Artificial Intelligence Approaches for GPS GDOP Classification

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ABSTRACT

Geometrical dilution of precision (GDOP) concept is a powerful and widespread quantify for determining the errors resulting from satellite configuration geometry. GDOP computation is based on the complicated transformation and inversion of measurement matrices that has a time and power burden. Also, basic back propagation neural network (BPNN) is easy to fall into local minima. To overcome this problem, in this study we propose an approach based on neural network (NN) and evolutionary algorithms (EAs) for GPS GDOP classification. In this article we use a number of EAs such as genetic algorithm (GA), particle swarm optimization (PSO), new PSO (NPSO), and imperialist competitive algorithm (ICA) to train an NN. Simulation results illustrate that the proposed methods have superiority performance.

General Terms

Signal Processing, and Pattern Recognition, Artificial Intelligence.

Keywords

GPS GDOP classification, neural network (NN), genetic algorithm (GA), particle swarm optimization (PSO), new PSO (NPSO), and imperialist competitive algorithm (ICA).

1. INTRODUCTION

Global positioning system (GPS) is a satellite based positioning system funded by the U.S. Department of Defense (DOD) in 1973. There are at least 24 satellites orbiting the earth in 6 planes transmitting signals which are received by a GPS receiver [1,2]. In order to verify the accuracy of the position, the user equivalent range error, the error of a component in the distance from receiver to a satellite, is multiplied by a factor which depends on the geometry of satellites constellation. This factor is called the geometric dilution of precision (GDOP) [2]. There are basically two main approaches employed to justify the satellite geometry based on the GPS GDOP including approximation and classification. Unlike the GPS GDOP approximate methods which compute the values in order to select the optimal subsets of the satellites, the GPS GDOP classification methods are employed to choose one of the acceptable subsets (four optimum satellites from 24 existing satellites) of satellites for navigation uses [1-3].

A subset of satellites with the value of GPS GDOP less than 2 is ideal, i.e. the attained measure of location and time by these satellites are trustworthy, but a higher GPS GDOP indicates poorer satellites positioning and an inferior measurement configuration [4]. The GPS GDOP ratings are represented in Table 1 [5].

Table 1. GPS GDOP ratings

Class number	GPS GDOP value	Ratings	
Class 1	1	Ideal	
Class 2	2-3	Excellent	
Class 3	4-6	Good	
Class 4	7-8	Moderate	
Class 5	9-20	Fair	
Class 6	21-50	Poor	

To reduce the computational burden for classification and approximation of the GPS GDOP, Simon and El-Sherief have used the basic back propagation (BP) approach to train the neural network (NN) [6]. Although the BP is the most popular algorithm to train an NN, it has two important problems: 1) the BP training is very slow in many applications including the GPS GDOP classification; and 2) the BP easily falls in local minima. To overcome these problems, Jwo and Lai [1] suggested utilizing the BP with momentum to train the NN (BPNN), the optimal interpolative network, general regression NN (GRNN), and probabilistic NN (PNN). In [5] and [7] Azami et al. proposed to use some improved NN algorithms, namely, BP with adaptive learning rate and momentum, Fletcher-Reeves conjugate gradient algorithm (CGA), Polak-Ribikre CGA, Powell-Beale CGA, scaled CGA, resilient BP (RBP), Levenberg-Marquardt (LM), modified LM, one step secant (OSS) and quasi-Newton. In addition, to have uncorrelated and informative features of the GPS GDOP, principal component analysis (PCA) was used as a preprocessing step [8].

To overcome this problem, in this study we propose an approach based on neural network (NN) and evolutionary algorithms (EAs) for GPS GDOP classification. In this article we use a number of EAs including genetic algorithm (GA), particle swarm optimization (PSO), new PSO (NPSO), and imperialist competitive algorithm (ICA) to train an NN.

The rest of the paper is organized as follows: Section 2 presents GPS GDOP briefly. In Section 3, EAs including genetic algorithm (GA), particle swarm optimization (PSO), new PSO (NPSO), and imperialist competitive algorithm (ICA) are explained. Section 3 describes the proposed methods. Simulation results and comparing with existing well-known methods are discussed in Section 5. Finally, Section 6 includes the conclusion of the study.

2. THE CONCEPT OF GPS GDOP

Basically, the GPS accuracy and performance are dependent on the GDOP. Figure 1 illustrates the geometry of the satellites and its affect on the GDOP values [1].





Figure 1: Satellite's diagram and its relation with DOP: (a) Bad GDOP and (b) Good GDOP

We assume the definitions of GDOP computation, useful for the sequel of the document and help understanding our contributions. The absolute distance between a satellite and a user is defined as follows [8,9]:

$$R_{i} = \rho_{i} + \Delta \rho^{iono,i} + \Delta \rho^{trop,i}$$
(1) where:

$$\rho_{i} = \sqrt{(X_{i} - X_{u})^{2} + (Y_{i} - Y_{u})^{2} + (Z_{i} - Z_{u})^{2}} - \delta_{t_{u}}$$

 $\Delta \rho^{iono,i}$ and $\Delta \rho^{trop,i}$ which are the errors induced by the ionospheric and the tropospheric propagation, are calculated from a model, $(X_u, Y_u, Z_u, \delta t_u)$ are the four system unknowns and δt_u is the correction the receiver has to apply to its own clock. To resolve this system we need four equations which mean four provide ranges from four different

equations which mean four pseudo-ranges from four different satellites. The pseudo-ranges can be approximated by a Taylor expansion. We obtain:

$$\hat{\rho}_{i} = \sqrt{(X_{i} - X_{u})^{2} + (Y_{i} - Y_{u})^{2} + (Z_{i} - Z_{u})^{2}} -\delta t_{u}$$

$$(3)$$

The Taylor expansion at the first order is:

$$\Delta \rho_{i} = \rho_{i} - \rho_{i} = a_{xi} \Delta x_{u} + a_{yi} \Delta y_{u} + a_{zi} \Delta z_{u}$$
$$-c\Delta t_{u}$$
(4)

where:

$$a_{xi} = \frac{X_{i} - X_{u}}{r_{i}}; a_{yi} = \frac{Y_{i} - Y_{u}}{r_{i}}; a_{zi} = \frac{Z_{i} - Z_{u}}{r_{i}};$$

$$a_{xi} = \sqrt{(X_{i} - X_{u})^{2} + (Y_{i} - Y_{u})^{2} + (Z_{i} - Z_{u})^{2}};$$

$$a_{zi} = \sqrt{(X_{i} - X_{u})^{2} + (Y_{i} - Y_{u})^{2} + (Z_{i} - Z_{u})^{2}};$$

Let assume N_{sat} be the number of visible satellites. The matrix H is as follows:

$$H = \begin{bmatrix} a_{x1} & a_{y1} & a_{z1} & 1 \\ a_{x2} & a_{y2} & a_{z2} & 1 \\ \vdots & \vdots & \vdots & \vdots \\ a_{xN_{sat}} & a_{yN_{sat}} & a_{zN_{sat}} & 1 \end{bmatrix}$$
(6)

Let define the $G_{\text{matrix:}}$

$$G = (H^T H)^{-1} \tag{7}$$

The GDOP is:

$$GDOP = \sqrt{trace[G]} \tag{8}$$

3. EVOLUTIONARY ALGORITHMS (EAs)

In this section we explain four mentioned EAs briefly.

3.1 Particle Swarm Optimization and New Particle Swarm Optimization

The idea of PSO was first proposed by J. Kennedy and R. Eberhart in 1995 [10]. PSO is an evolutionary computing algorithm inspired by nature and is based on repetition. The

social behavioral of animals such as birds and fish when they are together has been the inspiration source for this algorithm [10]. PSO, the same as other evolutionary algorithms, starts with a random matrix as a primary population. Unlike genetic algorithms (GA), standard PSO doesn't have evolutionary operators like mutation and breeding. Each member of the population is named a particle. As a matter of fact, in the PSO algorithm a certain number of particles that are formed randomly make the primary values. There are two parameters for each particle including position and velocity of the particle, which are defined by a space vector and a velocity vector, respectively. These particles form a pattern in an ndimensional space and move to the desired value. The best position of each particle in the past and the best position among all particles are stored separately. According to the experience from the preceding moves, the particles make a decision about how to decide the next move. In each iteration, all particles in the *n*-dimensional

problem space move to an optimum point. In each iteration, the position and velocity of each particle can be modified according to the following equations:

$$v_{i}(t+1) = wv_{i}(t) + C_{1}r_{1}(p_{best_{i}}(t) - x_{i}(t)) + C_{2}r_{2}(g_{best_{i}}(t) - x_{i}(t))$$
(9)

$$x_{i}(t+1) = x_{i}(t) + v_{i}(t+1)$$
(10)

where *n* shows the dimension $(1 \le n \le N)$, C_1 and C_2 are positive constants, generally considered 2.0. r_1 and r_2 are random numbers uniformly between 0 and 1; *w* is inertia weight that can be constant [11].

Equation (9) states that the velocity vector of each particle is updated $\binom{v_i(t+1)}{}$ and the new and previous values of the vector position $\binom{x_i(t)}{}$ create the new position vector (

 x_i (t+1)). As a matter of fact, the updated velocity vector influences both local and global values. The best response of the local positions is the best solution of the particle until current execution time (p_{best}) and the best global solution is the best solution of the entire particles until current execution time (g_{best}). Flowchart of PSO is presented in Figure 2.



Figure 2. Flowchart of PSO

Because PSO stays in local minima of fitness function, we use NPSO. In each iteration, as it was mentioned in PSO, global best particle and local best particle are calculated. NPSO strategy uses the global best particle and local "worst" particle, the particle with the worst fitness value of the particle until current execution time [12]. It can be defined as:

$$v_{i}(t+1) = wv_{i}(t) + C_{1}r_{1}(p_{worst_{i}}(t) - x_{i}(t)) + C_{2}r_{2}(g_{best_{i}}(t) - x_{i}(t))$$
(11)

3.2. Genetic Algorithm

A GA has three major steps. In the first step we create an initial population of m randomly selected individuals. The first generation is created by the initial population. The second step inputs m individuals and gives as output an assessment for each of them based on an objective function known as fitness function. This evaluation explains proximity value of our demands respect to each one of these m individuals. The third step is responsible for the formulation of the subsequently generation [8,13].

The fittest individuals of the previous one create a new generation. This evaluation procedure of generation N and production of generation N+1 is iterated until a performance criterion is joined. The creation of offspring based on the fittest individuals of the previous generation is named as breeding. The breeding process has three basic genetic

operations, namely, reproduction, crossover and mutation [8,13].

In each iteration of GA, some of the individuals with highest fitness value are eliminated and replaced by new members. New members are produced by so called parents. Crossover selects probabilistically two of fittest individuals of generation N; then in a random way selects a number of their characteristics and exchanges them in a way that the selected characteristics of the first individual would be obtained by the second and vice versa. Following this procedure creates two new children that both belong to the new generation. Mutation is a physiologically-inspired disturbance to the system and is often used to avoid possible local minima. To model this for real-world optimization, initially, a random number for each bit of chromosome, then, if the random number is greater than a pre-defined "mutation threshold", that bit is flipped. Finally, the flowchart of GA is presented in Figure 3 [8,13].



Figure 3. The flowchart of GA-based optimization cycle

3.3. Imperialist Competitive Algorithm

ICA is a new population based optimization algorithm proposed in 2007 by Atashpaz-Gargari and Lucas [14]. Nowadays ICA has numerous applications like designing controller for industrial systems, solving optimization problems in PID controller, communication systems, and training and analysis of artificial NNs [15-18].

Similar to every evolutionary algorithm, this algorithm starts with the preliminary population with random numbers that

each of them is called a "Country". A number of the members of the population that have best fitness values are chosen as imperialists. Each member of the remaining population is named a colony. Total fitness values of an empire depends on both the power of the imperialist country and the power of its colonies [13,15].

In each step, the countries go toward their related imperialist. If the fitness value of a colony achieves more than its related fitness value of imperialist then, this colony and its related imperialist transform to imperialist and colony, respectively. In each step, the weakest colony of the weakest empire moves toward the closest empire, and the empire without any colony is eliminated. After a while, all empires fall down except for the most powerful one and all the colonies go under the control of this unique empire. Finally the flowchart of the ICA is illustrated in Figure 4 [13-16].



Figure 4. Flowchart of the ICA

4. THE GPS GDOP CLASSIFICATION METHODS

First, in order to reduce the instruction time, all of the input and output variants become between 0 and 1 by normalizing.

Since
$$H^{T}H$$
 is a 4×4 matrix, it has four λ_{i} ($i = 1, 2, 3, 4$)

eigenvalues. It is clear that the four eigenvalues for the matrix $(H^T H)^{-1}$ will be λ_i^{-1} .

Based on the fact that the trace of a matrix is the sum of its eigenvalues, Equation (8) would be as below [4,5]:

$$GDOP = \sqrt{\lambda_1^{-1} + \lambda_2^{-1} + \lambda_3^{-1} + \lambda_4^{-1}}$$
(12)

Mapping with the definition of four variants would be done as below:

$$x_{1}(\vec{\lambda}) = \lambda_{1} + \lambda_{2} + \lambda_{3} + \lambda_{4} = trace(H^{T}H)$$

$$x_{2}(\vec{\lambda}) = \lambda_{1}^{2} + \lambda_{2}^{2} + \lambda_{3}^{2} + \lambda_{4}^{2} = trace[(H^{T}H)^{2}]$$

$$x_{3}(\vec{\lambda}) = \lambda_{1}^{3} + \lambda_{2}^{3} + \lambda_{3}^{3} + \lambda_{4}^{3} = trace[(H^{T}H)^{3}]$$

$$x_{4}(\vec{\lambda}) = \lambda_{1}\lambda_{2}\lambda_{3}\lambda_{4} = \det(H^{T}H)$$
(16)

The GDOP can be considered as a mapping directly from \overrightarrow{X} to GDOP classes. Mapping from \overrightarrow{X} to the GDOP classes is

very non-linear and cannot be determined analytically, but it can be determined exactly by the NN. In this paper, NN is

planned to do the mapping from \overline{X} to the GPS GDOP classes. Figure 5 shows the whole classification block diagram of GPS GDOP using NN. As mentioned before in this paper we propose four EAs including ICA, GA, PSO, and NPSO.



Figure 5: Classification block diagram of GPS GDOP using NN

5. SIMULATIONS AND DISCUSSIONS

Six classes or outputs for the proposed NNs are considered for GPS GDOP classification. When the GDOP is higher than the first threshold, NN outputs are [1 0 0] (for the large state) and when the GDOP is placed between the first and second threshold, NN outputs are [0 1 0] (for the average case) and when the GDOP is smaller than the second threshold, NN outputs are [0 0 1] (for the low state). Table 2 presents the accuracy numbers of GPS GDOP classification. Table 5 shows comparing accuracy between these three methods for GPS GDOP. The simulation results demonstrate that GPS GDOP classification using NN and ICA has better accuracy than other NN- based approaches including GA, PSO, and NPSO.

Classification Methods	NN trained by ICA	NN trained by GA	NN trained by NPSO	NN trained by PSO	BPNN [1]
Classification ratess	95.15%	94.31%	94.84%	94.17%	93.16

Table 2: Comparison of classification rates for proposed methods

6. CONCLUSIONS

GPS GDOP shows the geometrical relationship between positioning accuracy and satellite configuration. It can be calculated by using the positions of four visible satellites and the position of the receiver. The ideal GPS GDOP value is equal or less than one. Values of two to three are considered acceptable, whereas values greater than nine are considered high and should be discarded. Existing approaches for GPS GDOP approximation and classification are time-consuming and unreliable. In this paper, we propose four EAs, namely PSO, NPSO, GA and ICA, to train an NN. The simulation results demonstrate that GPS GDOP classification using NN and ICA has better accuracy than simple classification using NN and other NN-based approaches based on the EAs.

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