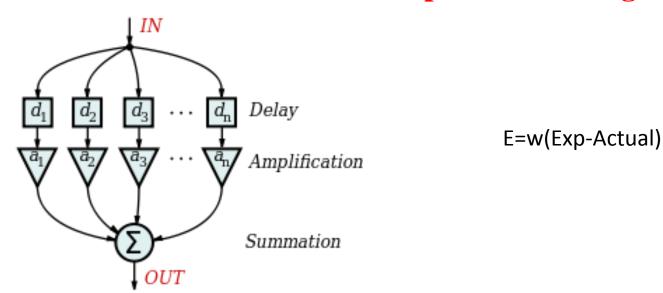
- Differential evolution (DE) was developed by Rainer Storn and Kenneth V. Price around 1995.
- DE is a unique evolutionary algorithm because it is **not** biologically motivate
- DE is used for multidimensional real-valued functions but **does not use the gradient** of the problem being optimized, which means DE **does not require** the optimization problem to be **differentiable**.
- DE can therefore also be used on optimization problems that are **not even continuous**, are noisy, change over time, etc.

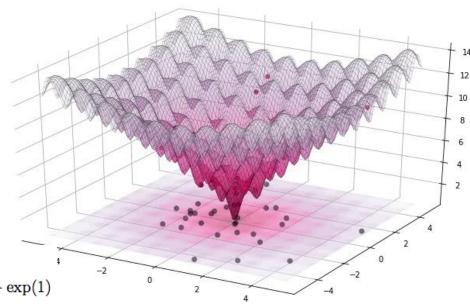
- Inspired from the real worlds problem of Digital filter coefficients
- In signal processing, a **digital filter** is a system that performs mathematical operations on a sampled, discrete-time signal to **reduce or enhance certain aspects of that signal**.



• DE optimizes a problem by maintaining a **population of candidate solutions** and **creating new candidate solutions** by **combining existing ones** according to its **simple formulae**, and then **keeping** whichever candidate solution has the **best score or fitness** on the optimization problem at hand.

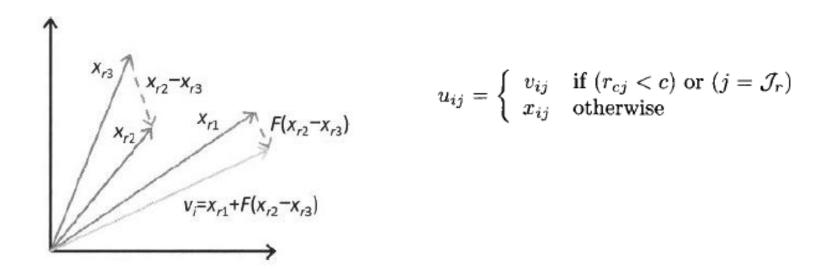
Global Minimum:

$$f(\mathbf{x}^*) = 0$$
, at $\mathbf{x}^* = (0, \dots, 0)$
 $x_i \in [-32.768, 32.768]$



$$f(\mathbf{x}) = -a \exp\left(-b\sqrt{\frac{1}{d}\sum_{i=1}^{d}x_i^2}\right) - \exp\left(\frac{1}{d}\sum_{i=1}^{d}\cos(cx_i)\right) + a + \exp(1)$$

A BASIC DIFFERENTIAL EVOLUTION ALGORITHM



• DE is based on the idea of **taking the difference vector** between two individuals, and **adding a scaled version** of the difference vector to a **third individual** to create a new candidate solution.

A BASIC DIFFERENTIAL EVOLUTION ALGORITHM

```
F = \text{stepsize parameter} \in [0.4, 0.9]
c = \text{crossover rate} \in [0.1, 1]
Initialize a population of candidate solutions \{x_i\} for i \in [1, N]
While not(termination criterion)
       For each individual x_i, i \in [1, N]
              r_1 \leftarrow \text{random integer} \in [1, N] : r_1 \neq i
              r_2 \leftarrow \text{random integer} \in [1, N] : r_2 \notin \{i, r_1\}
              r_3 \leftarrow \text{random integer} \in [1, N] : r_3 \notin \{i, r_1, r_2\}
              v_i \leftarrow x_{r1} + F(x_{r2} - x_{r3}) (mutant vector)
              \mathcal{J}_r \leftarrow \text{random integer} \in [1, n]
              For each dimension j \in [1, n]
                      r_{cj} \leftarrow \text{random number} \in [0, 1]
                      If (r_{cj} < c) or (j = \mathcal{J}_r) then
                             u_{ij} \leftarrow v_{ij}
                      else
                             u_{ij} \leftarrow x_{ii}
                      End if
              Next dimension
       Next individual
       For each population index i \in [1, N]
              If f(u_i) < f(x_i) then x_i \leftarrow u_i
       Next population index
Next generation
```

A BASIC DIFFERENTIAL EVOLUTION ALGORITHM

- This algorithm is often referred to as **classic DE**.
- It is also called **DE/rand/1/bin** because the base vector, x_{ri} , is **randomly chosen**; **one vector difference** (that is, $F(x_{r2} x_{r3})$) is added to x_{r1} and the number of mutant vector elements that are contributed to the trial vector closely follows a **binomial distribution**.
- In probability theory and statistics, the **binomial distribution** with parameters n and p is the discrete probability distribution of the number of successes in a sequence of n independent experiments, each asking a yesno question, and each with its own Boolean-valued outcome: yes or no (with probability q = 1 p).
- It would exactly follow a binomial distribution if not for the " $i=J_{\rm r}$ " test

Trial Vectors

- **DE/rand/1/L** works by generating a random integer L ϵ [l,n], copying L consecutive features from vi to ui, and then copying the remaining features from Xi to Ui

```
L \leftarrow \text{random integer} \in [1, n]
s \leftarrow \text{random integer} \in [1, n]
J \leftarrow \{s, \min(n, s + L - 1)\} \cup \{1, s + L - n - 1\}
For each dimension j \in [1, n]
If j \in J
u_{ij} \leftarrow v_{ij}
else
u_{ij} \leftarrow x_{ij}
End if
Next dimension
```

• For example, suppose that we have a seven-dimensional problem (n = 7). The DE/rand/1/L algorithm works by first generating a random integer L ϵ [l,n]; suppose that L = 3. We then generate a random starting point s ϵ [l,n]; suppose that s = 6.

```
u_{i1} \leftarrow v_{i1}
u_{i2} \leftarrow x_{i2}
u_{i3} \leftarrow x_{i3}
u_{i4} \leftarrow x_{i4}
u_{i5} \leftarrow x_{i5} \text{ (ending point)}
u_{i6} \leftarrow v_{i6} \text{ (starting point } s)
u_{i7} \leftarrow v_{i7}.
```

 $E(\text{number of } v_i \text{ elements copied}) = 1 + c(n-1)$ for DE/rand/1/bin.

 $E(\text{number of } v_i \text{ elements copied}) = n/2 \quad \text{for DE/rand/1/L}.$

Under what conditions is the expected number of mutant vector elements copied to the trial vector equal for the bin and L options?

 $E(\text{number of } v_i \text{ elements copied}) = 1 + c(n-1)$ for DE/rand/1/bin.

 $E(\text{number of } v_i \text{ elements copied}) = n/2 \quad \text{for DE/rand/1/L}.$

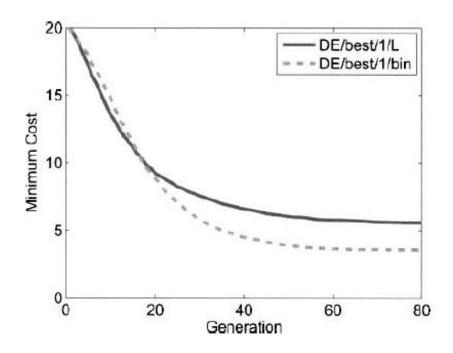
Under what conditions is the expected number of mutant vector elements copied to the trial vector equal for the bin and L options?

$$c = \frac{n-2}{2(n-1)}.$$

- Instead of randomly choosing the base vector x_{r1} , it may be beneficial to always use the best individual in the population as the base vector.
- That way the entire set of trial vectors U_i for $i \in [1, n]$ is comprised of mutations of the **best individual**.
- This approach is called **DE/best/1/bin.**

$$v_i \leftarrow x_b + F(x_{r2} - x_{r3})$$

• where x_b is the best individual in the population.



DE performance on the 20-dimensional Ackley function

 Another option is to use two difference vectors to create the mutant vector [Storn and sPrice, 1996]

$$r_{4} \leftarrow \text{random integer} \in [1, N] : r_{4} \notin \{i, r_{1}, r_{2}, r_{3}\}$$
 $r_{5} \leftarrow \text{random integer} \in [1, N] : r_{5} \notin \{i, r_{1}, r_{2}, r_{3}, r_{4}\}$
 $v_{i} \leftarrow \begin{cases} x_{r1} + F(x_{r2} - x_{r3} + x_{r4} - x_{r5}) & \text{DE/rand/2/?} \\ x_{b} + F(x_{r2} - x_{r3} + x_{r4} - x_{r5}) & \text{DE/best/2/?} \end{cases}$

- DE/rand/2/bin or DE/best/2/bin
- DE/rand/2/L or DE/best/2/L

• DE can also be implemented by using the current X_i as the base vector

$$v_i \leftarrow x_i + F\Delta x$$

where Δx is a difference vector.

- DE/target/1/bin, DE/target/2/bin,
- DE/target/1/L, or DE/target/2/L

- Yet another option is to create the difference vector by **using** the best individual in the population, x_b.
- This tends to create mutant vectors that all **move toward** \mathbf{x}_b . The vector that is subtracted from \mathbf{x}_b could be a random individual or the base individual.

$$v_{i} \leftarrow x_{i} + F(x_{b} - x_{i})$$

 $v_{i} \leftarrow x_{r1} + F(x_{b} - x_{r3})$
 $v_{i} \leftarrow x_{b} + F(x_{r2} - x_{r3} + x_{b} - x_{r5})$
 $v_{i} \leftarrow x_{i} + F(x_{b} - x_{i} + x_{r2} - x_{r3})$

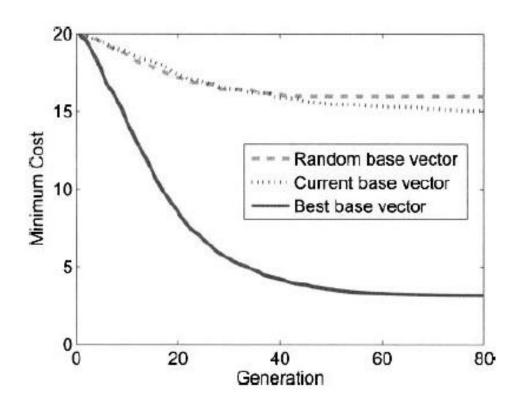
• If the last equation above is used to generate Vi, the algorithm is called **DE/target-to-best/1/bin**

- either-or algorithm
- We could **combine various methods** by randomly deciding how to generate the mutant vector.

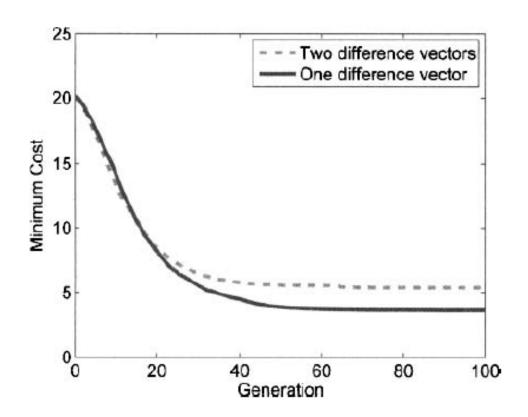
```
p_f = \text{mutation probability} \in [0, 1]
a \leftarrow \text{random number} \in [0, 1]
If a < p_f then
v_i \leftarrow x_{r1} + F(x_{r2} - x_{r3})
else
v_i \leftarrow x_{r1} + K(x_{r2} - x_{r1} + x_{r3} - x_{r1})
End if
```

Mutant vector generation that results in the DE/rand/1/either-or algorithm.

• K = (F + I)/2 gives good results in benchmark problems.



DE performance on the 20-dimensional Ackley function.



DE performance on the 20-dimensional Ackley function.

Scale Factor Adjustment

- DE's **scale factor F** determines the effect that difference vectors have on the mutant vector.
- So far we have assumed that F is a constant.
- We can vary the DE scale factor two different ways.
 - Dither
 - Jitter
- **Dither:** we can allow F to **remain a scalar** and randomly change it each time through the "for each individual"
- **Jitter**:-we can change **F** to an n-element vector and randomly change each element of F in the "for each individual" loop, so that each element of the mutant vector v is modified by a uniquely-scaled component of the difference vector.

Scale Factor Adjustment

Dither
$$F \leftarrow U[F_{\min}, F_{\max}]$$

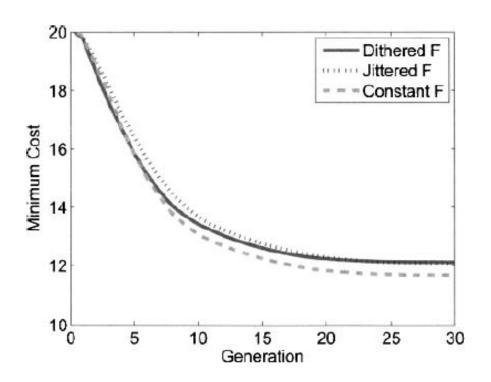
 $v_i \leftarrow x_{r1} + F(x_{r2} - x_{r3}).$

Jitter For each dimension
$$j \in [1, n]$$

$$F_j \leftarrow U[F_{\min}, F_{\max}]$$

$$v_{ij} \leftarrow x_{r1,j} + F_j(x_{r2,j} - x_{r3,j})$$
Next dimension

Scale Factor Adjustment



• DE performance on the 20-dimensional Ackley function with crossover rate c = 0.9. The traces show the cost of the best individual at each generation, averaged over 100 Monte Carlo simulations. The use of a constant scale factor F performs slightly better than dithering or jittering.

DISCRETE OPTIMIZATION

• The only place that discrete domains cause a problem in DE is in the generation of the **mutant vector**.

$$v_i \leftarrow x_{r1} + F(x_{r2} - x_{r3}).$$

Since F ∈ [0,1], V_i might not belong the problem domain D.

Mixed-Integer Differential Evolution

- One obvious approach to ensure that $V_i \in D$ is to simply project it onto D.
- For example, if *D* is the set of n-dimensional integer vectors, then

$$v_i \leftarrow \text{round}[x_{r1} + F(x_{r2} - x_{r3})]$$

round function operates element-by-element on a vector

A more general way to do this is

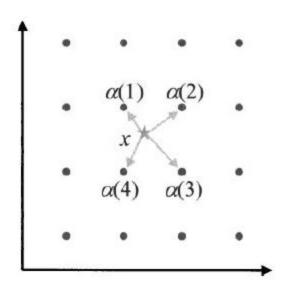
$$v_i \leftarrow P[x_{r1} + F(x_{r2} - x_{r3})]$$

where P is a projection operator such that $P(x) \in D$ for all x.

Mixed-Integer Differential Evolution

• P could be more complicated

$$P(x) = \arg\min_{\alpha} f(\alpha) : \alpha \in D, |x_j - \alpha_j| < 1 \text{ for all } j \in [1, n].$$



Projection of the continuous-valued vector *x onto a discrete-valued vector a.*

Discrete Differential Evolution

• Another way to modify DE for discrete problems is to **change the mutant vector generation method** so that it directly creates mutant vectors that lie in the discrete domain D.

$$v_i \leftarrow G(x_{r1}, x_{r2}, x_{r3})$$

$$v_i \leftarrow x_{r1} + \text{round}[F(x_{r2} - x_{r3})]$$

$$v_i \leftarrow x_{r1} + \operatorname{sign}(x_{r2} - x_{r3})$$

DIFFERENTIAL EVOLUTION AND GENETIC ALGORITHMS

```
Initialize a population of candidate solutions \{x_i\}, i \in [1, N]
While not(termination criterion)
      For each individual x_i, i \in [1, N]
             r_1 \leftarrow \text{random integer} \in [1, N] : r_1 \neq i
             v_i \leftarrow x_{r1}
             For each dimension j \in [1, n]
                    If rand(0,1) < c then
                           u_{ij} \leftarrow v_{ij}
                    else
                           u_{ij} \leftarrow x_{ij}
                    End if
             Next dimension
       Next individual
      For each i \in [1, N], If f(u_i) < f(x_i) then x_i \leftarrow u_i
Next generation
```

DIFFERENTIAL EVOLUTION AND GENETIC ALGORITHMS

```
Initialize a population of candidate solutions \{x_i\}, i \in [1, N]
While not(termination criterion)
       For each individual x_i, i \in [1, N]
              r_1 \leftarrow \text{random integer} \in [1, N] : r_1 \neq i
              r_2 \leftarrow \text{random integer} \in [1, N] : r_2 \notin \{i, r_1\}
               r_3 \leftarrow \text{random integer} \in [1, N] : r_3 \notin \{i, r_1, r_2\}
              v_i \leftarrow x_{r1} + F(x_{r2} - x_{r3})
               \mathcal{J}_r \leftarrow \text{random integer} \in [1, n]
               For each dimension j \in [1, n]
                      If (rand(0,1) < c) or (j = \mathcal{J}_r) then
                              u_{ij} \leftarrow v_{ij}
                      else
                             u_{ij} \leftarrow x_{ij}
                      End if
               Next dimension
       Next individual
       For each i \in [1, N], If f(u_i) < f(x_i) then x_i \leftarrow u_i
Next generation
```

Thank you