) case – that is the case of an electric field $E = (0, 0, E(x_1, x_2))$ with corresponding magnetic field $H = \frac{1}{i\omega\mu}(\frac{\partial}{\partial x_2}E, -\frac{\partial}{\partial x_1}E, 0)$ – and take $v = (0, 0, v(x_1, x_2))$ (with $\triangle v + \omega^2 \mu_0 q_0 v = 0$) then the analogue of (6.32) reads

$$\begin{aligned}
& \int_{\partial\Omega'} \frac{\partial E_\epsilon}{\partial \nu} v \, d\sigma - \int_{\partial\Omega'} \frac{\partial v}{\partial \nu} E_\epsilon \, d\sigma = \epsilon^2 \left[\omega^2 \mu_0 (q_0 - q_1) |B'| E_0(z') v(z') \right. \\
&\quad \left. + \left(\frac{\mu_0}{\mu_1} - 1 \right) M \left(\frac{\mu_0}{\mu_1} \right) R^T \nabla E_0(z') \cdot R^T \nabla v(z') \right] + o(\epsilon^2) \\
&= \epsilon^2 \left[\omega^2 \mu_0 (q_0 - q_1) |B'| E_0(z') v(z') \right. \\
&\quad \left. + \left(\frac{\mu_0}{\mu_1} - 1 \right) \frac{\mu_1}{\mu_0} M \left(\frac{\mu_1}{\mu_0} \right) \nabla E_0(z') \cdot \nabla v(z') \right] + o(\epsilon^2) \tag{6.33}
\end{aligned}$$

where B' , Ω' and z' denote the two dimensional base domains (or point) of B , Ω and z , respectively. In the last identity we used the second observation from Remark 1.7. The formula (6.33) is closely related to the boundary representations, rigorously derived in [16]. Indeed, the formula (6.33) may be obtained by appropriate integration of a slight variation of the representation given in Theorem 1 of [16].

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