

```

using LinearAlgebra

"""
    gradient_descent(f, ∇f, x0; α, tol, LIMIT)
Gradient descent algorithm for a function f, gradient ∇f, initial guess x0, learning rate α,
and tolerance tol. LIMIT is max iterations.
"""
function gradient_descent(f, ∇f, x0; α = 1.e-3, tol = 1.e-7, LIMIT = 5.e6)
    global counter = 0
    x = x0

    while norm(∇f(x)) > tol && counter < LIMIT
        x = x .- α .* ∇f(x)
        global counter += 1
    end

    println("Number of iterations: $counter.")
    if counter == LIMIT
        println("Max iterations reached.")
    end

    return x
end

"""
    newton_method(f, ∇f, ∇2f, x0; tol, LIMIT)
Newton method for a function f, gradient ∇f, Hessian ∇2f, initial guess x0, and tolerance
tol. LIMIT is max iterations.
"""
function newton_method(f, ∇f, ∇2f, x0; tol = 1.e-7, LIMIT = 500000)
    global counter = 0
    x = x0

    while norm(∇f(x)) > tol && counter < LIMIT
        x = x .- ∇2f(x) \ ∇f(x)
        global counter += 1
    end

    println("Number of iterations: $counter.")
    if counter == LIMIT
        println("Max iterations reached.")
    end

    return x
end

"""
    nonnegative_matrix_factorization(X, k; tol, LIMIT)

```

Performs nonnegative matrix factorization $X = W * H$ where X is $n \times p$, W is $n \times k$, and H is $k \times p$. The inner dimension k is a hyperparameter.

"""

```
function nonnegative_matrix_factorization(X, k; tol = 1.e-7, LIMIT = 1000)
```

```
    n, p = size(X)
```

```
    W = ones(n, k)
```

```
    H = ones(k, p)
```

```
    global counter = 0
```

```
    oldWH = X
```

```
    newWH = W * H
```

```
    while norm(newWH - oldWH) > tol && counter < LIMIT
```

```
        oldWH = W * H
```

```
        H_num = W' * X
```

```
        H_denom = W' * W * H
```

```
        H = H .* (H_num ./ H_denom)
```

```
        W_num = X * H'
```

```
        W_denom = W * H * H'
```

```
        W = W .* (W_num ./ W_denom)
```

```
        global counter +=1
```

```
        newWH = W * H
```

```
    end
```

```
    return W, H
```

```
end
```

"""

```
    adaboost(X, y)
```

Performs the AdaBoost algorithm on a data set with entries consisting of -1's and 1's.

The columns of X are weak classifiers, and the entries of y are the labels for the rows of X .

"""

```
function adaboost(X, y; tol = 1.e-5, LIMIT = 5)
```

```
    n, p = size(X)
```

```
    coeffs = zeros(p)
```

```
    f(R) = dot(coeffs, R)
```

```
    global counter = 1
```

```
     $\epsilon = 1$ 
```

```
    while counter < LIMIT && norm( $\epsilon$ ) > tol
```

```

current_min = Inf
current_min_index = 1
total_error = 0
polarity = 1

for c = 1:p
    class_error = 0
    total_error = 0

    for r = 1:n
        if X[r, c] != y[r]
            class_error += exp(-y[r] * f(X[r, :]))
        end

        total_error += exp(-y[r] * f(X[r, :]))
    end

    this_polarity = 1

    if class_error > total_error / 2
        class_error = total_error - class_error
        this_polarity = -1
    end

    if class_error < current_min
        current_min = class_error
        current_min_index = c
        polarity = this_polarity
    end
end

epsilon = current_min / total_error
alpha = 1/2 * log((1-epsilon)/epsilon + 1e-10)

coeffs[current_min_index] += polarity * alpha

global counter += 1
end

println("Number of iterations: $counter.")

return R -> dot(R, coeffs)
end

"""
    k_means(X, k; LIMIT)
Performs k-means clustering on the matrix X with k clusters.
"""
function k_means(X, k; LIMIT = 500)

```

```

n, p = size(X)
centroids = zeros(k, p)
used = []
nearest = zeros(n)

# generate initial centroids randomly
for i = 1:k
    rand_idx = rand(1:n)
    while rand_idx ∈ used
        rand_idx = rand(1:n)
    end

    centroids[i, :] = X[rand_idx, :]
    append!(used, [rand_idx])
end

# identify nearest centroids for each observation given initial centroids
for row = 1:n
    current_min = Inf
    current_min_idx = 1

    for ctrd in 1:k
        dist = norm(X[row, :] .- centroids[ctrd, :])
        if dist < current_min
            current_min = dist
            current_min_idx = ctrd
        end
    end

    nearest[row] = current_min_idx
end

global counter = 0
centroids_changed = true
while counter < LIMIT && centroids_changed
    centroids_changed = false

    # update centroids
    for ctrd in 1:k
        idxs = []
        for row = 1:n
            if nearest[row] == ctrd
                append!(idxs, [row])
            end
        end

        this_cluster = X[idxs, :]
        numrows, _ = size(this_cluster)
    end
end

```

```

        cluster_mean = zeros(p)
        for col in 1:p
            for row in 1:numrows
                cluster_mean[col] += 1/numrows * this_cluster[row, col]
            end
        end

        if centroids[ctrd, :] != cluster_mean[:]
            centroids_changed = true
        end

        centroids[ctrd, :] = cluster_mean[:]
    end

    #identify new nearest centroids for each observation
    for row = 1:n
        current_min = Inf
        current_min_idx = 1

        for ctrd in 1:k
            dist = norm(X[row, :] .- centroids[ctrd, :])
            if dist < current_min
                current_min = dist
                current_min_idx = ctrd
            end
        end

        nearest[row] = current_min_idx
    end

    counter += 1
end

#final nearest centroids
for row = 1:n
    current_min = Inf
    current_min_idx = 1

    for ctrd in 1:k
        dist = norm(X[row, :] .- centroids[ctrd, :])
        if dist < current_min
            current_min = dist
            current_min_idx = ctrd
        end
    end

    nearest[row] = current_min_idx
end

return centroids, nearest

```

end