

# An Efficient Forecasting Approach for the Real-Time Reduction of Boundary Effects in Time-Frequency Representations

Adrien Meynard, Hau-Tieng Wu

**Abstract**—Time-frequency (TF) representations of time series are intrinsically subject to the boundary effects. As a consequence, the structures of signals that are highlighted by the representations are garbled when approaching the boundaries of the TF domain. In this paper, for the purpose of real-time TF information acquisition for nonstationary oscillatory time series, we propose a numerically efficient approach for the reduction of such boundary effects. The solution relies on an extension of the analyzed signal obtained by a forecasting technique. In the case of the study of a class of locally oscillating signals, we provide a theoretical guarantee of the performance of our approach. Following a numerical verification of the performance of the algorithm, we validated its performance on various real-life biomedical signals.

**Index Terms**—Boundary effects, time-frequency, forecasting

## I. INTRODUCTION

IN any digital acquisition system, the study and the interpretation of the measured signals generally require an analysis tool, which enables researchers to point out the useful characteristics of the signal. The need for signal analysis arises from various signals, ranging from audio [1], [2], mechanical [3] or biomedical signals [4]. In this paper, we focus on oscillatory time series. For instance, biomedical signal, such as photoplethysmogram (PPG), contain several characteristics, such as respiratory rate or blood pressure, that cannot be interpreted from its running chart in the time domain. An analysis tool would make possible the extraction of these useful characteristics.

Usually, the measured signals exhibit nonstationary behavior, and the observed quantities might be interfered by transient phenomena that can vary rapidly and irregularly. For example, the signal might oscillate fast with large amplitude at one moment, and then oscillate slow with small amplitude at the next moment. In order to adapt the analysis to nonstationarities, local spectral analysis is generally performed [5], [6]. The short time Fourier transform [7] (STFT) is a typical tool which is build for this purpose, and enables the determination of the local frequency content of a nonstationary signal.

Windowing is a common method for performing local analysis. Among many others, STFT [8], continuous wavelet transform (CWT) [9], synchrosqueezing transform (SST) [10], and reassignment [11] (RS) are representations that fall back on the use of an analysis window. Let  $x : I \rightarrow \mathbb{R}$  denote the observed signal, where  $I$  denotes the finite interval where the signal is measured. Let  $g_s : \mathbb{R} \rightarrow \mathbb{R}$  denote the analysis window, where  $s$  is a shape parameter. The support of  $g_s$  is localized around the origin, and is small with respect to  $|I|$ . The translation operator is  $T_\tau$  defined as:

$$T_\tau f = f(t - \tau), \quad \forall f : \mathbb{R} \rightarrow \mathbb{R}.$$

Then, the local analysis of  $x$  around the instant  $\tau \in I$  rely on the evaluation of the following inner product:

$$V_x(s, \tau) = \langle x, T_\tau g_s \rangle_I. \quad (1)$$

A major shortcoming of this technique occurs when analyzing the signal  $x$  near the boundaries of the interval  $I$ . Clearly, at these points, half of the information is missing. Consequently the results of the inner product (1) is distorted. This phenomenon is usually understood as the *boundary effect*. We display on Fig. 1 the result of the SST of a PPG (see section IV-B3 for a comprehensive description). The distortion resulting from boundary effects is clearly visible on both sides of this representation. Indeed, while in the major central part of the image, clear lines stand out, they become blurred as they approach the left and right boundaries of the image. The right boundary is zoomed in and displayed on the right of Fig. 1, which emphasizes the result of boundary effects. Clearly, estimations of signal characteristics, like instantaneous frequencies [12] or amplitudes, from this TF representation appear to be imprecisely determinable (or even likely to fail) in the vicinity of the boundaries. Moreover, this boundary effect would unavoidably limit applying these window based TF analysis tools for the real-time analysis purpose. It is thus desirable to have a solution to eliminate this boundary effect.

Attempts to minimize the boundary effects generally consists in softening the discontinuity on signals edges. For instance, judicious choices of analysis windows whose support does not interfere with the boundary points can minimize the occurrence of aberrant patterns near the boundaries of the TF plane [13], [14]. Due to the

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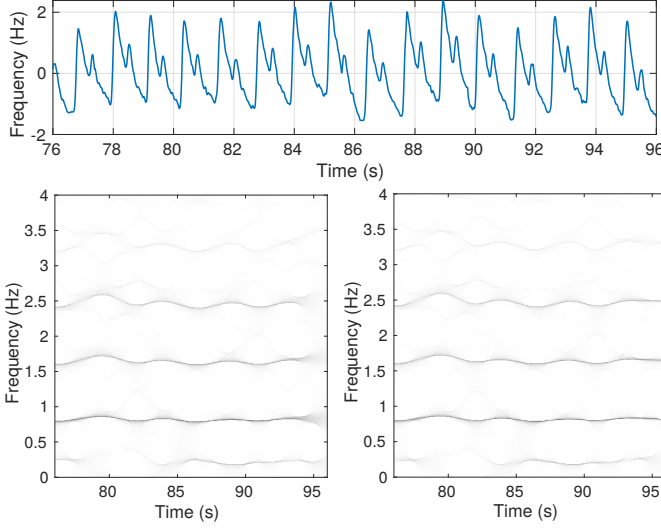


Figure 1. A segment of PPG signal (top) and the right boundary of a TF representation determined by the SST without extension (bottom left), and the right boundary of a TF representation determined by the SST with the proposed boundary effect reduction algorithm by forecasting (bottom right). The window length for the SST is 12 sec.

specific relationship of chosen analysis windows and TF analysis tool, these techniques do not make it possible to reduce the boundary effects of any TF representations, and for any analysis window. Another natural idea consists of carrying out a preliminary step of extending the signal beyond its boundaries, and due to its flexibility, various extension schemes have been proposed. For example, there exists simple extension schemes that do not take into account the dynamical behavior of the signal, such as zero-padding, periodic extension, symmetric extension [15], [16], or polynomial extrapolation [17]. There also exist extension schemes based on physically relevant dynamical models, such as the Extended Dynamic Mode Decomposition [18] (EDMD), the Gaussian process regression [19], [20] (GPR), the Trigonometric, Box-Cox transformation, ARMA errors, Trend and Seasonal components (TBATS) algorithm [21], the dynamic linear models [22], and long short-term memory model [23]. It is a far from exhaustive list, and we refer readers with interest to [24] for a friendly monograph on the general forecasting topic. While the second class of extension schemes give better extended signals than the first class, they generally have a great computational cost.

In this paper, we propose a fast extension algorithm based on a simple dynamical model, in order to optimize the trade-off between the extension quality and the computational cost. The proposed algorithm is composed of two steps.

- 1) *Extend the signal by forecasting it.* The aim is to use a dynamic model to predict the values taken by the measured signal outside the measurement interval. Then, once this operation is done, we have access to an extended signal defined on a larger interval  $I_\Delta$ ,

where  $\Delta$  denotes the size of the extension on both boundaries of  $I$ .

- 2) *Run the local analysis tool on the extended signal.* Assuming that the support of the analysis window is smaller than  $2\Delta$ , the local analysis near the boundary of  $I$  is now possible without lack of information thanks to knowledge brought by the extension on both sides.

Thus, assuming that the quality of the extension step is sufficient, the analysis results obtained that way will be less sensitive to the boundary effects than the result of the analysis tool applied directly to the non-extended signal. We claim and prove that forecasting oscillatory signals based on the simple dynamical model combined with the simple least square approach is sufficient for reducing the boundary effects for the TF analysis, or other kernel based analysis. See Figure 1 for a snapshot of the result. The main benefit of this simple approach is a numerically efficient solution with theoretical guarantee for the real-time analysis purpose.

The paper is organized in the following. In section II, we provide an extension method based on a linear dynamic model. We derive the corresponding algorithm for boundary effects reduction. In section III, we show that, even though we consider a simplistic dynamic model, it is sufficient to extend signals taking the form of sums of sine waves. An evaluation of the theoretical performance of our algorithm on a class of signals, the sums of sine waves, is given in section III. In section IV, we compare our extension method with more sophisticated methods such as EDMD, GPR, or TBATS. We show that our algorithm gives fast results of reasonable quality. Finally, we evaluate the performance of our boundary effects reduction algorithm on biomedical signals such as respiratory signals, and compare it to the theoretical results.

## II. ALGORITHM

As explained above, the algorithm for the reduction of boundary effects on TF representations relies on extending the signal by forecasting it before applying the TF analysis.

We start with the notation. Let  $x : \mathbb{R} \rightarrow \mathbb{R}$  denote a continuous-time signal. In this work, we consider a finite-length discretization of that one. Thus, the sampled signal  $\mathbf{x}$ , whose length is denoted by  $N$ , is such that

$$\mathbf{x}[n] = x\left(\frac{n}{f_s}\right), \quad \forall n \in \{0, \dots, N-1\},$$

where  $f_s$  denotes the sampling frequency. Let  $M$  and  $K$  be two integers such that  $M < N$  and  $K + M < N$ . Then, for all  $k \in \{0, \dots, K-1\}$ , we extract from  $\mathbf{x} \in \mathbb{R}^N$  the sub-signal  $\mathbf{x}_k \in \mathbb{R}^M$  given by:

$$\mathbf{x}_k = \begin{pmatrix} \mathbf{x}[N - K + (k-1) - (M-1)] \\ \vdots \\ \mathbf{x}[N - K + (k-1)] \end{pmatrix}. \quad (2)$$

These sub-signals are gathered into the matrix  $\mathbf{X} \in \mathbb{R}^{M \times K}$  such that:

$$\mathbf{X} = (\mathbf{x}_0 \quad \cdots \quad \mathbf{x}_{K-1}) .$$

Notice that these sub-signals are overlapping each other. Indeed,  $\mathbf{x}_{k+1}$  is a shifting of  $\mathbf{x}_k$  from one sample. We also consider the matrix  $\mathbf{Y} \in \mathbb{R}^{M \times K}$  given by:

$$\mathbf{Y} = (\mathbf{x}_1 \quad \cdots \quad \mathbf{x}_K) .$$

The boundary effect reduction algorithm is based on manipulating  $\mathbf{X}$  and  $\mathbf{Y}$ .

The pseudo-code of the proposed algorithm to reduce boundary effects on windowing-based TF representations is shown in the Algorithm 1. We coined the algorithm **BoundEffRed**. Below, we detail the algorithm, particularly the Signal extension algorithm **SigExt** in Algorithm 2.

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**Algorithm 1** Tackling boundary effects of a TF representation in real-time.  $\mathbf{F}_x = \text{BoundEffRed}(\mathbf{x}, M, K, L, \mathcal{F})$

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**Inputs:**  $\mathbf{x}_N, M, K, N_0, L, \mathcal{F}$

**while**  $N$  increases **do**

**Real-time input:**  $\mathbf{x}_N$

**Forecasting step.**

- Signal extension:  $\tilde{\mathbf{x}}_{N+L} = \text{SigExt}(\mathbf{x}_N)$ .

**Representation estimation step.**

- Extended representation evaluation:  $\mathcal{F}(\tilde{\mathbf{x}}_{N+L})$ .
- Restriction of  $\mathcal{F}(\tilde{\mathbf{x}}_{N+L})$  to the current time interval (see (9)) to obtain  $\mathbf{F}_x^{(N)} = \mathcal{F}^{\text{ext}}(\mathbf{x}_N)$ .

**Real-time output:** Signal representation  $\mathbf{F}_x^{(N)}$   
**end while**

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A. Step 1: Extension by forecasting

a) *Dynamical model and forecasting:* Establishing a dynamical model consists in determining the relation linking  $\mathbf{Y}$  to  $\mathbf{X}$ , that is finding a function  $f$  so that

$$\mathbf{Y} = f(\mathbf{X}) .$$

In a general framework, forecasting means estimating the function  $f$  from the observed values taken by the signal, in order to predict its future values. For instance, the dynamic mode decomposition [25], [18] or other more complicated models [20], [21], [22], [23], allow this by setting additional constraints on the behavior of  $f$ . We will see later, in section III, that it is not necessary to consider such a complex dynamic model for the study of the oscillatory signals of interest to us. That is why we consider here a naive dynamical model, assuming that we have the following relation:

$$\mathbf{Y} = \mathbf{A}\mathbf{X} , \quad (3)$$

where  $\mathbf{A} \in \mathbb{R}^{M \times M}$ . In other words, we adopt a classical strategy in the study of dynamical systems, the *linearization* of a nonlinear phenomenon, when the system is

sufficient regular. Notice that this linearized dynamical model can be written equivalently in function of the sub-signals  $\mathbf{x}_k$ , as:

$$\mathbf{x}_{k+1} = \mathbf{A}\mathbf{x}_k, \quad \forall k \in \{0, \dots, K-1\} . \quad (4)$$

The forecasting method consists in estimating the unknown matrix  $\mathbf{A}$ . Indeed, let  $\tilde{\mathbf{A}}$  denotes the estimate of  $\mathbf{A}$ . We then obtain the forecasting of the signal at time  $\frac{N-1+\ell}{f_s}$  by:

$$\tilde{\mathbf{x}}[N-1+\ell] = \boldsymbol{\alpha}^{(\ell)} \mathbf{x}_K , \quad (5)$$

where  $\boldsymbol{\alpha}^{(\ell)}$  denotes the last row of  $\tilde{\mathbf{A}}^\ell$ ; that is to say,

$$\boldsymbol{\alpha}^{(\ell)} = \mathbf{e}_M^T \tilde{\mathbf{A}}^\ell , \quad (6)$$

where  $\mathbf{e}_M$  is the unit vector of length  $M$  given by  $\mathbf{e}_M = (0 \quad \cdots \quad 0 \quad 1)^T$ .

b) *Model estimation:* To estimate the matrix  $\mathbf{A}$ , we consider the simple but numerically efficient least square estimator. That is, we solve the following problem:

$$\tilde{\mathbf{A}} = \arg \min_{\mathbf{A}} \mathcal{L}(\mathbf{A}) , \quad (7)$$

where the loss function  $\mathcal{L}$  is given by:

$$\mathcal{L}(\mathbf{A}) = \|\mathbf{Y} - \mathbf{A}\mathbf{X}\|^2 = \sum_{k=0}^{K-1} \|\mathbf{x}_{k+1} - \mathbf{A}\mathbf{x}_k\|^2 .$$

Therefore, solving the problem (7), i.e.  $\nabla \mathcal{L}(\tilde{\mathbf{A}}) = \mathbf{0}$ , gives the following estimate  $\tilde{\mathbf{A}}$  of the dynamical model matrix  $\mathbf{A}$ :

$$\tilde{\mathbf{A}} = \mathbf{Y}\mathbf{X}^T(\mathbf{X}\mathbf{X}^T)^{-1} . \quad (8)$$

**Remark 1.** This expression clearly shows that the matrix  $\tilde{\mathbf{A}}$  takes the following form:

$$\tilde{\mathbf{A}} = \begin{pmatrix} 0 & 1 & 0 & \cdots & 0 \\ \vdots & \ddots & \ddots & \ddots & \vdots \\ \vdots & & \ddots & \ddots & 0 \\ 0 & \cdots & \cdots & 0 & 1 \\ \alpha_1 & \cdots & \cdots & \cdots & \alpha_M \end{pmatrix} .$$

Then, except the row vector  $\boldsymbol{\alpha} = (\alpha_1 \cdots \alpha_M)$ , the matrix  $\mathbf{A}$  is fully determined by the dynamical model.

c) *Signal extension:* Since we are building a real-time algorithm, we consider that only the right "side" of the TF representation is affected by boundary effects, the left "side" being fixed since it only concerns the past values of the signal. We therefore construct the extended signal  $\tilde{\mathbf{x}} \in \mathbb{R}^{N+L}$  concatenating the observed signal  $\mathbf{x}$ , and the forward forecasting  $\tilde{\mathbf{x}}_{\text{fw}}$ . We summarize the extension step in Algorithm 2. Notice that we could, if necessary, handle the backward estimation using the same strategy than described above, but applying it to the reverse signal  $\mathbf{x}^r = (\mathbf{x}[N-1] \quad \cdots \quad \mathbf{x}[0])^T$ .

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**Algorithm 2** Signal extension.  $\tilde{\mathbf{x}} = \text{SigExt}(\mathbf{x}, M, K, L)$ 


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**Inputs:**  $\mathbf{x}_N, M, K, L$ 
**Forward forecasting.**

- LS estimation of the forward matrix  $\tilde{\mathbf{A}}_{\text{fw}}$  via equation (8).
- Forward forecasting  $\tilde{\mathbf{x}}_{\text{bw}}$  obtained applying equation (5) with  $\ell \in \{1, \dots, L\}$ .

**Output:** Extended signal  $\tilde{\mathbf{x}}_{N+L} = (\mathbf{x}_N \quad \tilde{\mathbf{x}}_{\text{fw}})^T$ .

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**B. Step 2: Extended Time-Frequency Representation**

Let  $\mathcal{F} : \mathbb{R}^N \rightarrow \mathbb{C}^{F \times N}$  generically denotes the TF representation of interest to us, which could be, for instance, STFT, CWT, SST, or RS. Here,  $F$  typically denotes the size of the discretization along the frequency axis. Due to the boundary effects, the representation  $\mathcal{F}(\mathbf{x}_N)$  shows undesired patterns when approaching its edges. To alleviate the boundary effect, we apply the representation to the estimated extended signal  $\tilde{\mathbf{x}}$ . This strategy moves the boundary effects out of the time interval  $I = [0, \frac{N-1}{f_s}]$ . Finally, the boundary-effects insensitive representation  $\mathcal{F}^{\text{ext}} : \mathbb{R}^N \rightarrow \mathbb{C}^{F \times N}$  of  $\mathbf{x}_N$  is given for all  $v \in \{0, \dots, F-1\}$ ,  $n \in \{0, \dots, N-1\}$  by:

$$\mathcal{F}^{\text{ext}}(\mathbf{x}_N)[v, n] = \mathcal{F}(\tilde{\mathbf{x}}_{N+L})[v, n]. \quad (9)$$

This amounts to restricting the representation  $\mathcal{F}(\tilde{\mathbf{x}}_{N+L})$  to the current measurement interval of  $\mathbf{x}_N$ . For the sake of simplicity, we denote the restriction operator by  $\mathcal{R}$ , where  $\mathcal{R} : \mathbb{C}^{F \times (N+L)} \rightarrow \mathbb{C}^{F \times N}$ . Consequently, we have:

$$\mathcal{F}^{\text{ext}}(\mathbf{x}_N) = \mathcal{R}(\mathcal{F}(\tilde{\mathbf{x}}_{N+L})). \quad (10)$$

We call  $\mathcal{F}^{\text{ext}}$  the *boundary-free TF representation*.

$\mathbf{F}_x^{(N)} \in \mathbb{C}^{F \times N}$  is the estimation of the boundary-free TF representation at iteration  $N$  in Algorithm 1. For numerical purpose, and to make the real-time implementation achievable, we do not perform a full re-estimation of  $\mathcal{F}^{\text{ext}}$ . Instead, the additional knowledge provided by  $\mathbf{x}[N+1]$  only influences the values of the last  $L_{\text{win}}$  columns of  $\mathcal{F}^{\text{ext}}(\mathbf{x}_{N+1})$ , where  $L_{\text{win}}$  denotes the window half-length used by the TF representation. Thus,  $\mathbf{F}_x^{(N+1)}$  is obtained by the concatenation of the first  $N - L_{\text{win}} + 1$  columns of  $\mathbf{F}_x^{(N)}$  with the last  $L_{\text{win}}$  columns of  $\mathcal{F}^{\text{ext}}(\mathbf{x}_{N+1})$ .

### III. THEORETICAL PERFORMANCE

#### A. Signal model

We model the deterministic part of the observed signal as a multicomponent harmonic signal; that is, a sum of sine waves,

$$\mathbf{z}[n] = \sum_{j=1}^J \Omega_j \cos\left(2\pi f_j \frac{n}{f_s} + \varphi_j\right), \quad (11)$$

where  $J$  denotes the number of components,  $\Omega_j > 0$  the amplitude of the  $j$ -th component,  $f_j$  its frequency, and

$\varphi_j \in [0, 2\pi)$  its initial phase. For the sake of simplicity, we make an additional assumption on the frequencies of each component. We assume that for all  $j \in \{1, \dots, J\}$ :

$$\exists p_j, p'_j \in \mathbb{N}^* : f_j = \frac{p_j}{M} f_s = \frac{p'_j}{K} f_s. \quad (12)$$

In addition, the observed signal is assumed to be corrupted by an additive Gaussian white noise. Therefore, the measured discrete signal  $\mathbf{x}$  is written as:

$$\mathbf{x} = \mathbf{z} + \sigma \mathbf{w}, \quad (13)$$

where  $\mathbf{z}$  follows model (11),  $\mathbf{w}$  is a Gaussian white noise, whose variance is normalized to one. Thus,  $\sigma^2$  denotes the variance of the additive noise  $\sigma \mathbf{w}$ .

#### B. Forecasting error

On the forecasting interval, we decompose the estimated signal  $\tilde{\mathbf{x}}$  as follows:

$$\tilde{\mathbf{x}}[n] = \mathbf{z}[n] + \epsilon[n], \quad (14)$$

where  $\epsilon$  is the forecasting error. When  $n \in I = \{0, \dots, N-1\}$ , this error is only containing the measurement noise, that is  $\epsilon[n] = \sigma \mathbf{w}[n]$ . Outside the interval  $I$ , the importance of the forecasting error  $\epsilon$  is also affected by the loss of information resulting from the linearization of dynamical model we consider in (3). To evaluate the actual behavior of the forward forecasting error  $\epsilon[n]$  when  $n \geq N$ , we determine its first two moments.

- 1) The mean, which is also the estimation bias, is such that:

$$\mu[n] \triangleq \mathbb{E}\{\epsilon[n]\} = \mathbb{E}\{\tilde{\mathbf{x}}[n]\} - \mathbf{z}[n].$$

Given the forecasting strategy, we have  $\mu[n] = 0$  when  $n \in I$  and:

$$\mu[n] = \mathbb{E}\{\alpha^{(\ell)}\} \mathbf{z}_K + \sigma \mathbb{E}\{\alpha^{(n-N+1)}\} \mathbf{w}_K - \mathbf{z}[n] \quad (15)$$

when  $n \geq N$ .

- 2) The covariance is given by:

$$\begin{aligned} \gamma[n, n'] &\triangleq \mathbb{E}\{(\epsilon[n] - \mu[n])(\epsilon[n'] - \mu[n'])\} \\ &= \mathbb{E}\{\tilde{\mathbf{x}}[n]\tilde{\mathbf{x}}[n']\} - \mathbf{z}[n]\mathbf{z}[n'] - \mu[n]\mathbf{z}[n'] \\ &\quad - \mu[n']\mathbf{z}[n] - \mu[n]\mu[n']. \end{aligned}$$

Thus by definition of the noise, we have  $\gamma[n, n'] = \sigma^2 \delta_{n, n'}$  when  $(n, n') \in I^2$ . When  $n \geq N$ , let us denote  $\ell = n - N + 1$ . Then, we have two cases.

- (i) If  $n' \in I$ :

$$\begin{aligned} \gamma[n, n'] &= \sigma \mathbb{E}\{\mathbf{w}[n']\alpha^{(\ell)}\} \mathbf{z}_K + \sigma^2 \mathbb{E}\{\mathbf{w}[n']\alpha^{(\ell)}\mathbf{w}_K\} \\ &\quad - \mathbf{z}[n]\mu[n'] - \mu[n]\mu[n']. \end{aligned} \quad (16)$$

- (ii) If  $n' = N - 1 + \lambda \geq N$ :

$$\begin{aligned} \gamma[n, n'] &= \mathbf{z}_K^T \mathbb{E}\{\alpha^{(\ell)T} \alpha^{(\lambda)}\} \mathbf{z}_K + \sigma \mathbb{E}\{\alpha^{(\ell)} \mathbf{w}_K \alpha^{(\lambda)}\} \mathbf{z}_K \\ &\quad + \sigma \mathbb{E}\{\alpha^{(\lambda)} \mathbf{w}_K \alpha^{(\ell)}\} \mathbf{z}_K + \sigma^2 \mathbb{E}\{\alpha^{(\ell)} \mathbf{w}_K \alpha^{(\lambda)} \mathbf{w}_K\} \\ &\quad - \mathbf{z}[n]\mathbf{z}[n'] - \mathbf{z}[n]\mu[n'] - \mathbf{z}[n]\mu[n'] - \mu[n]\mu[n']. \end{aligned} \quad (17)$$

Besides, we recall that  $\gamma[n, n'] = \gamma[n', n]$ .

Expressions (15), (16), and (17) show that these quantities depend on the behavior of the forecasting random vector  $\alpha^{(\ell)}$ . In Lemma 1, we specify the asymptotic behavior of the forecasting vector  $\alpha^{(\ell)}$  when the dataset size  $K$  is great.

**Lemma 1.** Let  $\mathbf{x}$  be a random vector defined by (13). Let  $\alpha^{(\ell)}$  be the associated forecasting vector for the estimation of  $\mathbf{x}[N-1+\ell]$ , given by (6) and obtained from the least square estimation (8). Let  $\alpha_0^{(\ell)}$  be the last row of the matrix  $\mathbf{A}_0^\ell$ , where  $\mathbf{A}_0$  is defined by:

$$\mathbf{A}_0 = \left( \frac{1}{K} \mathbf{Z}' \mathbf{Z}^T + \sigma^2 \mathbf{D} \right) \left( \frac{1}{K} \mathbf{Z} \mathbf{Z}^T + \sigma^2 \mathbf{I} \right)^{-1}, \quad (18)$$

where  $\mathbf{Z} = (\mathbf{z}_0 \cdots \mathbf{z}_{K-1})$ ,  $\mathbf{Z}' = (\mathbf{z}_1 \cdots \mathbf{z}_K)$ ,  $\mathbf{z}_k$  is the  $k$ -th sub-signal extracted from  $\mathbf{z}$  in the same way as  $\mathbf{x}_k$  is defined from  $\mathbf{x}$  in (2), and  $\mathbf{D} \in \mathbb{R}^{M \times M}$  is the Toeplitz matrix such that  $\mathbf{D}[m, m'] = \delta_{m+1, m'}$ .

Let  $\mathbf{h}^{(\ell)}$  be the error vector given by:

$$\mathbf{h}^{(\ell)} = \alpha^{(\ell)} - \alpha_0^{(\ell)}.$$

Then, the random vector  $\mathbf{h}^{(\ell)}$  converges in law to a zero-mean Gaussian random vector when  $K \rightarrow \infty$ , and we have:

$$\sqrt{K} \mathbf{h}^{(\ell)} \xrightarrow[K \rightarrow \infty]{\mathcal{D}} \mathcal{N} \left( \mathbf{0}, \Gamma^{(\ell, \ell)} \right), \quad (19)$$

with  $\Gamma^{(\ell, \ell)} = \mathbf{F}^{(\ell)T} \Gamma_0 \mathbf{F}^{(\ell)}$ , where  $\Gamma_0 \in \mathbb{R}^{M(M+1) \times M(M+1)}$  is a covariance matrix and  $\mathbf{F}^{(\ell)} \in \mathbb{R}^{M(M+1) \times M}$  is a Jacobian matrix. The expressions of  $\Gamma^{(\ell, \ell)}$  does not depend on  $K$  or  $\sigma$ .

*Proof.* See the Supplementary Material. The proof is based on the multivariate delta method (see paragraph 7.2 in [26]), which allows to asymptotically approximate a random vector normal as a Gaussian random vector. ■

Consequently, the covariance between  $\sqrt{K} \mathbf{h}^{(\ell)}$  and  $\sqrt{K} \mathbf{h}^{(\lambda)}$  remains bounded, i.e.:

$$K \mathbb{E} \left\{ \mathbf{h}^{(\ell)T} \mathbf{h}^{(\lambda)} \right\} \xrightarrow[K \rightarrow \infty]{} \Gamma^{(\ell, \lambda)} = \mathbf{F}^{(\ell)T} \Gamma_0 \mathbf{F}^{(\lambda)}.$$

**Theorem 1.** Let  $\mathbf{x} \in \mathbb{R}^N$  be a discrete-time random signal following model (13). Let  $\tilde{\mathbf{x}}$  denotes its forecasting, obtained using the extension Algorithm 2. Let  $n \geq N$  be a sample index. Then, the first-order moment of the forecasting error  $\epsilon[n]$  in (14) is approximated by:

$$\mu[n] \underset{K \rightarrow \infty}{\sim} o(\sigma^2) \quad (20)$$

Its second-order moment  $\gamma[n, n']$  verify the following approximation equations:

(i) if  $n' \in I = \{0, \dots, N-1\}$ :

$$\gamma[n, n'] \underset{K \rightarrow \infty}{\sim} \sigma^2 \alpha_0^{(n-N-1)} [n' - (N-M)] \mathbb{1}_{(n' \geq N-M)} \quad (21)$$

(ii) if  $n' \geq N$ :

$$\gamma[n, n'] \underset{K \rightarrow \infty}{\sim} \frac{1}{K} \mathbf{z}_K^T \Gamma^{(\ell, \lambda)} \mathbf{z}_K + \frac{\sigma^2}{K} \text{Tr} \left( \Gamma^{(\ell, \lambda)} \right) + \sigma^2 \left\langle \alpha_0^{(\ell)}, \alpha_0^{(\lambda)} \right\rangle, \quad (22)$$

where  $\ell = n - N + 1$  and  $\lambda = n' - N + 1$ .

*Proof.* See the Supplementary Material. The proof is mainly based on the results provided by Lemma 1, combined with the Isserlis' theorem [27], which provides a formula for the computation of higher-order moments of Gaussian random variables. ■

Ideally, the forecasting error would behave like the measurement noise  $\sigma \mathbf{w}$ , i.e. a zero-mean noise whose variance is of the order of  $\sigma^2$ . Theorem 1 shows that the forecasting error is asymptotically unbiased. Concerning the covariance of the forecasting error, although equations (21) and (22) are not easily readable, one can evaluate the dependence of the variance in function of the tuning parameters, that are adjusted by the user. Let us focus on the forecasting error variance  $\gamma[n, n]$  when  $n \geq N$ . First, as expected, the variance increases linearly with the noise variance  $\sigma^2$ . Second, it asymptotically depends linearly on the ratio  $\frac{1}{K}$ . This shows the need to use a sufficiently large dataset to obtain an accurate forecast. Third, the dependency on the sub-signals lengths  $M$  and the forecasting index  $\ell = n - N + 1$  is hidden in the expression of the covariance matrix  $\Gamma^{(\ell, \ell)}$ . We discuss this dependency in more detail in section IV-A1.

**Remark 2** (Adaptive Harmonic Model). One can extend the previous result to the case where the instantaneous frequencies and amplitudes of the components of the deterministic part of the observed signal are slowly varying. We therefore handle the AM-FM model which, in its continuous-time version, takes the following form:

$$z(t) = \sum_{j=1}^J a_j(t) \cos(2\pi \phi_j(t)), \quad (23)$$

where  $a_j$  and  $\phi_j'$  describe how large and fast the signal oscillate. Clearly, (11) is a special case satisfying the AM-FM model. It is also clear that if we apply SigExt for the forecasting, the forecasting error is additionally sensitive to the speed of variation of the instantaneous amplitudes  $a_j$  and frequencies  $\phi_j'$ .

In many practical signals, the amplitude and frequency do not change fast. It is thus reasonable to further restrict the regularity and variations of the instantaneous amplitudes and frequencies of these AM-FM functions so that the signal can be "locally" well approximated by a harmonic function in (11). We call thus functions satisfying the adaptive harmonic model (AHM) (see [28], [29] for mathematical details). In this case, we know that when  $K$  is not too large, the signal can be well approximated by (11) and hence Theorem 1 can still be applied to approximate the error. How to determine the optimal  $K$  based on the signal is out of the scope of this paper and will be explored in the future work.

### C. Performance of the boundary effects reduction

Since it is beyond the scope of the present paper, we do not provide here a generic proof of the reduction of the boundary effects on any TF representation. Instead, let us

discuss the particular case of SST. Since SST is designed to analyze signals satisfying the AHM, as is discussed in Remark 2, Theorem 1 ensures that the forecasting error  $\epsilon$  defined in (14) is controlled and bounded in terms of mean and covariance. Recall Theorem 3 in [28], which states that when the additive noise is stationary and remains small, the SST of the observed signal remains close to the ideal SST of an AM-FM signal, throughout the TF plane. In our case, while the additive noise is not stationary, the dependence only appears in the extended part of the signal. Thus, the proof of Theorem 3 in [28] can be generalized to explain the robustness of SST to the noise. We refer the reader to [28] for a precise quantification of the error made in the TF plane, which depends notably on the covariance of the additive noise and on the speed of variation of the amplitudes and instantaneous frequencies composing the signal. Therefore, in our case, this means that boundary effects are strongly reduced since the forecasting error does not impact SST. An immediate application is that the instantaneous frequencies can now be well estimated continuously up to the edges of the TF plane. While we do not provide a theoretical justification, we verify experimentally that algorithm 1 is efficient for a large number of representations in the following section IV.

#### IV. NUMERICAL RESULTS

For the reproducibility purpose, the MATLAB code and datasets used to produce the numerical results in this section are available online at <https://github.com/AdMeynard/BoundaryEffectsReduction>.

##### A. Evaluation the forecasting performance

In that section, we first evaluate the quality of the forecasting step and compare it the the theoretical results provided by Theorem 1. The level of the forecasting error depends on at least two parameters:

- The noise variance  $\sigma^2$ .
- The size of the training dataset  $K$ .

In subsections IV-A1 and IV-A2, we study the influence of these parameters. A comparison with the theoretical results of section III is also available.

1) *Sum of sine waves*: We proved that the linear dynamic model is sufficient to catch the dynamical behavior of signals taking the form (11). In order to validate this theoretical result, we apply the forecasting Algorithm 2 to a large number of realizations of the random vector  $\mathbf{x}$  following the model (13), and such that the deterministic component  $\mathbf{z}$  takes the form:

$$\mathbf{z}[n] = \cos\left(2\pi p_1 \frac{n}{M}\right) + R \cos\left(2\pi p_2 \frac{n}{M}\right), \quad \forall n \in \{1, \dots, N\},$$

with  $N = 10^4$ ,  $M = 150$ ,  $p_1 = 10$ ,  $p_2 = 33$  and  $R = 1.4$ . Besides, the additive noise is chosen to be Gaussian:  $\mathbf{w} \sim \mathcal{N}(\mathbf{0}, \mathbf{I})$ .

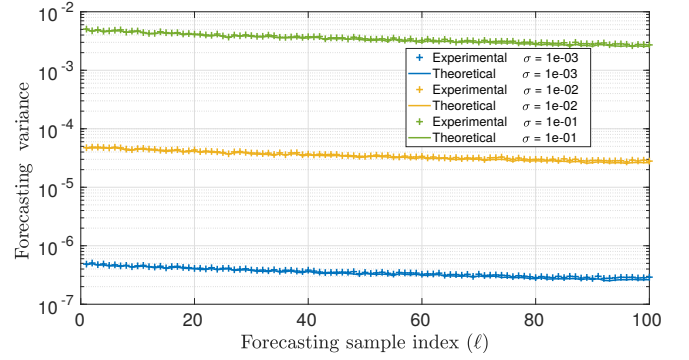


Figure 2. Evolution of the experimental and theoretical forecasting variance in function of the forecasting sample index for different values of  $\sigma$ .

a) *Influence of the noise variance  $\sigma^2$* : Here, the size of the training dataset is set to  $K = 450$ . Then, the forecasting algorithm is run on 1000 realizations on the discrete signal  $\mathbf{x}$  for three different values of  $\sigma$ , logarithmically equi-spaced from  $10^{-3}$  to  $10^{-1}$ . For each of these values, we determine the experimental bias, denoted as  $\mu_{\text{xp}}[N - 1 + \ell]$ , and experimental variance, denoted as  $\gamma_{\text{xp}}[N - 1 + \ell, N - 1 + \ell]$ , in function of the forecasting sample index  $\ell$  (going from 1 to  $L = 100$ ).

The experimental results show that the bias is neither depending on the noise variance  $\sigma^2$  nor the forecasting length  $\ell$ . Indeed, independently of  $\sigma$ , we always have  $\mu_{\text{xp}}[N - 1 + \ell] \in [-0.03\sigma, 0.03\sigma]$ , which is negligible with respect to the magnitude of  $\mathbf{z}$ . This result confirms the theoretical result (20).

On Fig. 2, we display the experimental variance  $\gamma_{\text{xp}}[N - 1 + \ell, N - 1 + \ell]$  for each value of  $\sigma$ . The associated theoretical asymptotic forecasting variance (22) is also displayed in solid line. As expected, this result highlights the fact that the forecasting variance increases linearly with respect to  $\sigma^2$ . Surprisingly, this result shows that the forecasting variance *slightly* decreases with  $\ell$ , what is counterintuitive.

b) *Influence of the training dataset size  $K$* : Here, the noise variance  $\sigma$  is set to  $\sigma = 10^{-2}$ . Then, the forecasting algorithm is run on 3000 realizations on the discrete signal  $\mathbf{x}$  for three different values of  $K$ , logarithmically equi-spaced from  $4.5 \times 10^2$  to  $2 \times 10^3$ . For each of these values, we determine the experimental bias  $\mu_{\text{xp}}[N - 1 + \ell]$  and variance  $\gamma_{\text{xp}}[N - 1 + \ell, N - 1 + \ell]$  in function of the forecasting sample index  $\ell$  (going from 1 to 500).

As in the previous study, the experimental bias vanishes when  $K$  increases, what confirms the approximation result (20). Besides, the experimental variance is displayed on Fig. 3, and compared with the associated theoretical variance (22). Each color corresponds to these experimental results obtained for a given value of  $K$ . This result validates the asymptotic behavior provided by (22), and we can show that the second order moment  $\gamma[\ell, \ell]$  is less dependent on  $K$  when  $K$  is large.

c) *Summary*: Both previous experimental results combined with the theoretical asymptotic equation (22)



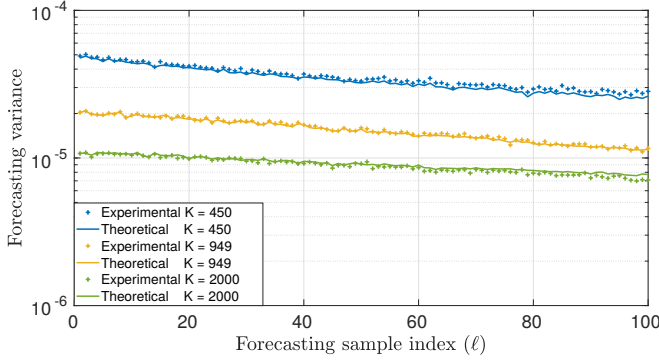


Figure 3. Evolution of the experimental and theoretical forecasting variance in function of the forecasting sample index for different values of  $K$ .

allow us to describe the influence of the noise variance and the size of the training dataset on the variance of the forecasting noise, empirically summarized as follows:

$$\gamma[N-1+\ell, N-1+\ell] \underset{K \rightarrow \infty}{\approx} \frac{\sigma^2}{K} g[\ell]. \quad (24)$$

where  $g$  is a bounded positive function. The empirical result is coherent with the theoretical result provided by Theorem 1.

This study neglects the analysis of the influence of the parameter  $M$ , whose influence on the value of the experimental variance is numerically not significant as long as  $M \ll 2K$ . The choice of this parameter is especially crucial when the deterministic component of the signal is no longer stationary. The AHM, discussed below, is an example.

2) *Adaptive harmonic model*: We now consider a signal satisfying the AHM so that the instantaneous frequencies and amplitudes of its components vary over time. The deterministic component  $\mathbf{z}$  of the random vector  $\mathbf{x}$  (constructed following the model (13)) takes the following form, for all  $n \in \{1, \dots, N\}$ :

$$\mathbf{x}[n] = \cos(2\pi\phi_1[n]) + R[n] \cos(2\pi\phi_2[n]),$$

where the instantaneous amplitude  $R$  is given by:

$$R[n] = 1.4 + 0.2 \cos\left(4\pi \frac{n}{N}\right),$$

and the instantaneous phases are such that:

$$\begin{aligned} \phi_1[n] &= \frac{p_1}{M} \left( n + \frac{0.01}{2\pi} \cos\left(2\pi \frac{n}{N}\right) \right) \\ \phi_2[n] &= p_2 \frac{n}{M} + \frac{20}{2Nf_s} n^2 \end{aligned}$$

Besides, the noise is chosen to be Gaussian:  $\mathbf{w} \sim \mathcal{N}(\mathbf{0}, \mathbf{I})$ , and we take:  $N = 10^4$ ,  $M = 750$ ,  $p_1 = 10$ ,  $p_2 = 23$ .

To highlight the fact that the linear dynamical model is sufficient to catch most of the dynamical behavior of signals following the AHM, we compare the performance of the Algorithm 1 with a simple extension obtained by pointwise symmetrization [15]. We also evaluate the performance of reference forecasting algorithm that could be used for extending such signals. These methods are:

Table I  
AHM SIGNAL. PERFORMANCE OF THE EXTENSION METHODS.

Algorithm	MSE		Computing time (sec.)
	Mean	Standard deviation	
SigExt	$1.433 \times 10^{-3}$	$4.361 \times 10^{-4}$	0.152
Symmetric	$1.019 \times 10^1$	$1.192 \times 10^2$	0.002
EDMD	$3.076 \times 10^{-2}$	$8.095 \times 10^{-2}$	2.537
GPR	$1.436 \times 10^{-3}$	$4.346 \times 10^{-4}$	146.331
TBATS	$1.732 \times 10^{-3}$	$4.924 \times 10^{-4}$	1837.120

- The EDMD has been developed by Williams *et al.* [18]. The proposed algorithm is a way to obtain an approximation of the so-called Koopman operator of the observed system, which theoretically allows to catch dynamic of nonlinear systems [30].
- The GPR [19] is a method relying on a probabilistic dynamical model. That one is based on the Gaussian process structure, and therefore offer more flexibility in the type of dynamic that could be modeled than the linear model (3).
- The TBATS method [21] is based on a classical decomposition of times series into a trend, a seasonal and an ARMA components, with a specific dynamic for the seasonal component. This model demands the estimation of numerous parameters and, by implication, may be slow.

To quantify the global quality (i.e. not depending on  $\ell$ ) of the forecasting approaches, we evaluate the Experimental Mean Square Error  $\text{MSE}_{\text{xp}}(\tilde{\mathbf{x}})$  of the forward forecast extended signals, namely:

$$\begin{aligned} \text{MSE}_{\text{xp}}(\tilde{\mathbf{x}}) &= \frac{1}{L} \|\tilde{\mathbf{x}} - \mathbf{x}^{\text{ext}}\|^2 \\ &= \frac{1}{L} \sum_{\ell=1}^L \mu_{\text{xp}}[N-1+\ell]^2 + \gamma_{\text{xp}}[N-1+\ell, N-1+\ell]. \end{aligned} \quad (25)$$

where  $\mathbf{x}^{\text{ext}}$  is the ground-truth extended signal, that is:  $\mathbf{x}^{\text{ext}} = (\mathbf{x}[-L] \ \dots \ \mathbf{x}[N-1+L])$ . Then, as long as the bias  $\mu[N-1+\ell]$  and the variance  $\gamma[N-1+\ell, N-1+\ell]$  of the forecasting estimator remain small for all  $\ell$ , the MSE takes small values either. Corresponding results are given in Table I. They show that the naive extension we propose gives satisfying results, in particular in comparison with the point-symmetric extension. Besides, even though the other more sophisticated methods, like GPR, give MSE values that have a slightly smaller standard deviation, these methods are substantially limited by the computing time they require, which prevent them from being used to exploit real-time data. Thus, SigExt is the extension method that optimize the trade-off between the forecasting quality and the computing time.

### B. Evaluation of the quality of the boundary effects reduction

1) *Metrics*: The quality of the boundary effects reduction is evaluated directly on the TF representation. To that aim, we compare the obtained representation to the optimal representation  $\mathcal{F}_N^{\text{opt}}(\mathbf{x})$ , defined as the restriction

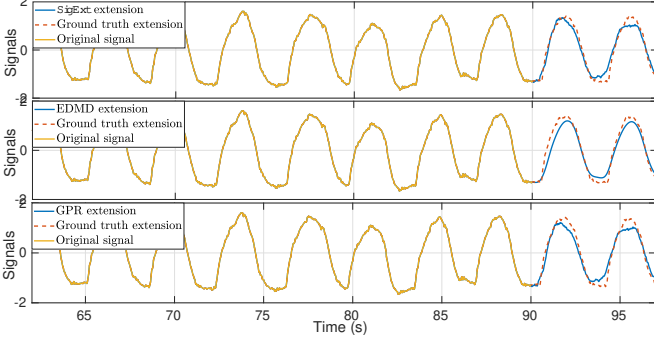


Figure 4. Extended THO signal (blue) obtained by the SigExt forecasting (top), the EDMD forecasting (middle), and the GPR forecasting (bottom), superimposed with the ground truth signal (red dash).

of the representation of the ground-truth extended signal  $\mathbf{x}^{\text{ext}}$ . Therefore, we have:

$$\mathcal{F}^{\text{opt}}(\mathbf{x}_N) = \mathcal{R}(\mathcal{F}(\mathbf{x}^{\text{ext}})) . \quad (26)$$

In the aim of comparing the different techniques, we use a criterion, proposed in [29], that quantify the distance between a given TF representation and the optimal one. It is built in analogy with the optimal transport distance, which enables quantifying the distance between two probability density functions. Let us generically denote a time frequency representation  $\mathcal{Q}$ . Then, for  $t$  fixed, we consider the following probability density function:  $p_{\mathcal{Q}}^t(\xi) = |\mathcal{Q}(\xi, t)|^2 / \int_{\mathbb{R}} |\mathcal{Q}(v, t)|^2 dv$ . At each instant  $t$ , we can then determine the optimal transport distance  $d_t$  between the two densities. It is given by the  $L^1$  norm of the difference between the associated distribution functions. In other words, we have:

$$d_t(\mathcal{Q}, \mathcal{F}_0) = \int_{\mathbb{R}} |\tilde{P}_{\mathcal{Q}}^t(\xi) - P_{\mathcal{F}_0}^t(\xi)| d\xi ,$$

where  $P_{\mathcal{Q}}^t(\xi) = \int_{-\infty}^{\xi} p_{\mathcal{Q}}^t(v) dv$  and  $\tilde{P}_{\mathcal{F}_0}^t(\xi) = \int_{-\infty}^{\xi} \tilde{p}_{\mathcal{F}_0}^t(v) dv$ . The *optimal transport distance* quantifies the proximity between the estimated and actual instantaneous frequencies while favoring the sparsity of the estimated TF representation. That is why, the performance index  $D(\mathcal{Q})$  for the reduction of boundary effects of a given TF representation  $\mathcal{Q}$  is given by the averaged ratio between the optimal transport distance to the optimal TF representation (26) and the optimal transport distance of to the optimal TF representation, that is:

$$D(\mathcal{Q}) = \frac{1}{|I|} \int_I \frac{d_t(\mathcal{Q}, \mathcal{F}^{\text{opt}})}{d_t(\mathcal{F}, \mathcal{F}^{\text{opt}})} dt . \quad (27)$$

Thus,  $D(\mathcal{Q}) < 1$  means a reduction of the boundary effects. Let us evaluate the quality of the boundary effects reduction on biomedical signals.

2) *Respiratory signal*: We first consider a respiratory signal of 6 hours 20 minutes. This signal, a small portion of which is displayed in Fig. 4, is sampled at  $f_s = 100$  Hz.

From that large signal, we build a dataset of 378 non-overlapping signals of 60 seconds, i.e.  $N = 6000$ . On

each of these pieces of signal, we implement the forecasting method introduced in section IV-A2, including the SigExt method detailed in Algorithm 2. However, the TBATS extension method is not implemented here because of its excessive computing time. The extensions of 7 seconds-long on each boundary, corresponding to  $L = 700$ . Thus, in order to catch slowly varying dynamical behaviors, the size of the training signal  $M$  is chosen so that  $M = \lfloor 1.5L \rfloor$ . As a result of section IV-A1, we take:  $K = \lfloor 2.5M \rfloor$ . The resulting MSE (25) averaged with respect to the simulations are given in Table II. The MSE of SigExt is higher than the MSE of the other methods. This is mainly caused by the presence of a few segments of the whole signal involving complicated behaviors, that are unpredictable via a too simple dynamical model like (3). The left of Fig. 5 illustrates one of those cases, where SigExt fails to catch the fast varying dynamic of the instantaneous amplitude to satisfactorily forecast the signal. The EDMD and Symmetric extensions are more robust to those situations, as shown on this example and in Table II. Nevertheless, SigExt provides a sufficiently relevant extension to give TF representations sparingly affected by boundary effects. On the right of Fig. 5, we display comparison between the right boundary of SST of the same segment of signal (top-right), and its boundary-free SST obtained after the SigExt forecasting (bottom-right). The extension of the instantaneous frequency visible on the right side of the image, illustrates the reduction of boundary effects produced despite an inaccurate signal forecasting.

We then apply BoundEffRed (i.e. Algorithm 1) for diverse TF representations: STFT, SST, RS, as well as concentration of frequency and time (ConceFT), a generalized multitaper SST-based representation introduced in [29]. In Table II, we give the averaged performance index (27), evaluated the whole TF representations (including boundaries). Even though SigExt performs somehow moderately, the boundary effects are dramatically reduced on the TF representations, in the same order of magnitude than with the forecastings given by EDMD or GPR. Notice that the extension length  $L$  has been set accordingly to the window length used by the TF analysis tool. For instance, the window length used to evaluate the STFT is of 1500 samples. To prevent the STFT from being sensitive to the boundaries, we set  $L = 750$ . In this way, the evaluation of the spectral content of the signal near its boundaries is not limited by a lack of information all along the window support. From now on, all results are given for  $L$  equal to the half of the width of the window used in the TF transform.

Finally, in order to verify the ability of BoundEffRed to be implemented in real time, we evaluate the computation time required for each iteration of Algorithm 1, corresponding to the update of the boundary-free TF representation. This check is performed on the real-time SST of a 10-minute sample of the respiratory signal. As the SST is temporally sub-sampled by a ratio of 10, the update should not exceed  $10/f_s = 0.1$  sec. This is indeed



Table II  
PERFORMANCE OF THE EXTENSION METHODS ON A RESPIRATORY SIGNAL.

Algorithm	Averaged MSE	Averaged performance index $D$			
		STFT	SST	RS	ConceFT
SigExt	0.046	0.705	0.698	0.694	0.776
Symmetric	0.039	1.280	0.984	9.063	1.295
EDMD	0.022	0.757	0.635	0.701	0.793
GPR	0.045	0.817	0.793	0.727	0.876

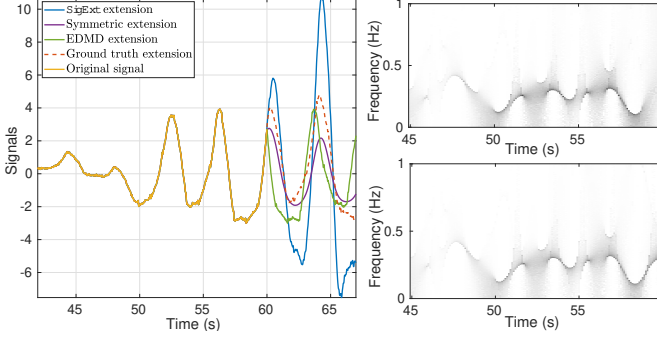


Figure 5. Extensions of a segment of the respiratory signal (left) where SigExt is outperformed by the EDMD and Symmetric extensions. Corresponding SST (top-right) and boundary-free SST obtained with SigExt (bottom-right).

the case, as no iteration lasted longer than xx sec on a 6-Core Xeon CPU running at 3.5 GHz and 64 GB of RAM.

3) *Photoplethysmogram*: We perform a study similar to the previous one on a 640 second-long photoplethysmogram (PPG) signal extracted from the Physionet dataset [31], [32], sampled at  $f_s = 125$  Hz. A 32 second-long piece of this signal is displayed on the top of Fig. 6. The estimated 2-second extension obtained by SigExt on both boundary of this signal is superimposed to the ground-truth signal in the bottom of Fig. 6.

We divide the signal into 32-second long segments, and apply Algorithm 1 on each piece. We provide in Table III the performance index  $D$  of the boundary-free TF representations averaged over the signals. For all the considered TF representations, the results clearly shows that our algorithm reduces the boundary effects. Even

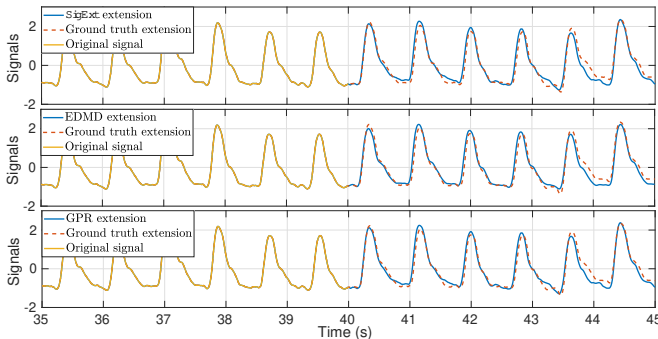


Figure 6. Extended PPG signal (blue) obtained by the SigExt forecasting (top), the EDMD forecasting (middle), and the GPR forecasting (bottom), superimposed with the ground truth signal (red dash).

Table III  
PPG SIGNAL: AVERAGED PERFORMANCE INDEX OF THE BOUNDARY-FREE TF REPRESENTATIONS FOR DIFFERENT EXTENSIONS.

Extension method	Performance index $D$		
	STFT	SST	ConceFT
SigExt	0.774	0.048	X
Symmetric	0.818	0.059	X
EDMD	0.566	0.038	X
GPR	0.735	0.045	X

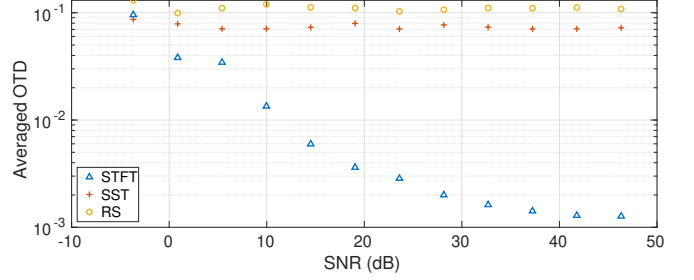


Figure 7. PPG signal. Averaged OTD of BoundEffRed in function of the SNR.

though, on this signal, the SigExt extension yields TF representations slightly more sensitive to boundary effects than the extensions given by EDMD or GPR, it is the only technique that allows a real-time implementation.

On the bottom-right panel of Fig. 1, we display the SST resulting from the BoundEffRed strategy, applied to the portion of PPG displayed on Fig. 6. We clearly observe an improvement of the quality of the SST near boundaries. Indeed, the blurring visible when zooming on the right boundary of the SST has almost vanished. The real-time tracking of the instantaneous frequencies contained in the measured signal is therefore largely facilitated.

To evaluate the influence of the noise level on the performance of BoundEffRed, we artificially add a Gaussian noise to the measured PPG signal. It is thus an additional noise to the measurement noise actually contained in the signal. Fig. 7 shows the averaged OTD of BoundEffRed for different values of the Signal to Noise Ratio (SNR). We notice that STFT is slightly more sensitive to noise than SST or RS, and STFT performs the best when the added noise is small, i.e. when the SNR is great.

## V. CONCLUSION

In this paper, we propose an algorithm, named BoundEffRed, for the reduction of boundary effects on TF representations in real-time. This method is based on an extension of the signal obtained by a simple-minded and numerically efficient forecasting. We have shown theoretically that the chosen dynamic model is sufficient for the extension of signals formed by a sum of sine waves. Moreover, the numerical results show that this strategy allows us to switch to a real-time implementation of BoundEffRed is possible, due to the low computation time, unlike other existing forecasting methods. The numerical results also confirmed the robustness to noise

of BoundEffRed, as well as its ability to be applied to many time-frequency representations.

Various improvements can be considered to make the algorithm more robust. In particular, we have noticed (see Fig. 5) that when the regular oscillations of the observed signal break, the forecasting is no longer relevant, and the TF representations are not ideal. A preliminary step should then be added to the algorithm to detect the portions to be rejected, for which a better forecasting scheme is needed. A more fundamental improvement would be to perform the prediction step directly in the TF plane, or even with the aim of specifically predicting the evolution of each instantaneous frequency. This would lead to seeing a TF representation as a multivariate time series (corresponding to each frequency band). We will explore these possibilities in our future work.

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