

Analysis of a Residual Model for the Decomposition of Polarimetric SAR Data

Xiaojing Bai, Binbin He^{*}, Xiaowen Li

School of Resources and Environment, University of Electronic Science and Technology of China, Chengdu, China Email: *binbinhe@uestc.edu.cn, 449196341@qq.com, 452672784@qq.com

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ABSTRACT

Accurate estimation of the double-bounce scattering f_d and surface scattering f_s coefficients with Freeman-Durden decomposition is still difficult. This difficulty arises because overestimation of the volume scattering energy contribution P_v leads to negative values for f_d and f_s . A generalized residual model is introduced to estimate f_d and f_s . The relationship between P_v and the residual model is analyzed. Eigenvalues computed from the residual model must be positive to explain physical scattering mechanisms. The authors employ a new volumetric scattering model to minimize P_v as calculated by several decomposition methods. It is concluded that decreasing P_v can help reduce negative energy. This conclusion is validated using actual polarimetric SAR data.

Keywords: Freeman-Durden Decomposition; Polarimetric SAR; Residual Model; Negative Scattering Energy

1. Introduction

Among the incoherent target decompositions for polarimetric SAR (PolSAR) data, Freeman-Durden decomposition (FDD), which is based on a physical model, consists of three types of scattering mechanisms: volumetric scattering, double-bounce scattering, and surface scattering [1]. FDD has been used for PolSAR data processing, such as speckle filtering [2], image classification [3] and soil moisture estimation [4].

However, there are three concerns about FDD model: the assumption of reflection symmetry in the azimuth, the elimination of the negative scattering power contribution from f_d and f_s , and the selection of an appropriate volumetric scattering model $\langle [C] \rangle_v$. In urban settings, azimuthally symmetric reflection occurs inferquently. To remove the assumption that the reflection is symmetric, a helix was added as the fourth scattering component in Yamaguchi decomposition (YD), and similar terms (volumetric scattering, double-bounce scattering, and surface scattering) were used to those in the FDD model [5].

A hybrid decomposition method was proposed by van Zyl *et al.* to overcome negative scattering energy [6]. Orientation angle compensation was employed by An *et al.* [7] and Lee and Ainsworth [8] to eliminate the nega-

tive scattering energy contribution. Subsequently, an adaptive model-based decomposition was proposed, which extended the method proposed by van Zyl *et al.* [9]. A hybrid Freeman/eigenvalues technique was used by Cloude for avoiding negative power problems [10]. Considering actual tree trunk and branch distribution, Yamaguchi *et al.* slightly modified the dipole model used in FDD model [5]. A unit matrix adopted as the new volumetric scattering model was introduced into the FDD method (abbreviated as UmFDD), and the major advantage of this new model was that it characterized the total randomness that was the prime feature of volume scattering [7].

In this paper, the authors attempt to adopt a new volumetric scattering model to minimize volume scattering energy for eliminating negative coefficients of doublebounce and surface scattering. Firstly, the authors review three decomposition methods: FDD, YD, and UmFDD. A generalized residual model is introduced to link these methods. Then, the authors analyze the relationship between the volume scattering energy contribution P_v and the residual model. The eigenvalues are extracted to analyze whether the generalized residual model can be used to interpret physical scattering mechanisms. The volume scattering contribution P_v is minimized by means of a new model, which confirms that decreasing the volumetric scattering contribution P_v can help re-

^{*}Corresponding author.

duce negative energy and gives the residual model an ability to explain physical scattering mechanisms.

2. Residual Model Extraction from Model-Based Decomposition

2.1. Residual Model Extracted from FDD Method

When the reciprocity assumption is valid for PolSAR imagery, each pixel of the image can be represented by a 3×3 covariance matrix $\langle [C] \rangle$,

$$\langle [C] \rangle$$

$$= \begin{bmatrix} \langle |S_{\rm HH}|^2 \rangle & \sqrt{2} \langle S_{\rm HH} S_{\rm HV}^* \rangle & \langle S_{\rm HH} S_{\rm HV}^* \rangle \\ \sqrt{2} \langle S_{\rm HV} S_{\rm HH}^* \rangle & 2 \langle |S_{\rm HV}|^2 \rangle & \sqrt{2} \langle S_{\rm HV} S_{\rm VV}^* \rangle \\ \langle S_{\rm VV} S_{\rm HH}^* \rangle & \sqrt{2} \langle S_{\rm VV} S_{\rm HV}^* \rangle & \langle |S_{\rm VV}|^2 \rangle \end{bmatrix}$$
(1)

where * denotes a conjugate operator and $\langle \cdot \rangle$ an assembled average. $\langle [C] \rangle$ is a positive semi-definite Hermitian matrix.

When a target is azimuthally symmetric, the reflection can be considered to be symmetric, and thus, the crossproducts of co- and cross-polarized returns are zero. Thus, (1) becomes,

$$\langle [\boldsymbol{C}] \rangle = \begin{bmatrix} \langle |\boldsymbol{S}_{\rm HH}|^2 \rangle & 0 & \langle \boldsymbol{S}_{\rm HH} \boldsymbol{S}_{\rm HV}^* \rangle \\ 0 & 2 \langle |\boldsymbol{S}_{\rm HV}|^2 \rangle & 0 \\ \langle \boldsymbol{S}_{\rm VV} \boldsymbol{S}_{\rm HH}^* \rangle & 0 & \langle |\boldsymbol{S}_{\rm VV}|^2 \rangle \end{bmatrix}$$
(2)

Freeman and Durden [1] decomposed the covariance matrix $\langle [C] \rangle$ in (2) as the sum of three components representing scattering mechanisms of volumetric scattering $f_v \langle [C] \rangle_v$, double-bounce scattering $f_d \langle [C] \rangle_d$, and surface scattering $f_s \langle [C] \rangle_s$:

$$\langle [\boldsymbol{C}] \rangle = f_{v} \langle [\boldsymbol{C}] \rangle_{v} + f_{d} \langle [\boldsymbol{C}] \rangle_{d} + f_{s} \langle [\boldsymbol{C}] \rangle_{s}$$
 (3)

where f_v, f_d , and f_s are scattering coefficients denoting the energy contribution for each scattering mechanism, respectively.

In FDD, the covariance matrix of the volumetric scattering component $\langle [C] \rangle_{v}$ is modeled as the scattering from a cloud of randomly oriented dipoles or long thin cylinders,

$$\langle [C] \rangle_{v} = \begin{bmatrix} 1 & 0 & 1/3 \\ 0 & 2/3 & 0 \\ 1/3 & 0 & 1 \end{bmatrix}$$
 (4)

 $\langle [C] \rangle_{d}$ is modeled as the scattering from a dihedral corner reflector,

$$\left\langle \begin{bmatrix} \boldsymbol{C} \end{bmatrix} \right\rangle_{\mathrm{d}} = \begin{bmatrix} \left| \boldsymbol{\alpha} \right|^2 & 0 & \boldsymbol{\alpha} \\ 0 & 0 & 0 \\ \boldsymbol{\alpha}^* & 0 & 1 \end{bmatrix}$$
(5)

and $\langle [C] \rangle_{s}$ is modeled as a first-order Bragg scattering,

$$\left\langle \left[\boldsymbol{C} \right] \right\rangle_{\mathrm{s}} = \begin{bmatrix} \left| \boldsymbol{\beta} \right|^2 & 0 & \boldsymbol{\beta} \\ 0 & 0 & 0 \\ \boldsymbol{\beta}^* & 0 & 1 \end{bmatrix}$$
(6)

The coefficient f_v can be directly computed from (3) to be,

$$f_{\rm v} = 3 \left\langle \left| \boldsymbol{S}_{\rm HV} \right|^2 \right\rangle \tag{7}$$

Then, the energy contribution P_v corresponding to the volumetric scattering component is,

$$P_{\nu} = \frac{8f_{\nu}}{3} = 8\left< \left| \boldsymbol{S}_{\rm HV} \right|^2 \right>$$

Inserting (4)-(6) into (3), and using the expression of f_v in (7) and $\langle [C] \rangle$ in (2), one can algebraically express the following equations:

$$\begin{cases} \left\langle \left| \boldsymbol{S}_{\mathrm{HH}} \right|^{2} \right\rangle - f_{\mathrm{v}} = f_{\mathrm{s}} \left| \boldsymbol{\beta} \right|^{2} + f_{\mathrm{d}} \left| \boldsymbol{\alpha} \right|^{2} \\ \left\langle \left| \boldsymbol{S}_{\mathrm{VV}} \right|^{2} \right\rangle - f_{\mathrm{v}} = f_{\mathrm{s}} + f_{\mathrm{d}} \\ \left\langle \boldsymbol{S}_{\mathrm{HH}} \boldsymbol{S}_{\mathrm{HV}}^{*} \right\rangle - f_{\mathrm{v}} / 3 = f_{\mathrm{s}} \boldsymbol{\beta} + f_{\mathrm{d}} \boldsymbol{\alpha} \end{cases}$$
(9)

(9) is the residual model derived from the FDD model. As shown in a previous study [11], the dominant scattering mechanism, that is, surface scattering or doublebounce scattering, can be determined by the sign of

 $\operatorname{Re}(\langle S_{\mathrm{HH}}S_{\mathrm{HV}}^*\rangle)$. When $\operatorname{Re}(\langle S_{\mathrm{HH}}S_{\mathrm{HV}}^*\rangle) > 0$, surface scattering dominates, and $\alpha = -1$. When $\operatorname{Re}(\langle S_{\mathrm{HH}}S_{\mathrm{HV}}^*\rangle) < 0$, the dominant mechanism is double-bounce scattering, and $\beta = 1$. Therefore, with the determination of the dominant scattering component, there are only three unknowns in (9): $f_{\mathrm{s}}, f_{\mathrm{d}}$, and α or β . There are three independent equations in (9), so all unknowns can be solved.

2.2. Residual Model Extracted from YD Method

Considering the special case, azimuthally symmetric in urban areas, Yamaguchi *et al.* added a helix as the fourth component [5]. The covariance matrix for helix scattering $\langle [C] \rangle_{h}$ is,

$$\left< [C] \right>_{h} = \frac{1}{4} f_{h} \begin{bmatrix} 1 & \pm j\sqrt{2} & -1 \\ \mp j\sqrt{2} & 2 & \pm j\sqrt{2} \\ -1 & \mp j\sqrt{2} & 1 \end{bmatrix}$$
(10)

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Then, the covariance matrix $\langle [C] \rangle$ of (1) is modeled as the sum of the four scattering components:

$$\langle [\boldsymbol{C}] \rangle = f_{v} \langle [\boldsymbol{C}] \rangle_{v} + f_{d} \langle [\boldsymbol{C}] \rangle_{d}$$

$$+ f_{s} \langle [\boldsymbol{C}] \rangle_{s} + f_{h} \langle [\boldsymbol{C}] \rangle_{h}$$
(11)

where f_v, f_d, f_s , and f_h denote the scattering coefficients for each scattering mechanism, respectively.

Yamaguchi *et al.* modified the volumetric scattering component [5] that was modeled as randomly distributed long thin cylinders [1] to reflect distribution patterns of tree trunks and branches. A general expression to represent the volume scattering model $\langle [C] \rangle_{v}$ is,

$$\left\langle \begin{bmatrix} \boldsymbol{C} \end{bmatrix} \right\rangle_{v} = f_{v} \begin{bmatrix} a & 0 & d \\ 0 & b & 0 \\ d^{*} & 0 & c \end{bmatrix}$$
(12)

With the value of $R = 10 \log \left(\frac{|S_{VV}|^2}{|S_{HH}|^2} \right)$ as a threshold,

three sets of *a*-*d* are chosen to represent three volumetric scattering models. In the first model, R < -2 dB, and *a*-*d* are sets such that $a = \frac{3}{15}, b = \frac{4}{15}, c = \frac{8}{15}$, and

 $d = \frac{2}{15}$. In this case, the volumetric scattering is modeled

as a cloud of horizontally oriented dipoles. When 3 2 3 1

$$|R| \le 2 \text{ dB}, a = \frac{3}{8}, b = \frac{2}{8}, c = \frac{3}{8}, \text{ and } d = \frac{1}{8}, \text{ and the co-$$

variance matrix of $\langle [C] \rangle_{v}$ is the same as that in the FDD model. When R > 2 dB, $a = \frac{8}{15}, b = \frac{4}{15}, c = \frac{3}{15}$, and

 $d = \frac{2}{15}$, and the volumetric scattering is modeled as a cloud of vertically oriented dipoles. In addition, the helix scattering coefficient $f_{\rm h}$ and the volume scattering coefficient $f_{\rm v}$ are directly computed from (11),

$$f_{\rm h} = 2 \left| \operatorname{Im} \left(\left\langle \boldsymbol{S}_{\rm HH} \boldsymbol{S}_{\rm HV}^* \right\rangle + \left\langle \boldsymbol{S}_{\rm HV} \boldsymbol{S}_{\rm VV}^* \right\rangle \right) \right|,$$

$$f_{\rm v} = \frac{2}{b} \left(\left\langle \left| \boldsymbol{S}_{\rm HV} \right|^2 \right\rangle - \frac{1}{4} f_{\rm h} \right)$$
(13)

The energy contribution P_v corresponds to the volumetric scattering such that,

$$P_{\rm v} = \begin{cases} \frac{15}{2} \left(\left\langle \left| \boldsymbol{S}_{\rm HV} \right|^2 \right\rangle - \frac{1}{4} f_{\rm h} \right) R < -2 \, \mathrm{dB} \\ 8 \left(\left\langle \left| \boldsymbol{S}_{\rm HV} \right|^2 \right\rangle - \frac{1}{4} f_{\rm h} \right) \quad |R| \le 2 \, \mathrm{dB} \\ \frac{15}{2} \left(\left\langle \left| \boldsymbol{S}_{\rm HV} \right|^2 \right\rangle - \frac{1}{4} f_{\rm h} \right) R > 2 \, \mathrm{dB} \end{cases}$$
(14)

Similar to the removal of f_v to obtain (9), and after the elimination of helix scattering and volumetric scattering related components, (11) can be expressed as:

$$\begin{cases} \left\langle \left| \boldsymbol{S}_{\mathrm{HH}} \right|^{2} \right\rangle - a f_{\mathrm{v}} - \frac{1}{4} f_{\mathrm{h}} = f_{\mathrm{s}} \left| \boldsymbol{\beta} \right|^{2} + f_{\mathrm{d}} \left| \boldsymbol{\alpha} \right|^{2} \\ \left\langle \left| \boldsymbol{S}_{\mathrm{VV}} \right|^{2} \right\rangle - c f_{\mathrm{v}} - \frac{1}{4} f_{\mathrm{h}} = f_{\mathrm{s}} + f_{\mathrm{d}} \\ \left\langle \boldsymbol{S}_{\mathrm{HH}} \boldsymbol{S}_{\mathrm{HV}}^{*} \right\rangle - d f_{\mathrm{v}} + \frac{1}{4} f_{\mathrm{h}} = f_{\mathrm{s}} \boldsymbol{\beta} + f_{\mathrm{d}} \boldsymbol{\alpha} \end{cases}$$
(15)

Also equivalently to (9), (15) is the residual model extracted from the YD model. With a given

 $\operatorname{Re}(\langle S_{HH}S_{HV}^*\rangle)$ and f_h and f_v derived from (13), the remaining unknown parameters f_s , f_d , α , and β in (15) can be solved.

2.3. Residual Model Extracted from UmFDD Method

Considering that the volumetric scattering mechanism has high polarimetric entropy H, An *et al.* introduced a new model [7], the unit matrix, which corresponds to totally random scattering. The coherency matrix for the new volume scattering model $\langle [T] \rangle_{vunit}$ is,

$$\langle [\mathbf{T}] \rangle_{\text{vunit}} = \frac{1}{3} f_{\text{vunit}} \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$
 (16)

The correlation between the coherency matrix and the covariance matrix is,

$$\left\langle \begin{bmatrix} \boldsymbol{C} \end{bmatrix} \right\rangle_{\text{vunit}} = \begin{bmatrix} D_3 \end{bmatrix} \left\langle \begin{bmatrix} \boldsymbol{T} \end{bmatrix} \right\rangle_{\text{vunit}} \begin{bmatrix} D_3 \end{bmatrix}^{-1} = \left\langle \begin{bmatrix} \boldsymbol{T} \end{bmatrix} \right\rangle_{\text{vunit}},$$

$$D_3 = \begin{bmatrix} 1/2 & 1/2 & 0 \\ 0 & 0 & \sqrt{2}/2 \\ 1/2 & -1/2 & 0 \end{bmatrix}$$

$$(17)$$

The new volumetric scattering model $\langle [C] \rangle_{vunit}$ is used to substitute for the original one in FDD [7], and the scattering coefficient f_{vunit} is,

$$f_{\text{vunit}} = 6 \left\langle \left| \boldsymbol{S}_{\text{HV}} \right|^2 \right\rangle$$
 (18)

The scattering energy contribution P_v corresponding to this model is,

$$P_{\rm v} = 6 \left\langle \left| \boldsymbol{S}_{\rm HV} \right|^2 \right\rangle \tag{19}$$

The residual model extracted from UmFDD is:

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$$\begin{cases} \left\langle \left| \boldsymbol{S}_{\rm HH} \right|^2 \right\rangle - \frac{1}{3} f_{\rm vunit} = f_{\rm s} \left| \boldsymbol{\beta} \right|^2 + f_{\rm d} \left| \boldsymbol{\alpha} \right|^2 \\ \left\langle \left| \boldsymbol{S}_{\rm VV} \right|^2 \right\rangle - \frac{1}{3} f_{\rm vunit} = f_{\rm s} + f_{\rm d} \\ \left\langle \boldsymbol{S}_{\rm HH} \boldsymbol{S}_{\rm HV}^* \right\rangle = f_{\rm s} \boldsymbol{\beta} + f_{\rm d} \boldsymbol{\alpha} \end{cases}$$
(20)

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2.4. A Generalized Residual Model

Three of decomposition models: FDD, YD, and UmFDD, all can decompose the whole backscattering into three basic scattering mechanisms: volumetric scattering, double-bounce scattering and surface scattering. The differences are that the choice for the volume scattering model is different and the result for the volume scattering energy contribution is distinct. The volume scattering energy computed from FDD is the largest, YD followed by, and the value extracted from UmFDD is lowest.

After a close examination of (9), (15) and (20), one should note that the three residual models are structurally similar, with the double-bounce scattering and surface scattering related components on the right side of the equals sign in both equations. Symbolically, the right sides in (9), (15) and (20) are identical. Thus, a general expression can be used to represent them:

$$\begin{cases} A = f_{\rm s} \left| \beta \right|^2 + f_{\rm d} \left| \alpha \right|^2 \\ B = f_{\rm s} + f_{\rm d} \\ C = f_{\rm s} \beta + f_{\rm d} \alpha \end{cases}$$
(21)

When $A = \langle |\mathbf{S}_{\text{HH}}|^2 \rangle - f_v, B = \langle |\mathbf{S}_{\text{VV}}|^2 \rangle - f_v$, and $C = \langle \mathbf{S}_{\text{HH}} \mathbf{S}_{\text{HV}}^* \rangle - f_v / 3$, (21) is the residual model of (9). When $A = \langle |\mathbf{S}_{\text{HH}}|^2 \rangle - af_v - \frac{1}{4}f_h, B = \langle |\mathbf{S}_{\text{VV}}|^2 \rangle - cf_v - \frac{1}{4}f_h$,

and $C = \langle S_{\text{HH}} S_{\text{HV}}^* \rangle - df_v + \frac{1}{4} f_h$, (21) becomes the residual model of (15). When

$$A = \left\langle \left| \boldsymbol{S}_{\text{HH}} \right|^2 \right\rangle - \frac{1}{3} f_{\text{vunit}}, B = \left\langle \left| \boldsymbol{S}_{\text{VV}} \right|^2 \right\rangle - \frac{1}{3} f_{\text{vunit}}, \text{ and} C = \left\langle \boldsymbol{S}_{\text{HH}} \boldsymbol{S}_{\text{HV}}^* \right\rangle, (21) \text{ becomes (20).}$$

3. Properties of the Generalized Residual Model

3.1. Relationship between Volumetric Scattering and Negative Energy

Because f_s , f_d , f_v , and f_h denote energy coefficients for different types of scattering mechanisms, they should all be positive, which is consistent with the actual physical scattering mechanisms observed in reality. Because surface scattering and double-bounce scattering have no contribution to cross-polarized return, the volume scattering coefficient f_v can be directly estimated from PolSAR data. The helix scattering coefficient f_h can be computed from (13) as well. Both f_v and f_h are positive, which can be seen from Equations (7), (13) and (18). The parameters f_s and f_d are computed from the data using the generalized residual model (21). The difficult and troublesome problems in (21) are that the scattering coefficients f_s and f_d aresometimes negative, which is not consistent with the physical scattering mechanisms. Negative results for the coefficients f_s and f_d must be reduced or eliminated.

Comparing (8) and (14), the volume scattering contribution P_v computed from FDD is more overestimated than the ones extracted from the YD and UmFDD methods. In addition, the negative scattering coefficients f_d and f_s computed with YD and UmFDD are reduced compared with the results obtained from the FDD method. That is to say, the value for volume scattering energy contribution P_v is strongly related to the scattering component coefficients f_d and f_s . We can reduce the negative coefficients f_d and f_s by decreasing the volume scattering energy contribution P_v . The following experiments will be used to validate that the emergence of negative values for the scattering mechanisms f_d and f_s should be attributed to the overestimation of the volume scattering contribution f_v .

3.2. Eigenvalues of the Generalized Residual Model

Based on the work of van Zyl *et al.* [12], the covariance matrix $\langle [C] \rangle$ is expressed as the sum of a determined scattering model $[C]_{model}$ and a residual model $[C]_{residual}$, which has the same meaning as (21) but with a different formulation:

$$\langle [C] \rangle = a [C]_{\text{model}} + [C]_{\text{residual}}$$
 (22)

where *a* is constant. In this paper, we set a = 1, so (22) can be rewritten as:

$$\begin{bmatrix} \boldsymbol{C} \end{bmatrix}_{\text{residual}} = \langle \begin{bmatrix} \boldsymbol{C} \end{bmatrix} \rangle - \begin{bmatrix} \boldsymbol{C} \end{bmatrix}_{\text{model}}$$
(23)

The residual model (23) can be expressed as,

$$\begin{bmatrix} \boldsymbol{C} \end{bmatrix}_{\text{residual}} = \begin{bmatrix} A & 0 & C \\ 0 & 0 & 0 \\ C^* & 0 & B \end{bmatrix}$$
$$= \begin{bmatrix} f_{\text{s}} |\beta|^2 + f_{\text{d}} |\alpha|^2 & 0 & f_{\text{s}}\beta + f_{\text{d}}\alpha \\ 0 & 0 & 0 \\ (f_{\text{s}}\beta + f_{\text{d}}\alpha)^* & 0 & f_{\text{s}} + f_{\text{d}} \end{bmatrix}$$
(24)

The variables A, B, and C in (24) have the same meaning as those in (21). Because the right-hand side of (24) represents physical scattering mechanisms, the eigenvalues computed from (24) must be positive [12]. Only positive eigenvalues correspond to actual physical scattering mechanisms, which has been discussed in the work of van Zyl *et al.* [12]. Because volumetric scattering and helix scattering have been eliminated, there are at most two non-zero eigenvalues computed from (24). Both of the two eigenvalues should be positive, but this is not true. The following experiment is carried out to

statistically count the percentage of negative eigenvalues computed from the residual model (24). It will show that the volumetric scattering contribution P_v has high correlation with the proportion of negative eigenvalues. P_v should be reduced to ensure that the residual model is able to represent the physical scattering mechanisms.

3.3. Volumetric Scattering Model Minimization

In the previous section's analysis, it was observed that an overestimated volumetric scattering contribution P_v will lead to negative coefficients f_d and f_s and negative eigenvalues computed from the residual model (24). Thus, we try to minimize the volumetric scattering energy contribution P_v , so that the scattering coefficients f_d and f_s will be accurately estimated.

We assume that volumetric scattering only has no contribution to co-polarized returns but to cross-polarized return. The covariance matrix for volumetric scattering, as shown in (25), corresponds to the minimum scattering energy contribution,

$$\left< [C] \right>_{\text{limit}} = \begin{bmatrix} 0 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 0 \end{bmatrix}$$
(25)

The corresponding volume scattering coefficient for the three decomposition models,

$$\begin{cases} f_{v} = 2\left\langle \left| \boldsymbol{S}_{HV} \right|^{2} \right\rangle, P_{v} = f_{v}, \text{ for } FDD \text{ and } UmFDD \\ f_{v} = 2\left\langle \left| \boldsymbol{S}_{HV} \right|^{2} \right\rangle - \frac{1}{2}f_{h}, P_{v} = f_{v}, \text{ for } YD \end{cases}$$
(26)

The modified residual model is written here:

$$\begin{cases}
A_m = f_s |\beta|^2 + f_d |\alpha|^2 \\
B_m = f_s + f_d \\
C_m = f_s \beta + f_d \alpha
\end{cases} (27)$$

When
$$A_m = \left\langle \left| \boldsymbol{S}_{\text{HH}} \right|^2 \right\rangle, B_m = \left\langle \left| \boldsymbol{S}_{\text{VV}} \right|^2 \right\rangle$$
, and

 $C_m = \langle S_{\rm HH} S_{\rm HV}^* \rangle$, (27) is the modified residual model extracted from FDD and UmFDD model. When

$$A_m = \left\langle \left| \boldsymbol{S}_{\text{HH}} \right|^2 \right\rangle - \frac{1}{4} f_{\text{h}}, B_m = \left\langle \left| \boldsymbol{S}_{\text{VV}} \right|^2 \right\rangle - \frac{1}{4} f_{\text{h}}, \text{ and}$$

 $C_m = \langle S_{\rm HH} S_{\rm HV}^* \rangle + \frac{1}{4} f_{\rm h}$, (27) becomes the modified re-

sidual model extracted from YD.

From (26), we learn that the volumetric scattering contribution P_v has been minimized. When P_v reaches its minimum, the negative scattering coefficients f_d and f_s will be greatly reduced, and the negative eigenvalues computed from the residual model (27) will be decreased as well. This view will be validated in the following experiments.

4. Experiment

The Advanced Land Observing Satellite (ALOS) L-band PolSAR dataset for San Francisco is used in this paper. The satellite image, including the ascending track, was acquired on 11 November 2009. The incidence angle is 22.74° in the near range and 24.95° in the far range. We chose a 512×512 sub-image for this experiment, and the sub-image is multilook processed in the azimuth direction.

First, the scattering component coefficients f_d and f_s are computed from the residual model (21). The statistical proportion of negative values for f_d and f_s is given in **Table 1**.

Some scattering coefficients f_d and f_s are negative. The percentage of negative values for f_d and f_s computed from FDD is higher than the ratios computed for the YD and UmFDD methods. The volume scattering energy contribution P_v extracted from FDD is overestimated compared with YD and UmFDD. This observation confirms that the overestimation of the volumetric scattering contribution P_v will lead to negative energy.

Second, the statistical percentages of negative eigenvalues computed from the residual model (24) are displayed in **Table 2**, where λ_1 denotes a relatively large eigenvalue, and λ_2 denotes a relatively small one.

It can be seen from **Table 2** that some eigenvalues λ_1 and λ_2 are negative. The YD and UmFDD methods perform better than FDD. The volume scattering model is strongly related to the residual model, as mentioned in the previous section. The residual model (24) represents physical scattering mechanisms. It is best to ensure that the eigenvalues computed from residual model (24) are positive. Only positive eigenvalues correspond to physical scattering mechanisms. The selection of an appropriate volumetric scattering model is a way to ensure positive eigenvalues.

Finally, a new volumetric scattering model (27) is introduced to the FDD, YD, and UmFDD methods. The

Table 1. Negative parameters $f_{\rm d}$ and $f_{\rm s}$ computed from (21).

Percentage negative	FDD method	YD method	UmFDD method
$f_{\rm d}$	13.08%	5.13%	3.08%
$f_{ m s}$	12.52%	0.91%	2.49%

Table 2. Negative eigenvalues λ_1 and λ_2 computed from (24).

Negative percentage	FDD method	YD method	UmFDD method
λ_{1}	1.40%	0.96%	0.35%
$\lambda_{_2}$	18.74%	11.13%	11.69%

scattering coefficients $f_{\rm d}$ and $f_{\rm s}$ are computed from the residual model (27). Eigenvalues are extracted from (27) as well. The results are shown in **Table 3**.

When the volume scattering energy contribution P_v is minimized, the scattering component coefficients f_d and f_s computed from the generalized residual model are almost all positive. The eigenvalues λ_1 and λ_2 are almost being positive. This observation indicates that the residual model (27) is able to represent surface scattering and double-bounce scattering.

The new volumetric scattering model (25) is adopted under the assumption that volumetric scattering has no contribution to co-polarized returns but only to crosspolarized return, which is not consistent with physical scattering in reality. In this situation, the volume scattering energy contribution is less overestimated. We should keep a balance between the volumetric scattering energy contribution and the negative scattering coefficients f_d and f_s so that the volumetric scattering model can reflect the physical scattering mechanisms and the scattering coefficients f_d and f_s are accurately estimated.

5. Conclusions

To reduce the negative scattering coefficients for doublebounce scattering f_d and surface scattering f_s , a general residual model is introduced, which is extracted from FDD, YD, and UmFDD methods. There is a high correlation between volume scattering energy P_v and scattering coefficients f_d and f_s . When the value of P_v is greater, the more negative f_d and f_s computed from the residual model.

In this paper, the authors adopt a new volume scattering model to minimize P_v . When P_v reaches to its minimum, the percentage of negative parameters f_d and f_s is greatly reduced, and the proportion of negative eigenvalue computed from residual model is rapidly decreased, which is validated using the actual Pol-SAR data. It is concluded that decreasing the volume scattering contribution P_v can help to reduce the presence of negative values for the scattering coefficients f_d and f_s .

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Table 3. Negative parameters $f_{\rm d}$, $f_{\rm s}$, λ_1 and λ_2 computed from (27).

Percentage negative	FDD/UmFDD method	YD method
$f_{ m d}$	0.00%	0.26%
$f_{ m s}$	0.00%	0.00%
λ_1	0.00%	0.00%
λ_2	0.00%	0.26%

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