PARAMETER ESTIMATION OF POLYNOMIAL PHASE SIGNAL BASED ON LOW-COMPLEXITY LSU-EKF ALGORITHM IN ENTIRE IDENTIFIABLE REGION

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ABSTRACT

Fast implementation of parameter estimation for polynomial phase signal (PPS) is considered in this paper. A method which combines the least squares unwrapping (LSU) estimator and the extended Kalman filter (EKF) is proposed. A small number of initial samples are used to estimate the PPS's parameters and then these coarse estimates are used to initial the EKF. The proposed LSU-EKF estimator greatly reduces the computation complexity of the LSU estimator and can work in entire identifiable region which inherits from the LSU estimator. Meanwhile, in the EKF stage its output is in point-by-point wise which is useful in real applications.

Index Terms— polynomial phase signal (PPS), parameter estimation, least squares unwrapping (LSU), extended Kalman filter (EKF), identifiable region

1. INTRODUCTION

Polynomial phase signals are common in fields including radar, sonar, geophysics, radio communication and biology [1, 2, 3]. The estimation of the parameters of polynomial phase signals has received considerable interest in recent years and several methods have been proposed to solve the problem. Among these methods, the maximum likelihood estimator (MLE) [4, 5] is an effective approach, being both computationally efficient and statistically accurate. However, this method becomes extremely complex when the order of PPSs is over 1.

In order to estimate the parameters of higher-order polynomial phase signals, researchers have developed many more efficient alternative methods based on 'multilinear transforms' or 'phase unwrapping'. Examples of these multilinear transform techniques include the cubic phase function (CPF) [6], the higher-order ambiguity function (HAF) [7] and its product version (PHAF) [8], which have good performance with acceptable computation and can be adapted to process multiple component signals. Their main limitation is the small identifiable region on which they can operate, owing to the error propagation. One alternative method is phase unwrapping, which utilizes the fact that the phase of a complex sinusoid is a linear function that is wrapped modulo 2π . A classical estimator based on phase unwrapping is the least squares unwrapping estimator [9, 10]. A significant advantage of this approach is that it works for polynomial phase parameters contained anywhere inside the identifiable region. The estimator also appears to perform well after thresholds at low signal-to-noise ratio (SNR) are reached. The major drawback of the LSU estimator is that computing a nearest lattice point is, in general, computationally difficult. Especially, the algorithm becomes prohibitively expensive as the number of samples is more than 1000 [9].

The parametric estimation for polynomial phase signals affected by Gaussian noise by extended Kalman filtering has been investigated in [11]. In this paper, we propose a PPS parameter estimation method which combines the LSU and the EKF. Firstly, the coarse estimates of the PPS's parameters are obtained by the LSU estimator which is performed on a small number of samples. Then, these coarse estimates are used to initial the EKF. The proposed method can work in entire identifiable region which is inherent from the LSU estimator and the estimation is output in point-by-point wise in the EKF stage. The numerical simulations show that the proposed method operating at a low SNR threshold exhibits high performance with small computation, and it is accurate over a far wider range of parameters than many popular existing estimators.

2. SIGNAL MODEL

A received polynomial phase signal can be written as

$$y(n) = Ae^{2\pi j\theta(n)} + w(n) \tag{1}$$

where A is the amplitude of the signal and w(n) is a complex additive white Gaussian noise (AWGN) with zero mean and variance σ^2 . $\theta(n)$ is the deterministic polynomial phase of order m, which is expressed by

$$\theta(n) = u_0 + u_1 n + u_2 n^2 + \dots + u_m n^m$$
(2)

where the coefficients u_i $(i = 0, \dots, m)$ are assumed to be real and unknown, and their identifiable regions are written as

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[12]

$$-\frac{1}{2i!} \le u_i < \frac{1}{2i!} (i = 0, 1, 2, \cdots m).$$
(3)

2.1. The State-Space Modelization

The discrete-time state vector X_k in the case of polynomial phase signals is given by the phase and the derivatives of the phase

$$\boldsymbol{X}_{k} = \begin{bmatrix} \theta(k) & \theta^{(1)}(k) & \theta^{(2)}(k) & \cdots & \theta^{(m)}(k) \end{bmatrix}^{T} \quad (4)$$

where $\theta^{(m)}(k)$ denotes m order derivation.

The model of polynomial phase signals can be described by a set of continuous-time differential equations cast in statespace form as

$$d\boldsymbol{X}(t)/dt = \boldsymbol{A}\boldsymbol{X}(t) + \boldsymbol{L}\boldsymbol{v}(t)$$
(5)

where the constant matrices $A_{m+1,m+1}$ and $L_{m+1,m}$, which characterize the behavior of the model, are written as

$$\boldsymbol{A} = \begin{bmatrix} 0 & 1 & 0 & \cdots & 0 & 0 \\ 0 & 0 & 1 & \cdots & 0 & 0 \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & 0 & \cdots & 1 & 0 \\ 0 & 0 & 0 & \cdots & 0 & 0 \end{bmatrix}, \boldsymbol{L} = \begin{bmatrix} 0 & 0 & \cdots & 0 \\ 1 & 0 & \cdots & 0 \\ 0 & 1 & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & 1 \end{bmatrix}$$
(6)

and v(t) is a white-noise process. There is a process-noise matrix Q_c that is related to the process-noise vector v(t) according to

$$\boldsymbol{Q}_c = E[\boldsymbol{v}(t)\boldsymbol{v}^T(t)]. \tag{7}$$

To be able to use the state (4), the model (5) must be discretized somehow [13], so that it can be described with a model of the form

$$\boldsymbol{X}_{k} = \boldsymbol{F}_{k} \boldsymbol{X}_{k-1} + \boldsymbol{v}_{k-1} \tag{8}$$

where F_k is the transition matrix of the dynamic model, which is evaluated as

$$\boldsymbol{F}_{k} = \boldsymbol{\vartheta}(T) = e^{\boldsymbol{A}T} = \boldsymbol{I} + \boldsymbol{A}T + \dots + \frac{(\boldsymbol{A}T)^{n}}{n!} + \dots \quad (9)$$

where T is the step-size of the discretization. On the other hand, v_{k-1} with variance Q_{k-1} is the discretized form of v(t)and Q_{k-1} satisfies

$$\boldsymbol{Q}_{k-1} = \int_0^T \boldsymbol{\vartheta}(\tau) \boldsymbol{L} \boldsymbol{Q}_c \boldsymbol{L}^T \boldsymbol{\vartheta}^T(\tau) d(\tau).$$
(10)

Therefore, the state-space modelization of polynomial phase signals is linear.

2.2. The Observation Equations

For the signal (1) is complex, it is necessary to express the measured signal as a 2×1 vector in terms of its real and imaginary part

$$\boldsymbol{Y}_{n} = \begin{bmatrix} \operatorname{Re}(y(n)) & \operatorname{Im}(y(n)) \end{bmatrix}^{T}.$$
 (11)

In this sense, the observation equation takes the form

$$\boldsymbol{Y}_k = h(\boldsymbol{X}_k) + \boldsymbol{\omega}_k \tag{12}$$

where the 2×1 nonlinear function $h(\mathbf{X}_k)$ is written as

$$h(\boldsymbol{X}_k) = \begin{bmatrix} h_{1k} \\ h_{2k} \end{bmatrix} = \begin{bmatrix} A\cos(2\pi\theta(k)) \\ A\sin(2\pi\theta(k)) \end{bmatrix}.$$
 (13)

The observation noise vector can be defined as $\omega_k = [w_R(k), w_I(k)]^T$, whose correlation matrix \mathbf{R}_k is given by

$$\boldsymbol{R}_{k} = \begin{bmatrix} \frac{\sigma_{\omega}^{2}}{2} & 0\\ 0 & \frac{\sigma_{\omega}^{2}}{2} \end{bmatrix}.$$
 (14)

In this paper, we assume that the process-noise v_k and the observation noise ω_k are not correlated, i.e.

$$\operatorname{cov}\left[\boldsymbol{\omega}_{k}, \boldsymbol{v}_{j}\right] = E\left[\boldsymbol{\omega}_{k}\boldsymbol{v}_{j}^{T}\right] = 0(k, j = 0, 1, 2, 3, \cdots).$$
(15)

Consequently, the observation model of polynomial phase signals is nonlinear.

3. THE EXTEND KALMAN FILTER ALGORITHM

As far as the observation is nonlinear, we use the extended Kalman filter algorithm to estimate the parameters of polynomial phase signals:

EKF Algorithm

Initial Conditions (k = 0):

$$\hat{\boldsymbol{X}}_{0} = E(\boldsymbol{X}_{0})$$

$$\hat{\boldsymbol{P}}_{0} = E\left[(\boldsymbol{X}_{0} - \hat{\boldsymbol{X}}_{0})(\boldsymbol{X}_{0} - \hat{\boldsymbol{X}}_{0})^{T}\right]$$
(16)

Predict Equations:

$$\begin{aligned} \mathbf{X}_{k}^{\prime} &= \mathbf{F}_{k} \hat{\mathbf{X}}_{k-1} \\ \mathbf{P}_{k}^{\prime} &= \mathbf{F}_{k} \hat{\mathbf{P}}_{k-1} \mathbf{F}_{k}^{T} + \mathbf{Q}_{k-1} \\ \mathbf{Y}_{k}^{\prime} &= \mathbf{H}_{k} \mathbf{X}_{k}^{\prime} \\ \mathbf{H}_{k} &= \left. \frac{\delta h}{\delta \mathbf{X}} \right|_{\mathbf{X}_{k}^{\prime}} \end{aligned}$$
(17)

Update Equations:

$$\boldsymbol{K}_{k} = \boldsymbol{P}_{k}^{\prime} \boldsymbol{H}_{k}^{T} \cdot \left[\boldsymbol{H}_{k} \boldsymbol{P}_{k}^{\prime} \boldsymbol{H}_{k}^{T} + \boldsymbol{R}_{k}\right]^{-1}$$
$$\hat{\boldsymbol{X}}_{k} = \boldsymbol{X}_{k}^{\prime} + \boldsymbol{K}_{k} \left[\boldsymbol{Y}_{k} - \boldsymbol{Y}_{k}^{\prime}\right]$$
$$\hat{\boldsymbol{P}}_{k} = \left[\boldsymbol{I} - \boldsymbol{K}_{k} \boldsymbol{H}_{k}\right] \boldsymbol{P}_{k}^{\prime}$$
(18)

The relationship between the parameters $\boldsymbol{u} = [u_0, u_1, \cdots, u_m]^T$ of polynomial phase signals and the initial state vector \boldsymbol{X}_0 is described by

$$\boldsymbol{u} = \boldsymbol{C}\boldsymbol{X}_0 \tag{19}$$

where the matrix C is derived as

$$\boldsymbol{C} = diag\left(\begin{bmatrix} 1 & 1 & \frac{1}{2!} & \cdots & \frac{1}{m!} \end{bmatrix}\right).$$
(20)

Therefore, we can evaluate the parameters of polynomial phase signals in the form

$$\boldsymbol{u} = \boldsymbol{C} \times \left(\prod_{m=1}^{k} \boldsymbol{F}_{m}^{-1}\right) \times \hat{\boldsymbol{X}}_{k}.$$
 (21)

The extended Kalman filter extends the scope of Kalman filter to nonlinear optimal filtering problems by forming a Gaussian approximation to the joint distribution of state X and measurements Y using a Taylor series based transformation. In this sense, the extended Kalman filter is sensitive to the initial conditions and if the filter is improperly initialized, the filtering may be failed. In this paper, in order to make the filter work effectively, we utilize the estimates of the least squares unwrapping estimator to initialize the EKF. Accounting for the computation load of the LSU estimator, only a few samples are used to obtain the initial estimates.

4. THE LEAST SQUARES UNWRAPPING (LSU) ESTIMATOR

The LSU estimator can estimate the parameters $\boldsymbol{u} = [u_0, u_1, \dots, u_m]^T$ from the phase of the observations $y(1), y(2), \dots, y(n)$. The phase θ_n of polynomial phase signals is written as

$$\theta_n = \angle y(n)/2\pi = \phi_n + \theta(n)(mod2\pi) \tag{22}$$

where \angle denotes the angle of a complex number, and ϕ_n are random variables representing the phase noise induced by w(n) and can be calculated as

$$\phi_n = \angle \left(1 + \frac{w(n)}{Ae^{2\pi j\theta(n)}} \right) / 2\pi \tag{23}$$

where ϕ_n is known as the projected normal distribution [14], if w(n) is complex Gaussian noise.

According to the fundamental of the least square method, the parameters $u_0, u_1, ..., u_m$ can be evaluated by

$$SS(\boldsymbol{u}) = \operatorname{argmin}\left(\sum_{n=1}^{N} \left(\theta_n - \sum_{k=0}^{m} u_k n^k\right)^2\right).$$
(24)

Due to the problem of the phase ambiguity in $\angle y(n)$, we may define

$$\left\langle \theta_n - \sum_{k=0}^m u_k n^k \right\rangle = \theta_n - \sum_{k=0}^m u_k n^k - \lfloor \varphi_n \rfloor$$
 (25)

where $\lfloor x \rfloor$ denotes the nearest integer to x. By considering $\lfloor \varphi_n \rfloor = \lfloor \theta_n - \sum_{k=0}^m u_k n^k \rfloor$ as nuisance parameters, (24) can be rewritten as

$$SS(\boldsymbol{u}) = \operatorname{argmin}\left(\sum_{n=1}^{N} \left\langle \theta_n - \sum_{k=0}^{m} u_k n^k \right\rangle^2\right)$$
(26)

where N represents the number of samples.

The LSU estimator is the minimum of SS(u) over the identifiable region of PPSs, which can be represented as a nearest lattice point problem [15]. When the number of samples is less than 60, the sphere decoder [16] can exactly computes a nearest lattice point to evaluate the parameters $u_0, u_1, ..., u_m$. In the other way, if the number of samples is more than 60, we can apply the K-best algorithm [17] having complexity $O(N^3 \log N)$ for approximating the LSU estimator. As a result, the LSU estimator is observed to occur at larger operations as the number of samples, N, grows. Especially, the LSU estimator is not suitable to run for the number of samples over 1000 [9].

The mean square error (MSE) of the LSU estimator can be obtained by

$$MSE = \frac{P_{signal}}{2\left[1 - f\left(-\frac{1}{2}\right)\right]^2 10^{\frac{SNR}{10}}} \times D^{-1}./G \qquad (27)$$

where P_{signal} denotes the power of the polynomial phase signal, $f(\cdot)$ is the probability density function of ϕ_n and A./B means element-by-element division between matrix A and matrix B. The superscript -1 denotes the inverse operation and \times denotes the product between a scalar and a matrix. Matrix D is a m + 1 by m + 1 Hilbert matrix with elements

$$D_{i,j} = 1/(i+j+1)(i,j \in [0,1,2,\cdots,m])$$
(28)

and the matrix G can be defined as

$$G = \begin{bmatrix} N & N^2 & \cdots & N^{m+1} \\ N^2 & N^3 & \cdots & N^{m+2} \\ \vdots & \vdots & \ddots & \vdots \\ N^{m+1} & N^{m+2} & \cdots & N^{2m+1} \end{bmatrix}.$$
 (29)

According to (27), to make the EKF work effectively at a low SNR, we should choose more samples for the LSU estimator. However, larger sample number means larger computation load. In this paper, the number of samples used for the LSU estimator is $N_1 = 64$.

5. SIMULATION

In this section, Monte-Carlo simulations are performed to verify the characteristic of our method. The numerical values are obtained over 2000 simulations. The number of samples used in the simulations is 199 and the number of initial samples used in our method is 64. Fig.1 shows the MSEs of the LSU estimator, the HAF estimator, the PHAF estimator, and the proposed estimator, where the values of coefficients are $\boldsymbol{u} = \begin{bmatrix} \frac{1}{4}, \frac{1}{4}, \frac{1}{8}, \frac{1}{24}, \frac{1}{96}, \frac{1}{480} \end{bmatrix}^T$ for the 5-order PPSs and $\boldsymbol{u} = \begin{bmatrix} \frac{1}{4}, \frac{1}{4}, \frac{1}{8}, \frac{1}{24} \end{bmatrix}^T$ for the 3-order PPSs. We can see that the HAF and PHAF estimators fail consistently, as expected. Although the ZW estimator [18] can operate on the entire identifiable region, its performance is poor, including its MSE and the threshold SNR.

The proposed LSU-EKF estimator reaches a threshold at 6 dB for the 3-order PPS and 17 dB for the 5-order PPS. The corresponding threshold SNRs for the LSU estimator are 2 dB for the 3-order PPS and 12 dB for the 5-order PPS respectively. The reason why the threshold SNRs of the LSU-EKF estimator are higher than those of the LSU estimator is that the number of samples used for the LSU estimator is larger than that for the LSU-EKF estimator in the parameter initialization step. Table 1 shows the running time of all the estimators, it is shown that the running time of the LSU-EKF estimator is approximately 24 times as much as that of the LSU-EKF estimator for the 3-order PPS and 15 times for the 5-order PPS.



Fig. 1: Sample MSEs for the coefficient u_3 for PPSs of order 3 and the coefficient u_5 for PPSs of order 5.

Table 1: Running time in seconds with different methods forPPSs of order 3 and PPSs of order 5.

	HAF	PHAF	ZW	LSU	LSU- EKF
m = 3	100.4s	180.5s	119.6s	25624.1s	1063.7s
m = 5	114.3s	174.1s	177.4s	44883.4s	3113.7s

In order to make the volume of the set of coefficients suit for the HAF and PHAF estimator, we chose the coefficients $u = \left[\frac{1}{4}, \frac{1}{4}, \frac{1}{8N}, \frac{1}{24N^2}, \frac{1}{96N^3}, \frac{1}{480N^4}\right]^T$ for the 5-order PPSs and $u = \left[\frac{1}{4}, \frac{1}{4}, \frac{1}{8N}, \frac{1}{24N^2}\right]^T$ for the 3-order PPSs. In Fig.2, we can see that the HAF estimator reaches a threshold at 6 dB, which is the same as our method for the 3-order PPSs and a threshold at 18 dB approximately a little higher than our method for the 5-order PPS. Meantime, the PHAF estimator has a good performance on the estimation of the coefficients u_3 for the 3-order PPS and u_5 for the 5-order PPS. As expected, the performance of our method remains unchanged for the set of small coefficients.



Fig. 2: Sample MSEs for the coefficient u_3 for PPSs of order 3 and the coefficient u_5 for PPSs of order 5.

6. CONCLUSION

The proposed LSU-EKF estimator greatly reduces the computation complexity of the LSU estimator while at the same time it can still work in entire identifiable region. The threshold SNR of our method is a few decibels higher than the LSU estimator because fewer samples are used in the LSU stage of the LSU-EKF estimator compared to the LSU estimator. Since the number of samples to initial the EKF is relative with the threshold SNR and the computation complexity, the size of the initial sample number should be careful chosen according to the tradeoff between these two factors.

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