

```

"""L-Shaped Algorithm

A method for solving two-stage stochastic linear program with recourse

Implemented for:

Math 6367 Optimization (Prof. Dr. Ronald H.W. Hoppe)
HW01
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Working with:
    python 3.6.6
    cvxpy 0.4.10
    numpy 1.15.1
"""

import cvxpy as cp
import numpy as np

class L_Shaped_Algorithm(object):
    """L-Shaped Algorithm for solving stochastic linear programs.

    The L-Shaped Algorithm is described in:

    Birge, J. R., & Louveaux, F. (2011). Introduction to stochastic
    programming. Springer Science & Business Media.

    Problems are in the form:

        minimize    c @ x + Q(x)
        subject to  A_eq @ x = b_eq
                   A_ineq @ x <= b_ineq
                   x >= 0

        where      Q(x) = E[Q(x, s(w))]
                   Q(x, s(w)) = min { q(w) @ y | W@y = h(w) - T(w)@x, y>=0}
        note: here we use s to denote the random variable

    Parameters
    -----
    c : array_like, required

    A_eq : array_like or matrix, or None

    b_eq : array_like, or None

    A_ineq : array_like or matrix, or None

    b_ineq : array_like, or None

    W : array_like or matrix, required

    h_func : function, required
        Should take two arguments (x, s) and return an array_like. See Notes.

    T_func : function, required
        Should take two arguments (x, s) and return an array_like. See Notes.

    q : list of array_likes, required
        Entries of q, realizations, and probabilities should have aligned
        indices. Each member of these lists should correspond to a particular
        realization of the random variables

    realizations : list of array_likes, required
        The value of the random variable(s) for each possible realization

```

probabilities : list of array_likes, required
 Probability for each random variable realization

max_iter : int, optional, default 100
 Maximum iterations before forced stopping

precision : float, optional, default 10e-6
 Precision to check for equality of floats. Also, the precision to round off the final solution. Full floating point precision is maintained during intermediate calculations.

verbose : bool, optional, default False
 Whether to print detailed information from intermediate steps. Basic information will always be printed.

debug : bool, optional, default False
 Whether to print advanced debug information

Attributes

solution : numpy ndarray
 Final optimal solution to the stochastic linear program

value : numpy float
 Final optimal objective value at the solution

print_precision : int
 Number of digits after the decimal point. Created from precision.

K_ : int
 Number of possible realizations of the random variable(s)

nu : int
 Iteration counter

r : int
 Counter for feasibility cuts

s : int
 Counter for optimality cuts

D_list : list of ndarrays
 List of matrices for feasibility cuts

d_list : list of ndarrays
 List of vectors for feasibility cuts

E_list : list of ndarrays
 List of matrices for optimality cuts

e_list : list of ndarrays
 List of vectors for optimality cuts

x_nu_list : list of ndarrays
 List containing the solution values obtained in step 1 from each iterate

theta_nu_list : list of ndarrays
 List containing the value of theta obtained in step 1 from each iterate

objective_value_list : list of ndarrays
 List of the objective values from step 1 in each iterate

Notes

In general, $h(w)$ and $T(w)$ may change depending on the value of the current

x iterate on the random variable value under consideration. Therefore, $h(w)$ and $T(w)$ are specified by functions which look like, for example:

```
def T_func(x, s):
    return np.array([[-1,0],
                     [0,-1],
                     [0,0],
                     [0,0],
                     [0,0],
                     [0,0]])

def h_func(x, s):
    return np.array([0, 0, -0.8*s[0], -0.8*s[1], s[0], s[1]])
```

Examples

 #The following example solves Example 2 from Page 188 of Birge & Louveaux.

```
import numpy as np
from l_shaped_algorithm.l_shaped_algorithm import L_Shaped_Algorithm
```

```
c = np.array([0])
```

```
A_ineq = [1]
b_ineq = [10]
```

```
W = np.array([1])
```

```
h = []
T = []
```

```
q = [[1],[1],[1]]
s = [1,2,4]
p = [1/3,1/3,1/3]
```

```
def T_func(x, s):
    if x <= s:
        return np.array([1])
    else:
        return np.array([-1])
```

```
def h_func(x, s):
    if x <= s:
        return np.array([s])
    else:
        return np.array([-s])
```

```
Solver = L_Shaped_Algorithm(c = c,
                             A_eq = None,
                             b_eq = None,
                             A_ineq = A_ineq,
                             b_ineq = b_ineq,
                             W = W,
                             h_func = h_func,
                             T_func = T_func,
                             q = q,
                             realizations = s,
                             probabilities = p,
                             max_iter = 100,
                             precision=10e-6,
                             verbose=False, debug=False)
```

```
x_opt = Solver.solve()
```

```
print (Solver.value)
print (Solver.solution)
```

```
"""
```

```
def __init__(self, c, A_eq, b_eq, A_ineq, b_ineq, W, h_func, T_func, q,
```

```

        realizations, probabilities,
        max_iter = 100, precision=10e-6,
        verbose=False, debug=False):

    self.c = c
    self.A_eq = A_eq
    self.b_eq = b_eq
    self.A_ineq = A_ineq
    self.b_ineq = b_ineq
    self.W = W
    self.h_func = h_func
    self.T_func = T_func
    self.q = q
    self.realizations = realizations
    self.p = probabilities

    self.max_iter = max_iter
    self.precision = precision

    self.debug = debug
    self.verbose = verbose

    np.set_printoptions(suppress=True)
    np.set_printoptions(precision=abs(int(np.log10(self.precision))))
    self.print_precision = abs(int(np.log10(self.precision)))

    # K is the number of possible realizations of the random variable(s)
    self.K_ = len(q)

    # Check that lengths match
    if self.K_ != len(self.p):
        raise ValueError("q and p should be same length")

    #Initialize counters and lists to store computed quantities

    self.nu = 0          # iteration counter
    self.r = 0          # counter for feasibility cuts
    self.s = 0          # counter for optimality cuts

    #Matrices and vectors which will form constraints pertaining to
    #feasibility cuts, ie:
    #  $D[i] @ x \geq d[i]$  where  $1 \leq i \leq r$ 

    self.D_list = []     # list of matrices for feasibility cuts
    self.d_list = []     # list of vectors for feasibility cuts

    #Matrices and vectors which will form constraints pertaining to
    #optimality cuts, ie:
    #  $E[i] @ x \geq e[i]$  where  $1 \leq i \leq s$ 

    self.E_list = []     # list of matrices for optimality cuts
    self.e_list = []     # list of vectors for optimality cuts

    #Lists to hold the values obtained in each iteration
    self.x_nu_list = []
    self.theta_nu_list = []
    self.objective_value_list = []

    self.value = None
    self.solution = None

def solve(self):
    """Solve the stochastic linear program as specified

    Returns

```

```

-----
solution : numpy ndarray
    The optimal solution, x, to the problem
"""

for _ in range(self.max_iter):
    self.nu += 1 # iterate step counter
    print()
    print( "=====")
    print(f"===== Iteration {self.nu} =====")
    print( "=====")

    _ = self._step_1()

    cut_made = self._step_2()

    if cut_made == 1:
        # A feasibility cut was made
        # Go back to step 1
        continue
    else:
        print("No feasibility cuts needed")

    cut_made = self._step_3()
    if cut_made == 0:
        # optimal solution found

        self.value = np.round(self.objective_value_list[-1],
                               self.print_precision)
        self.solution = np.round(self.x_nu_list[-1],
                                  self.print_precision)

        print()
        print("Optimal Solution Found")
        print()
        print("Objective Value = ", self.value)
        print("Optimal Solution = ", self.solution)
        return self.solution

# If no solution is found after max_iter steps, then return None
print()
print(f"Maximum iterations ({self.max_iter}) reached, and no ",
      "optimal solution found")
print("Try increasing max_iter or decreasing precision")
return None

def dot(self, a, b):
    """Return the dot product of two vectors
    Uses the numpy @ operator.
    If the expression involves a cvxpy variable which is actually a scalar,
    the @ operator doesn't work, so return the product instead.
    """
    try:
        return a @ b
    except ValueError:
        return a * b

def _step_1(self):
    """Solve the linear program with any constraints imposed by previous
    feasibility and optimality cuts.
    """

    print (f"----- Step 1 -----")

    n = len(self.c)

    x = cp.Variable(n)
    theta = cp.Variable(1)

```

```

if self.s == 0:
    # There are no optimality cuts, so set theta to -inf
    objective = cp.Minimize(self.dot(self.c, x))
else:
    objective = cp.Minimize(self.dot(self.c, x) + theta)

constraints = [x >= 0]

if self.A_eq is not None:
    # We must append the equality constraints on x
    constraints.append( self.dot(self.A_eq, x) == self.b_eq )
if self.A_ineq is not None:
    # We must append the inequality constraints on x
    constraints.append( self.dot(self.A_ineq, x) <= self.b_ineq )

for r in range(len(self.D_list)):
    # add constraints for each feasibility cut
    constraints.append( self.dot(self.D_list[r], x) >= self.d_list[r] )

for s in range(len(self.E_list)):
    # add constraints for each optimality cut
    constraints.append(
        self.dot(self.E_list[s], x) + theta >= self.e_list[s] )

prob = cp.Problem(objective, constraints)
result = prob.solve(verbose=self.debug)

if result is None and self.nu == 1:
    self.objective_value_list.append(0)
    self.x_nu_list.append(np.zeros(self.c.shape))
    self.theta_nu_list.append(-np.inf)
    return 1

# CVX sometimes makes the variables into funny size matrices, so we
# need to make them n-by-1 vectors
x_solution = np.array([x.value])
x_solution = x_solution.reshape(x_solution.size)

if self.s == 0:
    theta_solution = -np.inf
else:
    theta_solution = theta.value

print ("objective value = ", np.round(result,
                                         self.print_precision))
print ("x_nu          = ", np.round(x_solution,
                                         self.print_precision))
print ("theta_nu       = ", np.round(theta_solution,
                                         self.print_precision))

self.objective_value_list.append(result)
self.x_nu_list.append(x_solution)
self.theta_nu_list.append(theta_solution)

return 1

def _step_2(self):
    """Solve LPs for each possible realization of the random variables, and
    make feasibility cuts as appropriate.
    """

    print ()
    print (f"----- Step 2 -----")

    n = self.W.shape[0]
    if len(self.W.shape) > 1:

```

```

        m = self.W.shape[1]
    else:
        m = 1

    for k in range(self.K_):
        vp = cp.Variable(n)
        vm = cp.Variable(n)
        y = cp.Variable(m)

        objective = cp.Minimize(cp.sum_entries(vp) + cp.sum_entries(vm))

        # We use the user-specified driver functions to get the correct
        # matrix T and h for this particular realization of the random
        # variables
        T = self.T_func(self.x_nu_list[-1], self.realizations[k])
        h = self.h_func(self.x_nu_list[-1], self.realizations[k])

        constraints = [
            self.dot(self.W, y) + vp - vm == h - self.dot(T, self.x_nu_list[-1]),
            vp >= 0,
            vm >= 0,
            y >= 0]

        prob = cp.Problem(objective, constraints)
        result = prob.solve(verbose=self.debug)

        if np.abs(result) > self.precision:
            # Then we need to add a feasibility cut
            self.r += 1

            # Get the dual variables
            sigma = -1 * constraints[0].dual_value
            sigma = np.array(sigma).reshape(sigma.size)

            print (f"Feasibility cut identified for k={k},")
            if self.verbose:
                print (f"corresponding to realization {self.realizations[k]}")
                print ("objective      = ",
                        np.round(result, self.print_precision))
                print ("dual objective = ",
                        np.round((h - T @ self.x_nu_list[-1]) @ sigma,
                                self.print_precision+1))
                print ("dual variables = ",
                        sigma)

            D = sigma.T @ T
            d = sigma.T @ h
            print ("Dk = ", np.round(D, self.print_precision))
            print ("dk = ", np.round(d, self.print_precision))
            self.D_list.append(D)
            self.d_list.append(d)
            return 1 # cut was made

        # If we get through all realizations of the random variables, and no
        # infeasibilities were identified, then return 0
        return 0 # cut was not needed

def _step_3(self):
    """Solve LPs for each possible realization of the random variable, and
    make optimality cuts as appropriate.
    """

    #n = self.W.shape[0]
    if len(self.W.shape) > 1:
        m = self.W.shape[1]
    else:
        m = 1

```

```

print ()
print (f"----- Step 3 -----")

# Setup the variables E and e
E = np.zeros(len(self.x_nu_list[-1]))
e = 0

for k in range(self.K_):
    y = cp.Variable(m)

    # We use the user-specified driver functions to get the correct
    # matrix T and h for this particular realization of the random
    # variables
    T = self.T_func(self.x_nu_list[-1], self.realizations[k])
    h = self.h_func(self.x_nu_list[-1], self.realizations[k])

    # Define the objective function and constraints
    objective = cp.Minimize(self.dot(self.q[k], y[0:len(self.q[k])]))
    constraints = [
        self.dot(self.W, y) == h - self.dot(T, self.x_nu_list[-1]),
        y >= 0]

    prob = cp.Problem(objective, constraints)
    result = prob.solve(verbose=self.debug)

    # Get the dual variables
    pi = -1 * np.array(constraints[0].dual_value)
    pi = np.array(pi).reshape(pi.size)

    if self.verbose:
        print (f"Optimality cut variables for k={k}:")
        print ("objective      = ", result)
        print ("dual variables = ", pi)

    E += self.p[k] * pi.T @ T
    e += self.p[k] * pi.T @ h

w_nu = e - E @ self.x_nu_list[-1]

if np.abs(self.theta_nu_list[-1] - w_nu) <= self.precison:
    # The solution is optimal
    return 0 # no cut needed, solution is optimal

# Else append optimality cut
if self.verbose:
    print ("w_nu = ", w_nu)
    print ("theta_nu = ", self.theta_nu_list[-1])
print ("Optimality cut made")
print ("E = ", np.round(E, self.print_precision))
print ("e = ", np.round(e, self.print_precision))
self.s += 1
self.E_list.append(E)
self.e_list.append(e)
return 1 # a cut was made

```