Nonstationary signal extraction based on bat algorithm sparse decomposition technique

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For nonstationary time series i.e. natural electromagnetic field and acoustical signal, effective signal extraction always requires prior knowledge or hypothesis, and hardly do without artificial judgment. We proposed bat algorithm sparse decomposition (BASD) to realize adaptive recognition and extraction of nonstationary signal in a noisy background. We designed two general atoms for typical signals, and developed dictionary training method based on correlation detection and Hilbert transform. The sparse decomposition was turned into an optimizing problem by introducing bat algorithm with optimized fitness function. By contrast with variational modal decomposition, it was indicated that BASD can effectively extract short time target without inducing global aliasing of local feature, and no preset mode number and late screening were needed.

Signal extraction and signal-noise separation are always two of the research focuses in modern signal processing 1, which are commonly used in biomedical signal features extraction, vibration signal analysis, seismic signal detection 2, sound signals recognition 3, et al. Fourier transform is one of the most classic signal analysis and extraction method, but it cannot accurately describe nonstationary signals 4. In recent years, new theories and technologies continue to appear in signal extraction. For example, wavelet decomposition (WD) 5, 6, empirical mode decomposition 7, Hilbert-Huang Transform (HHT), variational mode decomposition (VMD) algorithm 8. These methods need to meet certain conditions to work, for example the decomposition levels, modal number, and termination thresholds.

To achieve a more flexible, concise and adaptive signal decomposition, researchers proposed sparse decomposition. This method represents the signal with as few atoms as possible in a given redundant dictionary by matching pursuit (MP) algorithm, which is a greedy algorithm for sparse decomposition. Various new evaluation criteria and basis pursuit, orthogonal matching pursuit algorithm (OMP), and time-frequency spectrum segmentation methods 9 were generated to select a set of optimal atoms from the constructed over-complete dictionary.

Aiming at two main research hotspots including sparse decomposition algorithm and over-complete atom dictionary of signal sparse decomposition 10, we designed two typical universal signal atoms, and proposed an adaptive feature-based atom construction method. Redundant dictionary was obtained by extending the feature-based atoms, which can balance the completeness and redundancy of the redundant dictionary. A signal matching tracking extraction algorithm was developed based on the bat algorithm, which could effectively realize nonstationary time domain signal extraction.

Methods

Bat algorithm sparse decomposition

General atomics designed for typical signals. For sinusoidal-like and one-sided decaying oscillatory signals, g-atoms were constructed:

\[
g(c, d, t_1, t_2, f, \phi) = \begin{cases} 
    c e^{(-|d|^4)} \sin(2\pi f (t-t_1) + \phi), & t \in [t_1, t_2] \\
    0, & \text{others}
\end{cases}
\]

(1)

where \( c \) was the normalization factor to ensure that the original signal had the same energy as its sparse decomposition results; \( d \) was the attenuation factor; \( t \) was the sampling time; \( t_1 \) was start point of atomic appearance; \( t_2 \) was the ending point; \( f \) was frequency and \( \phi \) was
The attenuation factor $d=0$, the g-atom degenerated to standard sine wave; when $d$ increased, the g-atom performed sinusoidal damped oscillation. Therefore, this atom had a strong match with sinusoidal signals, and single-sided oscillatory decay signals.

For kind of triangle waves, charge-discharge waves, and bilateral decay oscillation signals, $tr$-atoms were constructed:

$$t_r(c, d_1, d_2, t_0, t_1, t_2, f, \phi, \sigma) =\begin{cases} 
  c e^{-d_1(t - t_0)} \sin(2\pi f (t - t_0) + \phi), & t \in [t_0, t_1] \\
  0, & \text{others}
\end{cases} \quad (2)$$

where $d_1$ and $d_2$ were the bilateral damping factors; $t_0$ was the bilateral boundary of the atomic; $[t_1, t_2]$ was the atomic time range; and $\sigma$ was the bilateral scaling factor.

The time domain waveforms of $tr$-atoms with different parameters were shown in Fig. 2. When the bilateral scaling factor $\sigma=0$, the $tr$-atom degenerated to single-sided oscillating atom (reverse-order g-atom); when $0<\sigma<1$ and the atomic frequency was low enough, the $tr$-atom behaved as a charge-discharge triangle wave; when $\sigma=1$, the $tr$-atom with low-frequency behaved as a triangle-like wave, and behaved as bilateral oscillating decay signal with high-frequency.

The above analysis showed that the constructed g-atoms and $tr$-atoms were very flexible and could matched almost typical testing signals by parameter adjustment.

**Dictionary training algorithm.** A dictionary training algorithm based on feature parameters was studied to determine the key parameters of feature base-atom and to balance the completeness and redundancy for redundant dictionary library construction.

When constructing the redundancy dictionary, we firstly used the short-time Fourier transform to initially determine the frequency $f_s$ and phase $\phi_s$ of the target signal in the original data $x$. Then, we constructed standard sine atom $s_0=\sin(2\pi f_s t + \phi_s)$. The correlation detection technique was used by calculating the correlation function of the original data and the sine atom, and the upper and lower envelopes of the correlation function were obtained by searching the positive and negative peaks of the correlation function. The points of the maximum positive gradient between the positive and negative peaks were the extracted range of the target signal. Thus, the signal sample $E_i$ in the original observation sequence $x$ was extracted. The time information section obtained by the above calculation determined the time domain parameters such as $t_0$, $t_1$ and $t_2$ of the characteristic atoms.

Accurate time and frequency domain parameters were obtained by Hilbert transform of $E_i$:

$$R(t) = E_i(t)\ast h(t) \quad (3)$$

Where: $h(t)$ was the Hitch transform factor.
Complex analytic signal as follows was constructed:

$$z(t) = x(t) + iR(t) = Ae^{i\phi(t)}$$  \hspace{1cm} (4)$$

Where, \(A(t)\) was the amplitude function:

$$A(t) = \sqrt{E_i^2(t) + R^2(t)}$$  \hspace{1cm} (5)$$

And, \(\phi(t)\) was the phase function:

$$\phi(t) = \arctan \left( \frac{R(t)}{E_i(t)} \right)$$  \hspace{1cm} (6)$$

The instantaneous frequency of \(E_i\) was given by (7):

$$f(t) = \frac{d\phi(t)}{dt}$$  \hspace{1cm} (7)$$

The base-atom was obtained with the time information gained by the correlation detection and localization algorithm and the time-frequency parameter information obtained by the Hilbert transform as the reference. And the redundant dictionary of this feature atom was constructed by performing equal-step discrete expansion of the time-frequency parameters on both sides of the reference values.

$$\mathcal{G} = \{ g_j^i | j = 1, 2, \ldots, K \}$$  \hspace{1cm} (8)$$

Where, \(\mathcal{G}\) denotes the set of atoms consisting of extensions of the characteristic base-atom \(g_i^0\). Atoms had the same length as the corresponding sample \(E_i\).

### Signal extraction based on bat algorithm sparse decomposition

Signal sparse decomposition represented a signal by specific combinations of some atoms in a dictionary. For a given dictionary, the optimal combination can be accurately determined when all possible combinations were calculated. However, exhausting all combinations in a dictionary is a non-deterministic polynomial problem that is almost impossible to achieve for large dictionary bases.

So, the requirement was changed to finding a suboptimal combination from the dictionary with the lowest possible number of atoms and the smallest possible extraction error. This will reduce the computational complexity significantly, and the MP algorithm is one of the algorithms that can achieve this requirement.

Assume that the represented signal is \(x\) with length of \(N\). Let \(\mathbb{R}\) denote the Hilbert space in which a dictionary matrix \(D\) composed of a set of vectors \(\{g_1, g_2, \ldots, g_N\}\). Each vector is an atom with the same length \(N\) and these vectors have been treated as normalized as \(\|g_i\|^2 = 1\).

With \(\xi = x\), the MP algorithm selects one atom at a time from the dictionary matrix \(D\) that best matches \(x\), satisfying (9),

$$\mathcal{I}_{bsel} = \max_{a_i \in \mathbb{R}} \left[ x, g_i \right]$$  \hspace{1cm} (9)$$

Where \(\mathcal{I}_{bsel}\) is the index of the best matching atom in \(D\).

The signal \(x\) is then decomposed into two parts, a sparse approximation \(\hat{x}\) and an approximation residual \(\xi_\hat{x}\):

$$x = \hat{x} + \xi_\hat{x} = \left[ x, g_{\mathcal{I}_{bsel}} \right] g_{\mathcal{I}_{bsel}} + \xi_\hat{x}$$  \hspace{1cm} (10)$$

Continues to select the atoms that best matches \(\xi_\hat{x}\), iterating repeatedly and eventually the signal \(x\) can be approximated as a linear sum of these atoms:

$$\hat{x} = \sum_{i=1}^{n} \left[ x, g_i \right] g_i, \quad x - \hat{x} = \xi_\hat{x}$$  \hspace{1cm} (11)$$

For MP algorithm, the non-orthogonality between the vertical projection of the signal (or residuals) on the selected atoms and the residuals will lead to suboptimal iterative results instead of the best optimal, and convergence requires many iterations. The OMP algorithm is the orthogonalization of all selected atoms at each step of the decomposition, which makes the convergence faster with the same accuracy requirement. The convergence process of MP and OMP were described by a dictionary \(D\) with length of three, as shown in Fig. 3. However, although the OMP algorithm reduced iterations to some extent, it had to calculate the current residual and the inner product of all atoms within the current dictionary during each iteration, resulting in unsatisfied effectiveness. Therefore, this paper
introduced the bat algorithm (BA) to optimize the matching tracking algorithm.

![Figure 3. Convergence process of MP and OMP.](image)

**Bat algorithm optimized sparse decomposition algorithm**

The optimization-seeking process can be viewed as a global optimization problem. In order to solve the problems of large computation and low efficiency of existing matching tracking algorithms, the adaptive matching tracking algorithm, bat algorithm sparse decomposition (BASD), with fast convergence and accurate approximation was studied by introducing BA into the existing matching tracking extraction algorithm. The basic flow of BASD algorithm was as follows:

1. Initialize bat population number of bats $N_{pop}$, search space dimension $q$, initialize flight frequency $\hat{p} = \{p_i^0|i=1,2,...,q\}$, acoustic loudness $A^p = \{A_i^p|i=1,2,...,q\}$ and pulse emission frequency $\hat{r} = \{r_i^0|i=1,2,...,q\}$. The initial location of the randomly generated bat colony was randomly generated according to (12):

   $$P^0 = [p_i^0|i=1,2,...,q] = P_{max}^0 + \lambda_1^0 \times (P_{min}^0 - P_{max}^0)$$

   (12)

   where: $\xi_i$ was the random number satisfying the Gaussian observation matrix. $A_n$ was obtained by selecting $q$ atoms $g_i$ from the candidate dictionary $D$ according to the bat group position $P_i$.

   The best matching atom $g_{global}$ in $A_n$ was determined by the fitness function. In this paper, for noise-containing signal extraction, the fitness function of the traditional sparse decomposition was improved to take the $\ell_2$-norm of the residual as the fitness function, and the smaller the fitness indicated that the residual sequence contains smaller effective signal components and higher signal-to-noise separation, so the optimal individual bat position $P_b$ was determined and saved according to (13).

   $$\text{fit}_i(g_{max}) = \arg \min_v \| \xi_v - \tilde{x}_i - \{g_i*\tilde{x}_i\}g_{r_i} \|_2$$

   (13)

   Where

   $$\tilde{x}_i = \tilde{x}_i - \{g_i*\tilde{x}_i\}g_{r_i}, g_{r_i} \in A_i$$

   (14)

2. Update the velocity and position of the individual bat:

   $$f_i = f_{max} + \xi \times (f_{max} - f_{min})$$

   $$v_i = \nu_v + (P_i - P_{max}) \times f_i$$

   (15)

   where $\eta_i$ was a random number, satisfying $\eta_i \in [0,1]$, $f_i$ was the search pulse frequency of the $i$-th bat; $v_i$ denoted the velocity of the $i$-th bat at $j$-th iteration, $P_i$ denoted the position of the $i$-th bat at $j$-th iteration; and $P_0$ was the current global optimal solution.

3. Generate a random number $r_j \in [0,1]$ for each bat and update bat position according to (16).

   $$\begin{cases} 
   \text{global optimization:} & P_i^{t+1} = P_i + \nu_v, r_j \leq r_j^* \\
   \text{local optimization:} & P_i^{t+1} = P_i + \lambda_2 \times (P_{max} - P_{min}), r_j > r_j^* 
   \end{cases}$$

   (16)

   where: $\eta$ was a random number, satisfying $\eta \in [-1,1]$ and $\bar{A}$ was the mean value of fitness of the bat population at $t$ moment.

4. Update the fitness

   $$\text{fit}_i^{t+1} = \| \xi_v - \{g_i*\tilde{x}_i\}g_{r_i} \|_2$$

   (17)
(6) For each bat, a random number $r_j$ is generated, and update the position:

$$
\begin{cases}
    P_j^{n+1} = P_j^{n}, & r_j > A_j^n & \& \& \text{fit}^{n+1}_j < \text{fit}^n_j \\
    P_j^{n+1} = P_j^n, & \text{fit}^{n+1}_j = \text{fit}^n_j & \text{otherwise}
\end{cases}
$$

(18)

(7) The fitness and pulse emission frequency were updated by (19):

$$
\begin{align*}
    r_j^{n+1} &= r_j^n (1 - e^{-\lambda}) \\
    A_j^{n+1} &= \lambda A_j^n
\end{align*}
$$

(19)

where, $\lambda \in (0,1), \gamma > 0$, when $n \to \infty$, $A_j^n \to 0$, $r_j^n \to r_0$.

(8) Find the current matching atom based on the location of the optimal bat individual.

The random perturbation of the current optimal solution in step 4 can effectively avoid the iterative result from falling into a local optimal solution, which helped to find the global optimal solution faster compared to the particle swarm algorithm (PSO).

The Ackley function was used to test the BA and PSO. The expression of the Ackley function was as follows:

$$
\begin{aligned}
    f(x) &= -c_1 \exp \left( -0.2 \sqrt{\frac{1}{n} \sum_{i=1}^{n} x_i^2} \right) - \exp \left( \frac{1}{n} \sum_{i=1}^{n} \cos (2\pi x_i) \right) + c \\
    \text{(20)}
\end{aligned}
$$

In this study, $n=2$, $c=20$, $c=2.71289$. the Ackley function was taken as the fitness function and the global minimum of this function was searched using BA and PSO. The search paths and results were shown in Fig. 4.

**Figure 4.** the optimal trajectory of PSO and BA. (a) and (b) showed the 3D view and contour attempt of the optimal trajectory, respectively. The black and red solid lines in the figures were the merit-seeking trajectories of represented PSO and BA, respectively.

The detailed values were shown in Table 1. The comparison of the tracking trajectory and the optimization results showed that BA had higher convergence speed and computational accuracy than PSO because the gradient of the optimization deviation was larger and the optimization results were closer to the true value.

<table>
<thead>
<tr>
<th>Algorithm</th>
<th>Search results</th>
<th>Optimal solution</th>
</tr>
</thead>
<tbody>
<tr>
<td>PSO</td>
<td>[0.0011 -0.062]</td>
<td>0.2773</td>
</tr>
<tr>
<td>BA</td>
<td>[0.0082 -0.0088]</td>
<td>0.003789</td>
</tr>
</tbody>
</table>

The ideal search results were [0 0] and ideal solution is 0.

**Experiments**

We constructed a nonstationary signal $x$ to test the methods described above: $x = s + ns$, where $s$ is nonstationary target signal including pulse signal and partial discharge signal at random positions, and $ns$ is background noise.

The finesses of $x$ in the space $\mathbb{R}$ spread by different key parameters were calculated according to (13), as shown in Fig. 5.
Figure 5. Fitness values of the original signal over the entire dictionary.

Figure 5 showed that the global fitness was a multipolar, nonlinear function. When extracted s from x by BASD, the flow chart was shown in Table 2.

TABLE 2 Signal extraction based on BASD

<table>
<thead>
<tr>
<th>Input: x, f_s</th>
</tr>
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<tbody>
<tr>
<td>Defined the maximum iterations K, optimal threshold ε_thre, which was determined by the l-2 norm of the original segment without the target signal</td>
</tr>
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</table>

1: Initialization: K=1, n=1; z_0=x; fitness_{n=1}=1E6;

2: while (k<K&|z_n|^2>fitness_{n=1})

3: determined time-frequency parameters of \( \xi_n \) by dictionary training:

\[
pare=[f_0 \ \phi_0 \ d_0 \ \tau_1 \ \tau_2];
\]

4: produced the redundant dictionary within range of \([pare(i)*0.8 \ pare(i)*1.2] i=1,2,\ldots,6\);

5: initialize BA: q=6;

\[
P^q = \{ p_{1} \ldots p_{q} \} = pare + \lambda \times (1.2^* \ pare_0 - 0.8^* \ pare_1)
\]

6: Found the best matching atom \( g_{n, best} \) by BA and calculated the optimal fitness value fitness_{n, best}.

7: If (fitness_{n, best} < fitness_{n-1, best})

8: update \( \xi_{n+1} = \xi_n - (g_{n, best} \ * \ g_{n, best}^*) \ g_{n, best}; \)

9: n=n+1;

10: else

11: update fitness_{n, best} = fitness_{n-1, best}

12: end

13: k=k+1;

14: end

Parameters calculated results were shown in Fig.6.

Figure 6. Target segment extraction based on correlation detection technique. The upper and lower subplot denoted the signal components matched with t-atom and g-atom, respectively; black solid lines were cross-correlation functions; red and blue scatters described the envelopes of the effective correlated windows; red and black dotted lines were calculated and actual time range, respectively. The black cycle indicated the start-end of the target segment.

The first generation of the optimal trajectory was shown in Fig. 7. The final iteration number was 26 with a time of 10.7 s. Wavelet decomposition (WD), VMD, SD based on PSO were carried out to verify the performance of BASD. After several repeated attempts, sym3 wavelet basis and 3 decomposition layers were selected for WD. Intrinsic mode was set to be 4 for VMD.
Figure 7. BA algorithm optimization path. Where the blue balls represented the fitness values, the larger the diameter indicated the larger the residuals, and the weaker the fitness. The red ball was the result of the optimization solution.

The extraction results obtained by different methods were shown in Fig.8. Corresponding error curves were shown in Fig. 9. In Fig.8, there were local feature aliasing in the results of VMD and WD and serious errors will occur if the parameters were preset unsuitably. The PSO trapped into a locally optimal solution for some factors, leading to local maximum errors in Fig. 8. While results obtained by BASD is relatively ideal.

Figure 8. Signal reconstruction results. Colors: black: the original noisy signal $x$; cyan: the target signal; green: signal extracted by wavelet; blue: signal extracted by VMD; jasper: signal extracted by PSO; red: signal extracted by BASD.

Figure 9. Results error of different methods. The blue, green, jasper and red curves represented the errors between the target signal and the extracted signals obtained by VMD, WT, PSO and BASD.

Conclusion

For nonstationary signal extraction, the dictionary training algorithm based on feature parameters is firstly used to determine the key parameter range of feature atoms, which can effectively reduce the redundancy while ensuring the completeness of the redundant dictionary; the bat algorithm is utilized to transform the signal sparse decomposition problem into an optimization problem with the $\ell_2$-norm of residual as the fitness, which effectively improves the efficiency of the sparse decomposition algorithm. The experimental results showed that compared with other methods, the BASP algorithm is flexible and auto-adaptive, and can extract nonstationary signals from noise background without over constrained prior knowledges, and ensures maximum retention of other signal components, which is very important for weak signal detection under strong sparse interference.

References


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Author contributions

S.C. Ge wrote the main manuscript text, and S. D. Zhou participated in the experiment. All authors reviewed the manuscript.

Competing interests

The authors declare no competing interests.

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