# 3WC-GBNRS++:A novel three-way classifier with granular-ball neighborhood rough sets based on uncertainty (Supplementary materials)

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# S1 THE ORIGINAL GRANULAR-BALLS GENERATION METHOD

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Algorithm S1: The original granular-balls generation method [1]
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```
Input: Dataset \mathbb{D}, purity threshold p;
   Output: The set of granular balls GB list.
 1 Initialize GB list = \{\};
 2 Function Split_GB (GB):
 3
       if the purity of GB > p then
 4
           Add GB to the GB\_list;
 5
       end
 6
       else
 7
           GB_1, GB_2 \leftarrow GB is subdivided into two
             granular-balls by 2-means clustering algorithm;
 8
            Split\_GB(GB_1);
            Split_GB(GB_2);
10
       end
11 return
12 Function Main (\mathbb{D}):
       Turn \mathbb{D} to a granular ball GB;
13
14
       Split_GB(GB);
15
       return GB\_list;
16 return
```

In the process of generating granular-balls, Algorithm S1 introduces a purity threshold as a parameter to control the granularity. The purity threshold plays a crucial role in controlling the degree of granularity in the division of these granular-balls. The purity of a granular-ball can be determined by calculating the proportion of the majority label. If the purity of a granular-ball falls below the purity threshold, it will be divided again. Therefore, the higher the purity threshold, the higher the purity of the final generated granular-balls.

The process of threshold acquisition is shown in Algorithm S2. First, the quality index of each granular-ball in GBNRS++ is computed. Subsequently, the average fuzziness of  $\mathbb{X}_{\mathbb{R}}^{\mathbb{J}}$  is determined. Next, based on the objective function, we calculate  $H_{\psi(\mathbb{X}_{\mathbb{D}}^{\mathbb{J}})}$  and  $\left|\bar{H}_{\psi(\mathbb{X}_{\mathbb{D}}^{\mathbb{J}})} - \bar{H}_{\mathbb{X}_{\mathbb{D}}^{\mathbb{J}}}\right|$  under varying thresholds by ad-

```
Algorithm S2: The process of threshold acquisition
      Input: DBGB list, step size L
      Output: The decision thresholds \alpha^* and \beta^*
 1 for GB \in DBGB\_list do
             Compute the quality index m(GB);
 3 end
 4 Computing H_{\mathbb{X}_{\mathbb{R}}^{\mathbb{J}}}=\frac{1}{|U|}\sum_{GB\in DBGB\_list}\xi(GB), where \xi(GB)=4m(GB)(1-m(GB));
 5 Assuming \beta = 1 - \alpha and Min = H_{\mathbb{X}^{\mathbb{J}}};
 6 for \beta = 0 to 0.5 do
               \alpha = 1 - \beta;
             Computing H_{\psi(\mathbb{X}_{\mathbb{R}}^{\mathbb{J}})} and \left|\bar{H}_{\psi(\mathbb{X}_{\mathbb{R}}^{\mathbb{J}})} - \bar{H}_{\mathbb{X}_{\mathbb{R}}^{\mathbb{J}}}\right|; if \left|\bar{H}_{\psi(\mathbb{X}_{\mathbb{R}}^{\mathbb{J}})} - \bar{H}_{\mathbb{X}_{\mathbb{R}}^{\mathbb{J}}}\right| < Min then \left|\begin{array}{c} Min = \left|\bar{H}_{\psi(\mathbb{X}_{\mathbb{R}}^{\mathbb{J}})} - \bar{H}_{\mathbb{X}_{\mathbb{R}}^{\mathbb{J}}}\right|; \\ \beta^* = \beta; \\ \alpha^* = 1 - \beta^*; \end{array}\right|
10
11
12
13
               \beta = \beta + L;
15 end
16 return \alpha^* and \beta^*;
```

justing the step size. When the minimum value is achieved, the optimal thresholds are obtained. The main steps in Algorithm 2, from step 6 to step 14, involve searching for optimal thresholds  $(\beta^*, \alpha^*)$  by exploring all possible thresholds.

Algorithm S3 gives the details of choosing the turning point. To compute the set DIS, all GBs must be traversed, resulting in a time complexity of O(N). To compute dV and AllMaxDiff each takes O(N) time since it involves traversing the elements of DIS. The time complexity of this loop over i can be considered as O(N).

### REFERENCES

[1] M. Ester, H. Kriegel, J. Sander, and X. Xu, "A density-based algorithm for discovering clusters in large spatial databases with noise," in *Proc.* 2nd Int. Conf. Knowledge Discovery and Data Mining (KDD'96), 1996, p. 226–231.

## **Algorithm S3:** Construction of adaptive granular-ball neighborhood

```
Input: DBGB\_list, parameters Local R and
              Global R, predicted object x_{test}.
   Output: The \delta_R^*(x_{\text{test}}) of x_{\text{test}}.
1 Compute the set DIS = \{\gamma_1, \gamma_2, \gamma_3, \dots, \gamma_n\};
2 Len = length(DIS);
3 dV = \{\gamma_i - \gamma_{i+1}\}_{i=1}^{\text{Len}-1};
4 AllMaxDiff = \max(dV);
5 FitLen = 20;
6 for i = Len - FitLen; i \geq 2; i = i - 1 do
        CV = \{\gamma_j\}_{j=i+1}^{i+\mathrm{FitLen}};

Index = \{i+1,\ldots,i+\mathrm{FitLen}\};
        [a, b] = LinearFit(Index, CV);
        PredictV = a \times i + b;
10
        CDiff = \gamma_i – PredictV;
11
        dCV = \{CV_j - CV_{j+1}\}_{j=1}^{\text{FitLen}-1};
12
        CMAllMaxDiff = \max(dCV);
13
        if CDiff > LocalR × CMAllMaxDiff and
14
          CDiff > GlobalR \times AllMaxDiff then
15
             break:
        end
16
17 end
18 \delta_R^*(x_{\text{test}}) = \{ \gamma \text{ values less than } \gamma_i \} return \delta_R^*(x_{\text{test}});
```

### **S2 THE BIOGRAPHIES OF AUTHORS**



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