

Environmental Selection

```
function [Population,FrontNo,CrowdDis] = EnvironmentalSelection(Population,N)
% The environmental selection of NSGA-II

%% Non-dominated sorting
[FrontNo,MaxFNo] = NDSort(Population.objs,Population.cons,N);
Next = FrontNo < MaxFNo;

%% Calculate the crowding distance of each solution
CrowdDis = CrowdingDistance(Population.objs,FrontNo);

%% Select the solutions in the last front based on their crowding distances
Last      = find(FrontNo==MaxFNo);
[~,Rank] = sort(CrowdDis(Last),'descend');
Next(Last(Rank(1:N-sum(Next)))) = true;

%% Population for next generation
Population = Population(Next);
FrontNo    = FrontNo(Next);
CrowdDis    = CrowdDis(Next);
end
```

polyMutateCore

```
function result = polyMutateCore(genome, lb, ub, eta)
% Ported from the polynomial mutation in Pymoo, circa early 2022

% This function is written by Ian Meyer Kropp

delta1 = (genome - lb) / (ub - lb);    % Should be between 0 and 1
delta2 = (ub - genome) / (ub - lb);    % Should be between 0 and 1

exp = (eta + 1) ^ -1;

ran = rand(size(genome));
deltaq = zeros(size(genome));

leftMask = ran < 0.5;
rightMask = ran >= 0.5;

xy = 1 - delta1;
val = 2.0 * ran + (1.0 - 2.0 * ran) .* (xy .^ (eta + 1.0));
d = (val .^ exp) - 1.0;
deltaq(leftMask) = d(leftMask);

xy = 1.0 - delta2;
val = 2.0 * (1.0 - ran) + 2.0 * (ran - 0.5) .* (xy .^ (eta + 1.0));
d = 1.0 - (val .^ exp);
```

```

    deltaq(rightMask) = d(rightMask);

    muted_genome = genome + deltaq .* (ub - lb);

    muted_genome = min(max(muted_genome,lb),ub);

    result = muted_genome;
end

```

sbx

```

function Offspring = sbx(Parent, lb, ub, Parameter)
% Migrated from the PlatEMO OperatorGA module for convenience

% This function is written by Ian Meyer Kropp

    [proC,disC] = deal(Parameter{:});

    Parent1 = Parent(1:floor(end/2),:);
    Parent2 = Parent(floor(end/2)+1:floor(end/2)*2,:);
    [N,D]    = size(Parent1);

    beta = zeros(N,D);
    mu    = rand(N,D);
    beta(mu<=0.5) = (2*mu(mu<=0.5)).^(1/(disC+1));
    beta(mu>0.5)  = (2-2*mu(mu>0.5)).^(-1/(disC+1));
    beta = beta.*(-1).^randi([0,1],N,D);
    beta(rand(N,D)<0.5) = 1;
    beta(repmat(rand(N,1)>proC,1,D)) = 1;
    Offspring = [(Parent1+Parent2)/2+beta.*(Parent1-Parent2)/2
                 (Parent1+Parent2)/2-beta.*(Parent1-Parent2)/2];

    Lower = repmat(lb,2*N,1);
    Upper = repmat(ub,2*N,1);

    % Put everything back in bounds
    Offspring = min(max(Offspring,Lower),Upper);
end

```

sm2target

```

function newPop = sm2target(Pop, lb, ub, newSparsities)

% This function is written by Ian Meyer Kropp

% Sparse Mutate to a target
if numel(Pop) == 0
    newPop = Pop;
    return;

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end

[~,D] = size(Pop);

sparsities = sum(Pop == 0, 2) / D;

mutateMask = newSparsities ~= sparsities;

indv2mut = find(mutateMask);

% determine the non-zero increase/decrease for each indiv
nz2add = round(D * (sparsities(mutateMask) - newSparsities(mutateMask)));

% find where the non-zeros are
[zIndvs, zGenes] = find(Pop(indv2mut,:) == 0);
% find where the zeros are
[nzIndvs, nzGenes] = find(Pop(indv2mut,:) ~= 0);

newNzs = false(size(Pop));
newZs = false(size(Pop));

for i = 1:size(indv2mut,1)
    % gather relevant info
    indv_i = indv2mut(i);

    % Case where more non-zeros are needed
    if nz2add(i) > 0
        % find where there are zeros to flip
        zeroLocs = zGenes(zIndvs == i);

        % determine how many of them to flip
        numToFlip = nz2add(i);

        % Determine which of these possible zero positions to flip
        toFlip = zeroLocs(randperm(length(zeroLocs), numToFlip));

        % Record these positions
        newNzs(indv_i, toFlip) = true;

    % Case where more zeros are needed
    else

        % find where there are non-zeros to flip
        nZeroLocs = nzGenes(nzIndvs == i);

        % determine how many of them to flip
        numToFlip = -nz2add(i);

        % Determine which of these possible non-zero positions to flip
        toFlip = nZeroLocs(randperm(length(nZeroLocs), numToFlip));

        % Record these positions
        newZs(indv_i, toFlip) = true;
    end
end

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        end
    end

    %% Make the mutations
    % Find the min/max of the genome positions to mutate
    [~, newNzsCols] = find(newNzs);

    newNzsLb = lb(newNzsCols);
    newNzsUb = ub(newNzsCols);

    Pop(newNzs) = newNzsLb + rand(1,sum(newNzs, 'all')) .* (newNzsUb - newNzsLb);
    Pop(newZs) = zeros(sum(newZs, 'all'), 1);

    newPop = Pop;
end

```

SNSGAI

```

classdef SNSGAII < ALGORITHM
% <multi> <real> <large/none> <constrained/none> <sparse>
% Sparse nondominated sorting genetic algorithm II

% This function is written by Ian Meyer Kropp

    methods
        function main(Algorithm, Problem)

            [ sampling_method, mutation_method, crossover_method ] = ...
                Algorithm.ParameterSet( ...
                    {@vssps, 0.75, 1}, ...
                    @spm, ...
                    @ssbx ...
                );

            %% Generate random population

            sampler = sampling_method{1};
            lowerBound = sampling_method{2};
            upperBound = sampling_method{3};
            Population = sampler(Problem, lowerBound, upperBound);

            [~,FrontNo,CrowdDis] = EnvironmentalSelection(Population,Problem.N);

            %% Optimization
            while Algorithm.NotTerminated(Population)
                MatingPool = TournamentSelection(2,Problem.N,FrontNo,-CrowdDis);

                Offspring = sparseOperatorGA(Problem, Population(MatingPool), ...
                    {1,20,1,20,1,20,mutation_method,crossover_method});
            end
        end
    end
end

```

```

                [Population,FrontNo,CrowdDis] =
EnvironmentalSelection([Population,Offspring],Problem.N);
            end
        end
    end
end

```

sparseOperatorGA

```

function Offspring = sparseOperatorGA(Problem, Parent, Parameter)
% Adapted from OperatorGA in PlatEMO by Ian Meyer Kropp

% This function is written by Ian Meyer Kropp

%% Parameter setting
if nargin > 1
    [proC,disC,proM,disM,proSM,disSM,mutation_method,crossover_method] =
deal(Parameter{:});
else
    [proC,disC,proM,disM,proSM,disSM,mutation_method,crossover_method] =
deal(1,20,1,20,true,true);
end

calObj = false;

if isa(Parent(1),'SOLUTION')
    calObj = true;
    Parent = Parent.decs;
end

% Check if any of the decision variables are non-real values
if any(ones(size(Problem.encoding)) ~= Problem.encoding)
    error('Only real encoding supported.');
```

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end

Offspring = crossover_method(Parent, Problem.lower, Problem.upper,
{proC,disC});

mutation_params = {proM,disM, proSM,disSM};

Offspring = mutation_method(Offspring, Problem.lower, Problem.upper,
mutation_params);

if calObj
    Offspring = Problem.Evaluation(Offspring);
end
end

```

```

function newPop = spm(Pop, lb, ub, Parameter)

% This function is written by Ian Meyer Kropp

% Each row is a different population member
% Each column is a different genome

if nargin > 3
    [probMut,distrMut, probSMut, distrSMut] = deal(Parameter{:});
else
    [probMut,distrMut, probSMut, distrSMut] = deal(1,20,1,20);
end

[N,D] = size(Pop);

% Determine where the zeros are
nonZeroMask = Pop ~= 0;

%% Value mutations
ran = rand(size(Pop(nonZeroMask)));

toMutateNZ = ran < (probMut/D);

toMutate = false(size(Pop));

toMutate(nonZeroMask) = toMutateNZ;

[~, genomesToMutate] = find(toMutate);

lb_pm = lb(genomesToMutate)';
ub_pm = ub(genomesToMutate)';

% mutate values
Pop(toMutate) = polyMutateCore( Pop(toMutate), ...
                                lb_pm, ub_pm, distrMut);

%% Sparsity mutations

% Determine which population members to mutate sparsity
ran = rand(N, 1);

mutateMask = ran < probSMut/D;

% Figure out the individual sparsities of each individual
sparsities = sum(Pop == 0, 2) / D;

lb_sp = zeros(sum(mutateMask), 1);
ub_sp = ones(sum(mutateMask), 1);

newSparsities = sparsities;

newSparsities(mutateMask) = polyMutateCore(sparsities(mutateMask), lb_sp,

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ub_sp, distrSMut);

newSparsities = min(max(newSparsities,0),1);

% check if there's anything to do
if newSparsities == sparsities
    newPop = Pop;
else
    newPop = sm2target(Pop, lb, ub, newSparsities);
end
end

```

sssbx

```

function newPop = sssbx(Parent, lb, ub, Parameter)

% This function is written by Ian Meyer Kropp

%% Fetch paramters/setup
if nargin > 3
    [proC,disC] = deal(Parameter{:});
else
    [proC,disC] = deal(1,20);
end

Parent1 = Parent(1:floor(end/2),:);
Parent2 = Parent(floor(end/2)+1:floor(end/2)*2,:);

% figure out where are zeros/non-zeros
zMaskP1 = Parent1 == 0;
zMaskP2 = Parent2 == 0;

nzMaskP1 = ~zMaskP1;
nzMaskP2 = ~zMaskP2;

% figure out which positions are both zero or both non-zero
matching = (zMaskP1 & zMaskP2) | (nzMaskP1 & nzMaskP2);
not_matching = ~matching;

% empty template for results
Offspring1 = ones(size(Parent1))*-99;
Offspring2 = ones(size(Parent2))*-99;

%% Step 1: crossover on positions are both non-zero or both zero

[~,genes2sbx] = find(matching);

sbx_results = sbx([Parent1(matching)';Parent2(matching)'], lb(genes2sbx),
ub(genes2sbx), {proC, disC});

Offspring1(matching) = sbx_results(1,:);

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Offspring2(matching) = sbx_results(2,:);

%% Step 2: swap values that are mismatches between zero and non-zero

% empty mask of which positions to swap
swap_mask = ones(sum(not_matching, 'all'), 1);

% generate random number to determine how many zeros/non-zeros will go
% to each child
z2nzRatio = unifrnd(0,1);

swap_mask = sm2target(swap_mask', 0, 1, z2nzRatio);

swap_mask = swap_mask == 1;

% swap
not_matching_p1 = Parent1(not_matching);
not_matching_p2 = Parent2(not_matching);

not_matching_p1_temp = not_matching_p1;
not_matching_p1(swap_mask) = not_matching_p2(swap_mask);
not_matching_p2(swap_mask) = not_matching_p1_temp(swap_mask);

Offspring1(not_matching) = not_matching_p1;
Offspring2(not_matching) = not_matching_p2;

% return result
newPop = [Offspring1; Offspring2];
end

```

vssps

```

function Population = vssps(prob, sLower, sUpper)
% Randomly generate an initial population

% This function is written by Ian Meyer Kropp

% Nomenclature example
% N = 8
% D = 14
%
%           Cycle length of 14
%           |
% |-----|-----|
% 1 1 1 1                                     -
%           1 1 1 1                             |-- One full cycle
%                   1 1 1                         |
%                           1 1 1                     -
% |-----|-----|-----|----- Cycle count of 4
% |---|---|---|----- Cycle count of 4
% 1 1

```



```

%      1 1
%      1
%      1
% |----|----|
%      |
%      Cycle length of 6

%% Result set up
pop = prob.Initialization();
varCount = size(prob.lower,2);
mask = false(prob.N, varCount);

%% Determine the positioning of each stripe per individual
densityVector = 1 - linspace(sLower, sUpper, prob.N);

widthVector = round(densityVector.*prob.D);

% Put widths back into bound if rounding error occurred
lb = floor((1- sLower)*prob.D);
widthVector(widthVector > lb) = lb;

cumulativeWidths = cumsum(widthVector);

% if all sparsities are 100%, then skip processing, since everything
% will be zeros
if sum(widthVector == 0) == prob.N
    processedIndvs = prob.N;
else
    processedIndvs = 0;
end

cycle_count = 0;
cycles = zeros(prob.N, prob.D);
while processedIndvs < prob.N
    % Figure out how many stripes will fit in this cycle
    cycle_count = cycle_count + 1;
    spotsThatFitMask = cumulativeWidths <= prob.D & cumulativeWidths ~= 0;
    numThatFit = sum(spotsThatFitMask);
    largestFit = max(cumulativeWidths(spotsThatFitMask));

    cumulativeWidths = cumulativeWidths - largestFit;
    cumulativeWidths(cumulativeWidths<0) = 0;
    processedIndvs = processedIndvs + numThatFit;
    spotsThatFit = find(spotsThatFitMask);
    cycles(cycle_count,1:numThatFit) = spotsThatFit;

end

%% Create density mask

% Mask out non-zero values cycle-by-cycle
currentIndv = 1;
for c = 1:cycle_count
    cycle = cycles(c, cycles(c, :) ~= 0);

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widths = widthVector(cycle);

gapToFill = prob.D - sum(widths);
gapSize = ceil((prob.D - sum(widths))/numel(widths));
% For each individual in the cycle
position = 1;
for i = 1:numel(widths)

    width = widths(i);

    % Determine if a gap is needed
    gapWidth = 0;
    if gapToFill > 0
        gapWidth = gapSize;
        gapToFill = gapToFill - gapWidth;
    end

    % Determine the position of the stripe
    startPoint = position;

    if c == cycle_count
        endPoint = position+width-1;
    else
        endPoint = position+width-1+gapWidth;
    end

    % Prevent overflow from a gap calculation
    if endPoint > prob.D
        endPoint = prob.D;
    end

    % Mask out stripe
    mask(currentIndv, startPoint:endPoint) = true;

    % Go to the next individual

    position = position+width+gapWidth;
    %position = position+width;

    currentIndv = currentIndv + 1;

end

end

%% Mask off population according to stripe position
sparse_pop = pop.decs;
sparse_pop(~mask) = 0;

% Recalculate objective and constraints
popDec = prob.CalDec(sparse_pop);
popObj = prob.CalObj(sparse_pop);
popCon = prob.CalCon(sparse_pop);

```

```
Population = SOLUTION(popDec, popObj, popCon);  
end
```