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M-NL: Robust NL-means approach for PolSAR images denoising

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Abstract—This paper proposes a new method for polarimetric synthetic aperture radar (PolSAR) denoising. More precisely, it seeks to address a new statistical approach for weights computation in non-local (NL) approaches. The aim is to present a simple criterion using M-estimators and to detect similar pixels in an image. A binary hypothesis test is used to select similar pixels which will be used for covariance matrix estimation together with associated weights. The method is then compared to an advanced state of the art PolSAR denoising method, NL-SAR method [1]. The filter performances are measured by a set of different indicators, including relative errors on incoherent target decomposition parameters, coherences, polarimetric signatures, and edge preservation on a set of simulated PolSAR images, as in [2]. Finally, results for RADARSAT-2 PolSAR data are presented.

Index Terms—PolSAR, NL means, M-estimators, Wishart distribution, Detection

I. INTRODUCTION

N the past decades, there has been a growing interest for polarimetric synthetic aperture radar (PolSAR) images and their use for terrain classification, target detection etc. Speckle in PolSAR data significantly degrades the image quality as well as the application performances. In PolSAR images each pixel is given by a complex-valued matrix (or vector) formed out of backscattered signals in different combinations of the linear received and transmitted polarizations. Since the scatterers are distributed and due to the coherency of PolSAR systems, this matrix has a random nature and is refereed to as speckle noise. Therefore, in order to determine the physical parameters of interest, a speckle reduction step is usually applied, aiming at reducing the parameter fluctuations. Different methods have been proposed to reduce speckle in PolSAR data [3]. Recent approaches of image processing and computer vision rely on patch-based comparison to select similar samples. The idea was born with the first works of Buades et al. [4], [5] and then extended and adopted to (Pol)SAR images by several authors such as in [6], [7]. In these methods, the covariance matrices (CMs) of scattering vectors (SVs) are estimated using carefully selected samples in order to reduce the speckle. The SVs are usually modelled by complex circular Gaussian vectors, completely characterized by their CM. The CM, unknown in practice, is then obtained by averaging, i.e. using the well-known sample CM (SCM)

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which, in the Gaussian context, is known to be the Maximum Likelihood (ML)-estimators (MLE) and Wishart distributed.

Nevertheless, the Gaussian distribution is not always a good approximation for PolSAR data since, for instance, this model is not able to describe textured scenarios. Therefore, it is necessary to increase the complexity of the underlying distribution model. This has been successfully done by introducing the so-called Spherically Invariant Random Vectors (SIRV) [8]. The SIRV displays an observation as a product of a Gaussian random process with the square root of a nonnegative random scalar variable that contains the information about the texture variation. Consequently, the SIRVs encompass a wide class of well-known distributions such as t-distribution, K-distribution or inverse Gaussian texture distribution etc. In a non-Gaussian environment, the SCM may not provide accurate results for the CM estimate. For this reason, M-estimators have been introduced and broadly employed and analysed in robust statistics and signal processing [9]-[11]. M-estimators are a wide class of scatter matrix (SM)¹ estimators robust to the data model. These estimators are given by fixed-point equations which makes the direct analysis of their statistical properties very difficult. Nevertheless, it has been recently shown that the behaviour of M-estimators can be well approximated by Wishart distribution [12], [13] which is of great importance in various applications.

In this paper, thanks to the new properties of M-estimators, we introduce a new NL-means algorithm to estimate CM in PolSAR data. Then, we evaluate performances of this method and compare it to NL-SAR method. NL-SAR is a state of the art method composed of several steps: pre-estimation, weights computation, bias reduction and best estimate selection. In this work, the focus is on the two first steps:

- Pre-estimation: Instead of using a Gaussian kernel, we propose to use an *M*-estimator in order to pre-estimate matrices in each pixel and compare patches.
- Weight computation: In this step, instead of learning kernels to weights similarities, we first select similar pixels based on the statistics of a robust similarity test and, then, compute weights using only selected pixels.

These steps are repeated for a set of different parameters which results in several estimates in each pixel. Finally, we perform bias-reduction and select the best estimate as in NL-SAR.

This paper is organised as follows. Section II introduces some background about PolSAR images and M-estimators. Section III presents the proposed method and the correspond-

¹SM is equal to the CM up to a scale factor when the latter exists.

ing algorithm, with discussions. In Section IV the results for simulated and RADARSAT-2 data are presented. Finally, some conclusions are drawn in Section V.

Notations. Vectors (resp. matrices) are denoted by bold-faced lowercase letters (resp. uppercase letters). T and H respectively represent the transpose and the Hermitian operator. $\text{Tr}(\cdot)$ represents the trace of a matrix. Finally, $|\cdot|$ stands for the determinant of a matrix.

II. THEORETICAL BACKGROUND

A. PolSAR

Polarimetric SAR sensors measure the amplitude and phase of backscattered signals in four combinations of the linear received and transmitted polarizations: horizontal-horizontal (HH), horizontal-vertical (HV), vertical-horizontal (VH) and vertical-vertical (VV). These signals form the complex scattering matrix S

$$\mathbf{S} = \begin{bmatrix} S_{HH} & S_{HV} \\ S_{VH} & S_{VV} \end{bmatrix}$$

where S_{IJ} denotes the complex scattering amplitude for the received polarization I and the transmitted polarization J. The reciprocity theorem says that the cross-pol channels of the scattering matrix are equal, that is $S_{HV}=S_{VH}$. Therefore, there are only three independent complex coefficients required to characterize the SV s = $\begin{bmatrix} S_{HH} & \sqrt{2}S_{HV} & S_{VV} \end{bmatrix}^T$ or alternatively, the SV is replaced by the linear transformation $\mathbf{k} = \frac{1}{\sqrt{2}} \begin{bmatrix} S_{HH} + S_{VV} & S_{HH} - S_{VV} & 2S_{HV} \end{bmatrix}^T$ known as the Pauli representation of the SV [14].

B. PolSAR modeling

So far the most used model for underlying distribution of PolSAR data is the complex Gaussian circular distribution. To reduce the effect of inherent speckle noise, PolSAR images are often spatially averaged, and the data are represented by the SCM $\widehat{\Sigma}_{SCM} = \frac{1}{N} \sum_{n=1}^{N} \mathbf{s}_n \mathbf{s}_n^H$ (or the sample coherency matrix $\widehat{\Sigma}_{SCM} = \frac{1}{N} \sum_{n=1}^{N} \mathbf{k}_n \mathbf{k}_n^H$). Under the assumption that N > 3 the estimate is complex Wishart distributed with N degrees of freedom (DoF) around the expectation value Σ (true CM of Gaussian data).

Since the Gaussian distribution fails to give a good approximation in high textured scenarios an alternative is to use SIRV model: $\mathbf{k} = \sqrt{\tau} \mathbf{n}$, where τ is a texture parameter whose distribution is unspecified, with mean value equal to 1, and \mathbf{n} is the speckle vector, following a multivariate Gaussian distribution. SIRV represent a subclass of complex elliptically symmetric (CES) distributions. In this context, the performances of the SCM can be really degraded and one needs other estimators that are better adapted to a non-Gaussian framework.

C. M-estimators

Let $(\mathbf{k}_1, \dots, \mathbf{k}_N)$ be an N-sample of m-dimensional complex independent CES-distributed vectors (with the same PDF). An M-estimator, is defined by the solution of the following M-estimating equation

$$\widehat{\Sigma} = \frac{1}{N} \sum_{n=1}^{N} \varphi \left(\mathbf{k}_{n}^{H} \widehat{\Sigma}^{-1} \mathbf{k}_{n} \right) \mathbf{k}_{n} \mathbf{k}_{n}^{H}$$
(1)

where φ is any real-valued weight function on $[0,\infty)$ (see [9]) that does not need to be related to the PDF of the underlying distribution. Existence and uniqueness of the solution of (1) as well as the convergence of the corresponding recursive algorithm have already been shown, provided the function φ satisfies a set of general assumptions [11]. In particular, the resulting estimators are consistent estimators of the SM (up to a scale factor). Some of widely used M-estimators are: Tyler's M-estimator [10], [15], Huber's M-estimator, Student's M-estimator.

III. ROBUST M-NLMEANS METHOD

In this section we present step-by-step the proposed method for weights computation and discuss about its benefits.

A. Robust pre-estimation

In order to compute the (dis)similarity between two pixels, one needs to compute a pre-estimation of the CM. In NL-SAR, this pre-estimation is done using truncated Gaussian on the patch of the size $S=(2s+1)\times(2s+1)$. In this method we propose to use a Student's M-estimator given as the solution of

$$\widehat{\mathbf{\Sigma}}_t = \frac{m + \nu/2}{S} \sum_{n=1}^{S} \frac{\mathbf{k}_n \mathbf{k}_n^H}{\nu/2 + \mathbf{k}_n^H \widehat{\mathbf{\Sigma}}_t^{-1} \mathbf{k}_n}$$
(2)

where m is the vector dimension and ν is the DoF parameter. This estimator is very interesting since it represents a sort of trade-off between the SCM $(\nu \to \infty)$ and the Tyler's estimator $(\nu=0)$, the least and the most robust estimators. The SCM can be completely degraded with only one outlier while, on the other hand, the Tyler's estimator is entirely resistant to aberrations but gives the estimation of so-called "shape" matrix, i.e. loses the information about matrix power. Therefore, the Student's M-estimator represents a good compromise that simultaneously can be robust and preserves the power information.

B. Pixel selection

Using these pre-estimated values, neighbouring samples are selected around each pixel. The central pixel at location l is compared to all pixels in a circular window following a spiral path (see [1] for more details). To compute the dissimilarities between two pixels instead of classical generalized likelihood ratio tests (GLRT), we propose to use the Box's M-test defined

$$\mathcal{L} = rac{\left|\widehat{oldsymbol{\Sigma}}_1
ight|^{S/2}\left|\widehat{oldsymbol{\Sigma}}_2
ight|^{S/2}}{\left|\widehat{oldsymbol{\Sigma}}
ight|^S}$$

where $\widehat{\Sigma}_1$ is obtained with a sample $\mathbf{k}^{(1)} = (\mathbf{k}_1, \dots, \mathbf{k}_S)$, $\widehat{\Sigma}_2$ with $\mathbf{k}^{(2)} = (\mathbf{k}_{S+1}, \dots, \mathbf{k}_{2S})$ and $\widehat{\Sigma}$ with $\mathbf{k} = (\mathbf{k}^{(1)}, \mathbf{k}^{(2)})$. This statistic has values between 0 and 1, where the values close to 0 reject the hypothesis that the matrices $\widehat{\Sigma}_1$ and $\widehat{\Sigma}_2$ are equal and values close to 1 accept it. By modifying the statistic \mathcal{L} , Box has obtained the approximated χ^2 distribution:

$$u = -2(1-\beta)\ln(\mathcal{L}) \sim \chi^2(m(m+1)/2)$$

with $\beta = \frac{3}{2S} \frac{2m^2 + 3m - 1}{6(m+1)}$, where m is the size of the SV, usually 3.We then propose to compute the similarity between two patches centred in pixels l and l' as

$$\Delta(l, l') = \sum_{\tau} u\left[(l + \tau), (l' + \tau) \right], \tag{3}$$

where $\tau \in [-p,p]$ is a 2-D shift indicating the location within each patch of size $P=(2p+1)\times(2p+1)$. Dissimilarities are then compared to the threshold in order to select similar pixels. Under the hypothesis \mathcal{H}_0 that two patches follow the same ditribution, Δ has the χ^2 distribution with $d=\frac{1}{2}m(m+1)P$ DoF. The critical region of the test is the given by

$$\left\{ R_c = \Delta, \Delta > \chi^2_{P_{fa}}(d) \right\}$$

with P_{fa} the probability of false alarm and $\chi^2_{P_{fa}}(d)$ the quantile of order $1-P_{fa}$ of the $\chi^2(d)$.

C. Weights computation

Once similar pixels are chosen, we proceed to the weights computation. To define the weights from the dissimilarity measure $\Delta(l, l')$, we propose to use an exponential kernel

$$\omega(l, l') = \begin{cases} e^{-\frac{|\Delta(l, l') - c|}{\lambda}} & \text{if } l \neq l'\\ 1 & \text{if } l = l'. \end{cases}$$
 (4)

The parameter $c=\mathbb{E}\left[\Delta(l,l')|\mathcal{H}_0\right]$ is the expected dissimilarity of two patches under the \mathcal{H}_0 and the threshold λ can be computed as $\lambda=F_{\chi^2(d)}^{-1}(1-P_{FA})$. This mapping from the (dis)similarities to the weights prevents any pixel from having a larger weight than the central pixel. The parameter c has been introduced in order to give a weight close to 1 when the compared pixels comes from the same distribution while preventing the noise enforcing. We normalize the quantity with λ in order to obtain comparable weight values for different values of p.

Finally, the weighted maximum likelihood estimator is given by the weighted means

$$\widehat{\mathbf{\Sigma}}_{NL}(l) = \frac{\sum_{l'} \omega(l, l') \mathbf{k}' \mathbf{k}'^{H}}{\sum_{l'} \omega(l, l')}.$$
 (5)

The method is recapped in Algorithm 1: First, the maximum sizes of search windows W, patch P and pre-estimation scale S are set together with the threshold λ and the constant c that differs for each patch size. Then, the pre-estimation is performed for all pixels in the image and for all values of $0 \le s \le S$ where s = 0 means the pre-estimation is off, i.e. the matrix is equal to $\mathbf{k}\mathbf{k}^H$ for the pixel at location lwith coordinates (x, y). Then, for all window sizes the central pixel is compared to all pixels in the window using the preestimations corresponding to different values of s and different sizes of patches to perform the patch comparison. Then, for each triple of (s, p, w) an estimate Σ_{NL} is computed. Afterwards, for each $\widehat{\Sigma}_{NL}$ the new one $\widehat{\Sigma}_{NLRR}$ is obtained in order to reduce the bias introduced in the first one and finally, the best one (with highest number of looks) is selected for each pixel giving the final filtered image. This final part, enclosed within the box, is the same as in NL-SAR and will not be detailed because of lack of space.

```
Algorithm 1: M-NL method
 Initialization: W, P, S, \lambda, c, \nu
 forall x, y do
      for s = 0 : S (scale size) do
      Pre-estimation with (2)
 forall x, y (coordinates of pixel l) do
      for w = 1 : \mathcal{W} (search window size) do
          Compute \Delta x and \Delta y
          x' = x + \Delta x
          y' = y + \Delta y;
                                        ⊳ coordinates of pixel l'
          for s = 0 : \mathcal{S} do
               for p = 1 : \mathcal{P} (patch size) do
                    Compute \Delta(l, l') with (3)
                    if \Delta(l, l') \leq \lambda[p] then
                        Compute \omega(l, l') with (4)
                    else
                     \mid \omega(l, l') \leftarrow 0
       forall s, p, w do
            Compute \widehat{\Sigma}_{NL} with (5)
             Bias-reduction step \rightarrow \widehat{\Sigma}_{NLRB}
```

IV. EXPERIMENTS: IMPLEMENTATION AND EVALUATION

return The best estimate

A. Simulated data

In this section the results obtained for simulated and Pol-SAR data are presented. The simulated images have been generated using a Markov Random Field (MRF) following a Gibbs distribution as in [2]. Then, a polarimetric behavior has been assigned to the different parts of the designed images. The polarimetric signatures have been sampled from the Pol-SAR data as explained in the reference. The procedure is the following: First, a random number C of polarimetric classes is chosen between 3 and 5, C-1 classes for distributed scatterers and the last class for point scatters (targets). A ground truth is generated using a MRF and the targets correspond to squares of sizes varying between 2×2 and 5×5 pixels. For each distributed scatterer, one of the seven possible polarimetric signatures is randomly assigned and the Gaussian speckle noise is generated according to them. Finally, the targets generated using the remaining eighth polarimetric signature are added to the speckle noise. The exact values of CMs can be found in [2]. After the denoising, the set of following parameters has been evaluated:

- Radiometric parameters σ : diagonal elements of the CM (power information)
- Complex correlation parameters ρ: derived from the three complex off-diagonal terms (channels correlation)
- Incoherent decomposition parameters: Entropy (H), Anisotropy (A) and the mean Alpha angle $(\overline{\alpha})$ scattering mechanism
- Co-polar and cross-polar polarization signatures (PS)[16]
- Edge preservation (EP).

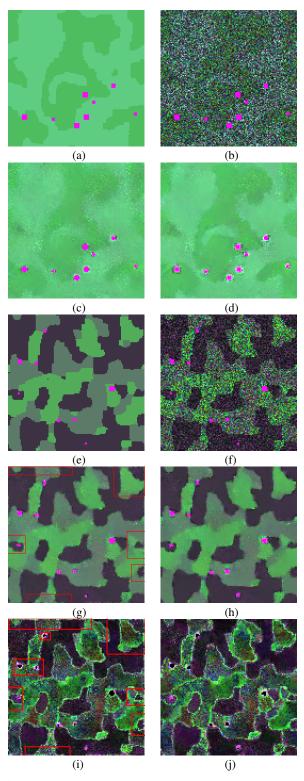


Figure 1: Application to simulated data: results for two different realizations. (a), (e) - ground truth, (b), (f) - speckle, (c), (g) - results obtained with NL-SAR, (d), (h) - results obtained with M-NL, (i), (j) - difference between the results and ground truth for NL-SAR and M-NL, respectively

For the first three groups of parameters the estimated value is obtained as the corresponding mean value from the pixels for every scattering class (SC) and for every simulated image,

given the filter. Then, the absolute relative bias of the estimated parameter is computed. The final (average) value is chosen as the median across all the simulated images and the various SCs. Then, the co- and cross-polar signatures of the average CM of the pixels belonging to a given SC of every simulated image are obtained (91x181 values). The average value for each class and image is obtained as a median value of absolute relative bias for every orientation and ellipticity angle. Then, as in previous case, the final value is obtained as the median across all the images and classes. Finally, edge preservation is measured on the boundary positions between extended targets. First, the gradient preservation (GP) is obtained as the average ratio between the observed gradient values on the power bands of the filtered power band image to the gradient values on the ground truth image. Then, the simple mapping from GP to EP is performed in order to give a measure close to 0 in the case of edge oversmoothing or undersmoothing and values close to 1 for good edge preservation.

In order to perform the comparison, we have simulated one hundred 128×128 artificial PolSAR images as described previously. The set of parameters used in both methods is: window size: $\{3^2, 5^2, \dots, 25^2\}$, patch size: $\{3^2, 5^2, \dots, 11^2\}$ and scale: $\{0, 1, 2\}$. Since the speckle is Gaussian, we have chosen ν to be big enough $(\nu = 100)$ in order to preserve the information about the texture and ensure the convergence of the solution in the pre-estimation step. The values for λ have been computed using the corresponding formula from Section III-C.

Figure 1 shows the results for two different simulated images. The images are presented in the following order, from left to right and from top to bottom: original image (ground truth), speckle, NL-SAR results and M-NL results. The first image Figure 1(a) contains two classes for distributed scatterers. As can be seen from Figure 1(c) et Figure 1(d) the homogeneous areas are much better smoothed with M-NL than with NL-SAR while the edges are better captured and less blurred. The second image Figure 1(e) consists of three SCs. In this case we display the filtered images together with the difference of the filtered image and corresponding ground truth in order to better visualize the results. From Figure 1(i) and Figure 1(j) one can see that in some parts of the images the noise is apparently more reduced with M-NL. This is the most visible (green points) at the image borders, one can look for instance the top border and corners of the images. Obviously, the M-NL gives an estimation closer to the ground truth in both cases. The visualization of the difference also reveals that most of the targets are better estimated with M-NL, which can not be seen directly from Figure 1(g) and Figure 1(h). Some of them are marked in red squares. Table I lists the

Filters	σ	$ \rho $	$\angle \rho$	H
NL-SAR	2.21	7.47	11.96	14.51
M-NL	1.56	9.10	14.47	14.49
	$\overline{\alpha}$	A	PS	EP
NL-SAR	35.84	10.51	1.15	0.45
M-NL	33.96	10.97	1.05	0.56

Table I: Filtering results for simulated data: all measures but EP (EP \in [0, 1]) are absolute relative errors in %.

evaluation parameters defined above. Numerical results have been computed over the set of simulated PolSAR images and the final values are compared. One can note that M-NL outperforms NL-SAR in almost all measures except ρ and A. Thus, it could be more convenient to use NL-SAR for terrain classification based on correlation coefficient or when measuring surface roughness implementing anisotropy parameter. On the other hand, M-NL gives better estimation of radiometric parameters, almost all incoherent decomposition parameters, polarisation signatures and edge preservation parameters. A significant improvement in edge preservation is also visually noticeable on the simulated images, thus these results are not surprising.

B. RADARSAT-2 PolSAR data

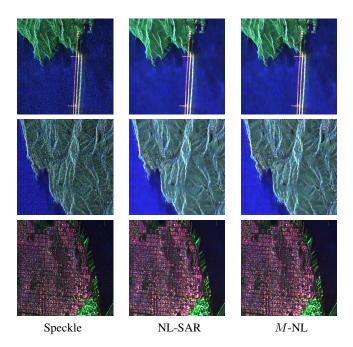


Figure 2: Real data: San Francisco Bay - 512×512 PolSAR images. From left to right: speckle, results obtained with NL-SAR and results obtained with M-NL.

Finally, the results for real data are given. Three different parts of San Francisco Bay are presented from top to bottom representing different scenarios in PolSAR images such as water, vegetation and urban areas. In this case, we do not dispose of any information about the ground truth, thus one can analyze the results only visually. First, one can note that M-NL better smoothes the homogeneous areas, while preserving well the edges in textured scenarios. It can also be noted that, as in the case of simulated data, M-NL gives results with higher contrast.

Last but not least, we would like to mention that an advantage of this method is also its simplicity, since after the preestimation it immediately starts to measure the dissimilarities and computes weights using a single, simple formula. On the other side, NL-SAR needs first to analyze noise in a homogeneous area in order to learn its distribution. Then, when the denoising step starts, a binary search is necessary to find the corresponding quantile and then a mapping is used to evaluate the weight, which obviously requires more steps.

V. CONCLUSION

This paper has investigated a new statistical approach for NL estimation in PolSAR imagery. The proposed method relies on M-estimators and has been compared to the NL-SAR method showing better results. It should be noted that the proposed method has common steps with NL-SAR. Yet, the estimation of NL estimates, the crucial part of any NL denoising method, is based on the statistical behavior of M-estimators, which improves the method simplicity.

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