

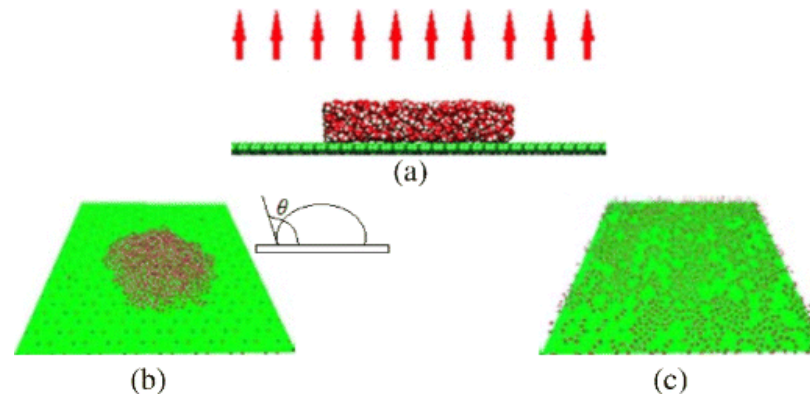
# Water Evaporation Optimization

**Soft Computing  
Final Report**

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# 演算法概念

- 仿照大自然中水分子的蒸散行為
- 隨著代次的演化，基底表面濕度逐漸下降
- Iteration < half of limit iteration -> **Monolayer evaporation phase**
- Iteration >= half of limit iteration -> **Droplet evaporation phase**
- 好的水分子留下，不好的蒸發



substrate with **low** wettability ( $q = 0$  e)

substrate with **high** wettability ( $q = 0.7$  e)

# 演算法步驟

- 取得起始解
- 根據使用者所設定之水分子數量，以及求解問題之上下屆，隨機亂數產生起始解，並記錄最好及最差值，以便下一步驟計算。

## 演算法步驟

- 計算表面能量( $E_{sub}$ ) / 接觸角度(contact angle)
- 上文有提到，濕度較大之表面影響水蒸發因子為表面本身所提供之能量，而此步驟就是將上一步之目標值，轉換成相同權重之表面能量：

$$E_{sub}(i)^t = \frac{(E_{max} - E_{min}) \times (Fit_i^t - \text{Min}(Fit))}{(\text{Max}(Fit) - \text{Min}(Fit))} + E_{min}$$

- ( $E_{max}$ 、 $E_{min}$ 為能量之上下界，此篇論文定為[-3.5,-0.5]， $Fit$ 為目標函式值)

## 演算法步驟

- 計算Monolayer evaporation Probability Matrix / Droplet evaporation Probability Matrix
- 蒸發量(Evaporation flux)，定義為水分子離開表面之大小，被認為一種更新演算法個體之計算量度，而此蒸發量根據上一步所計算之表面能量及接觸角改變

```
//1.2 Generate MEP matrix using Eq. (6)
double[,] mep = new double[numberOfWM, space];
for (int i = 0; i < mep.GetLength(0); i++)
{
    for (int j = 0; j < mep.GetLength(1); j++)
    {
        double rnd = RandomDouble(0, 1);
        if (rnd < Math.Exp(esub[i]))
        {
            mep[i, j] = 1;
        }
        else
        {
            mep[i, j] = 0;
        }
    }
}
```

更新

不更新

Exp(esub[i])  
Min:0.03  
Max:0.6

## 演算法步驟

- 產生新的水分子
- 根據Probability Matrix，以及亂數排列水分子，取得一組新位置新目標函式值之新一組水分子。

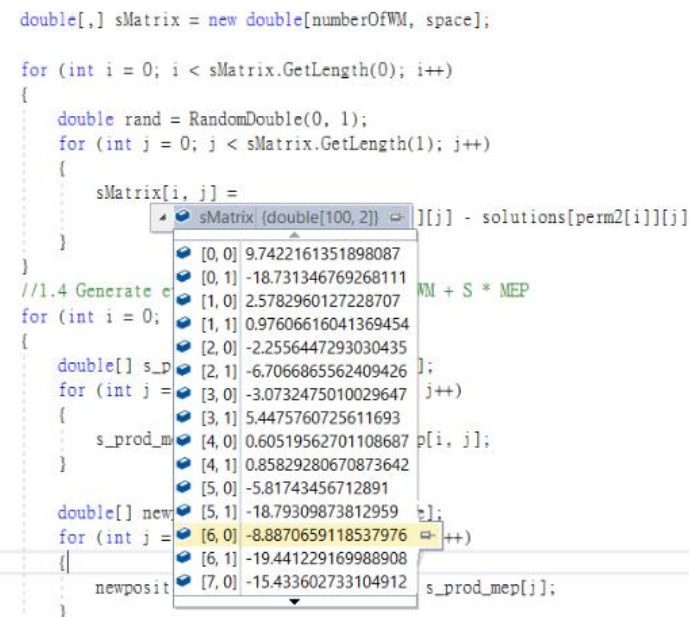
//1.3 Generate S matrix using Eq. (10)

```
int[] perm1 = GenerateRandomIntegerNumbers(numberOfWM);  
int[] perm2 = GenerateRandomIntegerNumbers(numberOfWM);
```

```
double[,] sMatrix = new double[numberOfWM, space];
```

```
for (int i = 0; i < sMatrix.GetLength(0); i++)  
{  
    double rand = RandomDouble(0, 1);  
    for (int j = 0; j < sMatrix.GetLength(1); j++)  
    {  
        sMatrix[i, j] =  
            rand * (solutions[perm1[i]][j] - solutions[perm2[i]][j]);  
    }  
}
```

```
double[,] sMatrix = new double[numberOfWM, space];  
  
for (int i = 0; i < sMatrix.GetLength(0); i++)  
{  
    double rand = RandomDouble(0, 1);  
    for (int j = 0; j < sMatrix.GetLength(1); j++)  
    {  
        sMatrix[i, j] =  
            sMatrix[i, j] - solutions[perm2[i]][j]  
    }  
}  
  
//1.4 Generate c  
for (int i = 0; i < sMatrix.GetLength(0); i++)  
{  
    double[] s_p  
    for (int j = 0; j < sMatrix.GetLength(1); j++)  
    {  
        s_prod_m  
    }  
    double[] new  
    for (int j = 0; j < sMatrix.GetLength(1); j++)  
    {  
        newposit  
    }  
}
```



## 演算法步驟

```
//1.4 Generate evaporated molecules:  $WM = WM + S * MEP$ 
for (int i = 0; i < numberOfWM; i++)
{
    double[] s_prod_mep = new double[space];
    for (int j = 0; j < s_prod_mep.Length; j++)
    {
        s_prod_mep[j] = sMatrix[i, j] * mep[i, j];
    }

    double[] newposition = new double[space];
    for (int j = 0; j < newposition.Length; j++)
    {
        newposition[j] = solutions[i][j] + s_prod_mep[j];
    }

    double newCost = objFun(newposition);
    newsolution[i] = newposition;
    newobjectiveValues[i] = newCost;
}
```

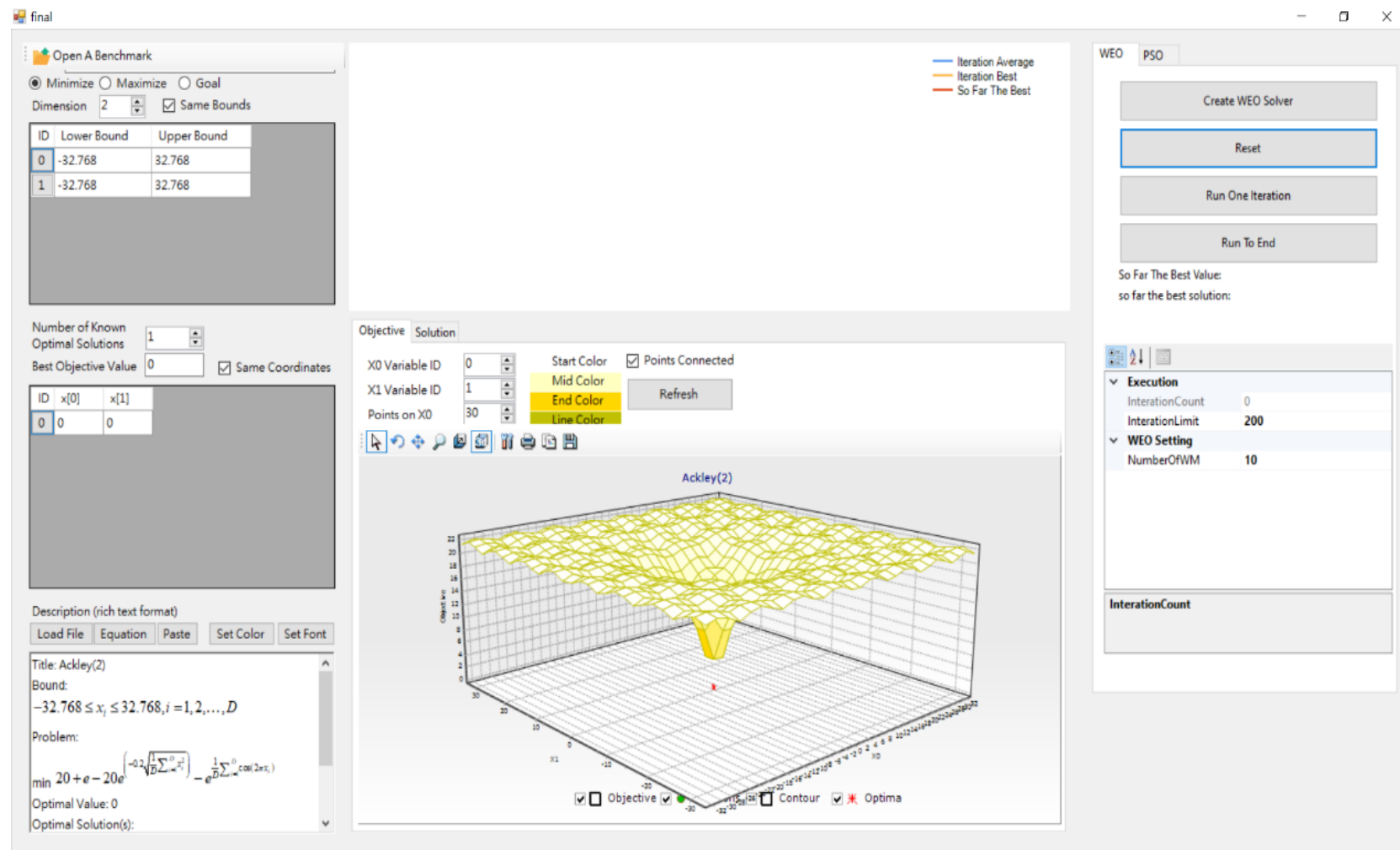
## 演算法步驟

- 更新水分子(即下一世代水分子)
- 比較新舊水分子，根據目標函式，如新水分子目標值比舊的值好，即取代，最後得的水分子，即為下一世代之水分子。

```
//3. Comparing and updateing water molecules
for (int i = 0; i < numberOfWM; i++)
{
    if (newobjectiveValues[i] < objectiveValues[i])
    {
        objectiveValues[i] = newobjectiveValues[i]; //t+1代的水分子
    }
}
```



# 介面介紹



## 實際結果

✚  $f_1(x) = \sum_{i=1}^n x_i^2$  ↵

✚ 1.Bound:[-100,100]    2.Limit iteration=100    3.dimension=2 ↵

➤ Number of water molecules=10 ↵

Objective value: ↵

0.000221 ↵	0.001348 ↵	0.000887 ↵	0.0034 ↵	0.003391 ↵
------------	------------	------------	----------	------------

➤ Number of water molecules=30 ↵

Objective value: ↵

0.002246 ↵	0.000744 ↵	<u>1.11E-05</u> ↵	0.000202 ↵	0.00015 ↵
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➤ Number of water molecules=100 ↵

Objective value: ↵

0.000221 ↵	<u>3.09E-05</u> ↵	0.000139 ↵	0.00092 ↵	0.000318 ↵
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## 實際結果

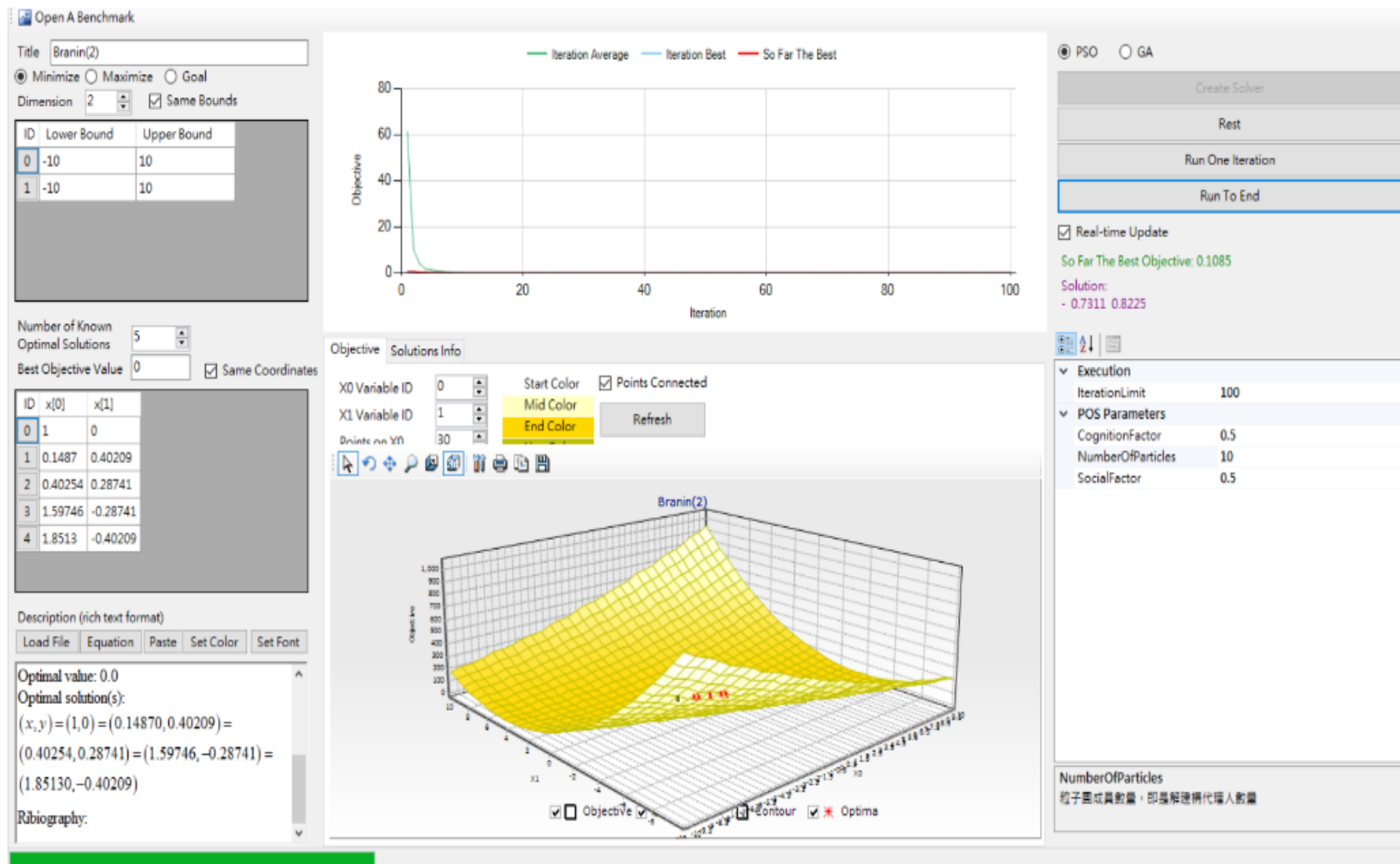
➤ 比較結果

	average	<u>svd</u>	best
<u>Nwm=10</u>	0.001237	0.00122	0.000221
<u>Nwm=30</u>	0.000659	0.000754	<u>1.11E-05</u>
<u>Nwm=100</u>	0.000238	0.000286	<u>3.09E-05</u>

水分子越多，解越好

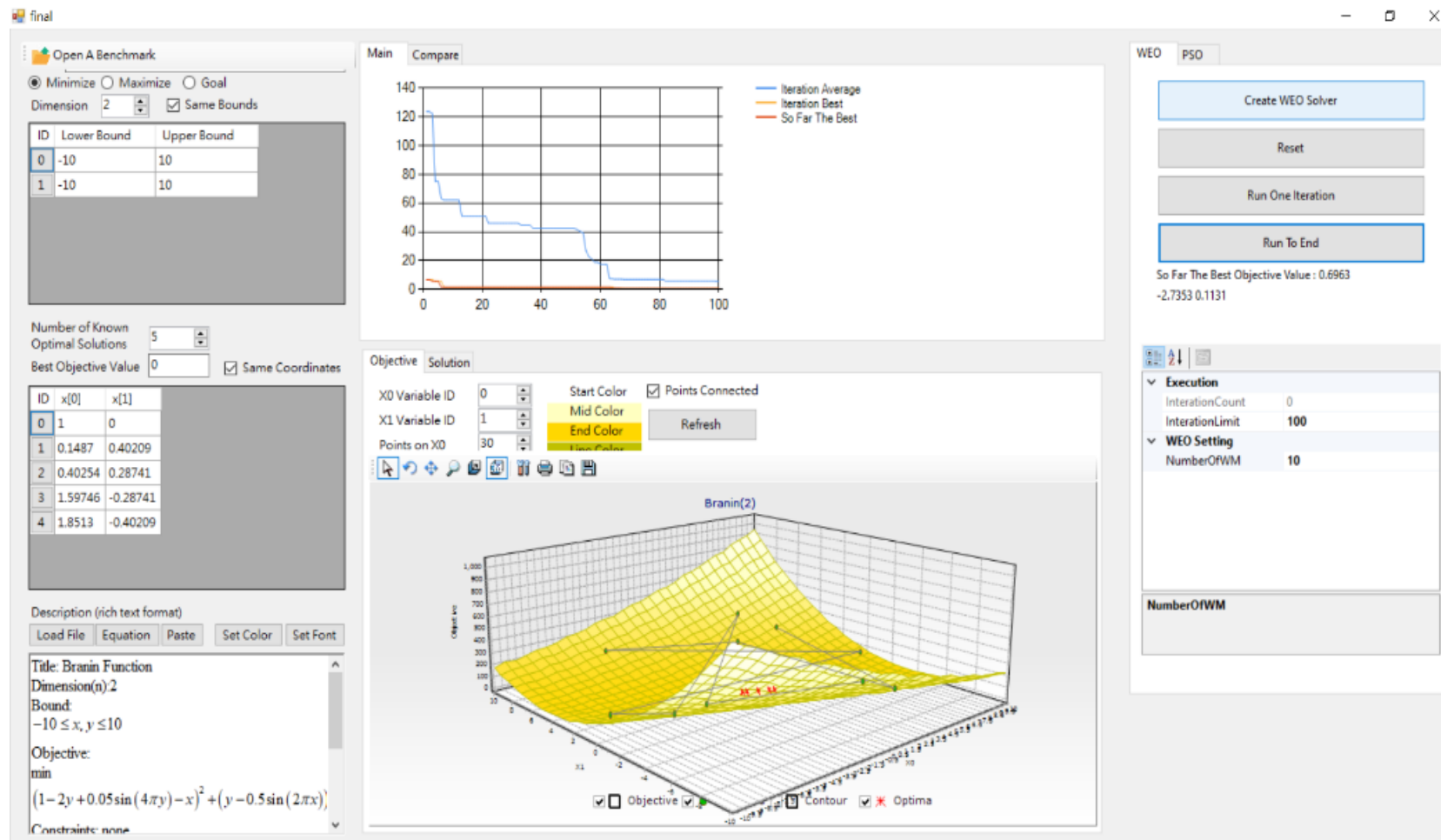
# 比較 psa

## 收斂快



# 比較 weo

## 收斂慢

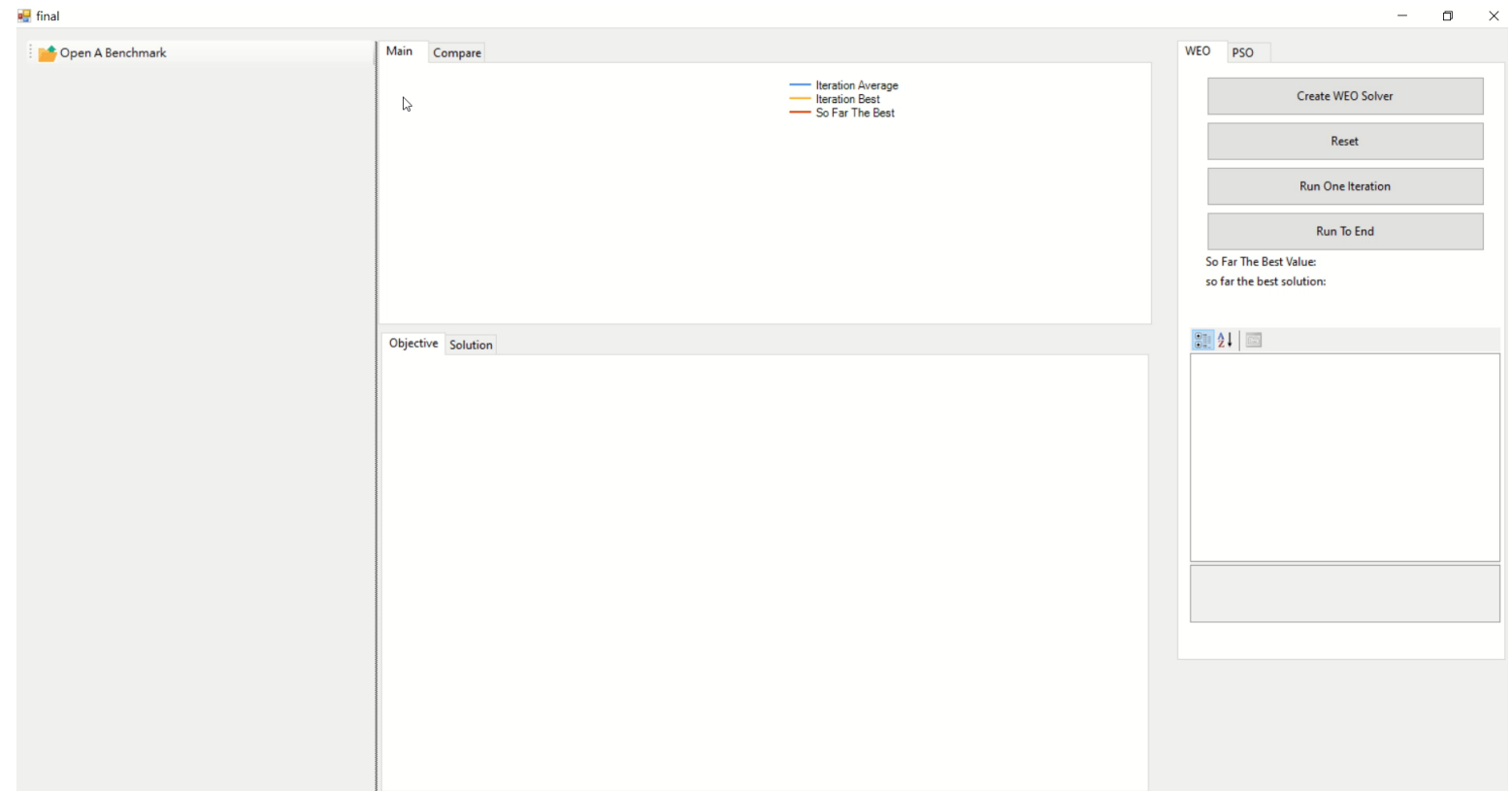


## 討論

- 收斂速度:PSO>WEO
- 物理性質之參數固定，讓使用者調整的參數少
- 低維問題效果比高維佳

	WEO	PSO
Branin Function(2)	0.0184	0.1085
Easom Function(2)	-0.9821	-0.5993
Ackley(2)	3.4875	1.8233
Ackley(30)	19.5180	16.2355
Girewank (30)	384.4991	69.3407

# Demo



THANKYOU!

