pypgapack Documentation

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CONTENTS

CHAPTER

ONE

GETTING STARTED WITH PYPGAPACK

1.1 Background

`pypgapack` is a Python wrapper for the parallel genetic algorithm library pgapack, written in C by David Levine. The source and documentation for pgapack can be found at http://ftp.mcs.anl.gov/pub/pgapack/. The motivation for wrapping the code is ultimately to support a class project aiming to optimize loading patterns of nuclear reactor cores, which is a rather large and difficult combinatorial problem. Lots of researchers have applied genetic algorithms (and many other algorithms) to the problem, and the class project aims to provide a flexible test bench in Python to investigate various ideas. Wrapping pgapack is one step toward that goal. pgapAck was chose largely due to limited but positive past experience with it.

It should be pointed out that a similar effort to wrap pgapack in Python was made called `pgapy` (see http://pgapy.sourceforge.net/), but I actually couldn't get it to work, probably because I didn't know a thing about building Python modules before I started this (and my minimal C knowledge didn't help matters). Hence, I decided to "roll my own" using SWIG in combination with a C++ wrapper around pgapack instead of interfacing directly with pgapack as `pgapy` does.

The `PGA` class wraps almost all of pgapack's functionality, including allowing user functions for several operations (like initialization, crossover, etc.) for the `PGA.DATATYPE_BINARY`, `PGA.DATATYPE_REAL` and `PGA.DATATYPE_INTEGER` alleles. No such support is currently offered for other allele types, including user-specified types. The intended way to use `pypgapack` is to derive classes from `PGA`, with objective and other functions as members.

Parallel functionality is supported with the help of mpi4py.

Note: `pypgapack` is currently in beta mode, so there may be many things that look wrapped but are not. Testing is a future goal, but not a priority—I need a grade! Feedback is welcome at robertsj@mit.edu.

1.2 Building pypgapack

Included in ./pypgapack are the required source files and a simple script `build_pypgapack` which generates the Python module. To build, do the following:

1. Build PGAPack with the patches in ./patches. The major difference is a slight change to allow use with C++. The Makefile template also is set to produce shared and static libraries.

2. Modify the paths and variables in `build_pypgapack` below to suit your needs.
3. The source as distributed is set for serial. To use in parallel, do the following:
    - Uncomment `PARALLEL` in `build_pypgapack`
    - Set `CXX` to the appropriate compiler (e.g. mpic++) in `build_pypgapack`
    - Delete or move the dummy `mpi.h` included with PGAPack to avoid redefinitions. There's probably a better approach.
    - This assumes PGAPack was built in parallel; if not, do so. Refer to the PGAPack documentation. You need an MPI-enabled compiler.
    - Get mpi4py (e.g. easy_install mpi4py). You need an MPI-enabled compiler. Note, a few files from mpi4py are included in ./pypgapack/mpi4py. These *may* need to be updated.
4. Execute `build_pypgapack` and set PYTHONPATH accordingly.

1.3 Next Steps

The user is encouraged to read the pgapack documentation thoroughly before using pypgapack, as the shared API is *not* covered in this documentation (and neither are the many PGAPack defaults). It's helpful to go through their examples in C/C++ or Fortran if you know the languages.

Thereafter, see the collection of *Examples*, which include several of the original pgapack examples along with a few additional ones that demonstrate how to use user-defined functions for a variety of operations. Reference output is included, though don't expect to reproduce the numbers exactly for the small number of generations used, as they'll be sensitive to compilation, etc.

For a quick refresher, the basic gist of genetic algorithms is discussed briefly in *Methods*, which lists a few references that may be of use.

Documentation for the relatively small number of additional methods not explicitly in pgapack can be found in the *API Reference*.

CHAPTER

TWO

EXAMPLES

All examples are located in the `pypgapack/examples` and the reference output for all examples is in `pypgapack/examples/output`. Aside from small floating point differences, the values should be the same given the use of a fixed random number generator seed in all the examples. A utility script `run_examples.py` is included to test user output to the included reference cases. (Note, the above might actually be untrue, as compilation can and will change how a fixed pseudo-random number sequence is generated.)

Also, the maximum generation count is limited to 50 for all cases to produce short output. Experiment with that limit to see better solutions.

2.1 Basic Examples

The following are some simple examples that illustrate the basic PGAPack functionality.

2.1.1 Example 1: MAXBIT

`pypgapack` is pretty easy to use, and to demonstrate, we'll solve the maxbit problem, the first example in the PGAPack documentation.

```
"""
pypgapack/examples/example01.py  --  maxbit
"""
from pypgapack import PGA
import sys
class MyPGA(PGA) :
    """
    Derive our own class from PGA.
    """
    def maxbit(self, p, pop) :
        """
        Maximum when all alleles are 1's, and that maximum is n.
        """
        val = 0
        # Size of the problem
        n = self.GetStringLength()
```

```
        for i in range(0, n) :
            # Check whether ith allele in string p is 1
            if self.GetBinaryAllele(p, pop, i) :
                val = val + 1
        # Remember that fitness evaluations must return a float
        return float(val)

# (Command line arguments, 1's and 0's, string length, and maximize it)
opt = MyPGA(sys.argv, PGA.DATATYPE_BINARY, 100, PGA.MAXIMIZE)
opt.SetRandomSeed(1)        # Set random seed for verification.
opt.SetMaxGAIterValue(50)   # 50 generations (default 1000) for short output.
opt.SetUp()                 # Internal allocations, etc.
opt.Run(opt.maxbit)         # Set the objective.
opt.Destroy()               # Clean up PGAPack internals
```

Running it yields the following output:

```
***Constructing PGA***
Iter #     Field      Value
10         Best       6.900000e+01
Iter #     Field      Value
20         Best       7.200000e+01
Iter #     Field      Value
30         Best       7.600000e+01
Iter #     Field      Value
40         Best       7.700000e+01
Iter #     Field      Value
50         Best       8.000000e+01
The Best Evaluation: 8.000000e+01.
The Best String:
[ 01111101111011110111111001110111101111011111110111111110100111111 ]
[ 01111111011111011111111010111101011011 ]
***Destroying PGA context***
```

2.1.2 Example 2: MAXINT

This is a similar problem, but the alleles are integers ranging from -100 to 100. Note that when the integer ranges are set, a cast to `intc` is used. Python uses high precision datatypes, and there doesn't seem to be a safe implicit conversion between the Python integer type and the C integer type behind the scenes (in SWIG land). Casting explicitly circumvents the issue.

```
"""
pypgapack/examples/example02.py  --  maxint
"""
from pypgapack import PGA
import numpy as np
import sys
class MyPGA(PGA) :
    """
    Derive our own class from PGA.
    """
    def maxint(self, p, pop) :
```

```
        """
        The maximum integer sum problem.

        The alleles are integers, and we solve
            max f(x) = x_1 + x_2 + ... + x_N
        subject to
            |x_i| <= 100 .
        That maximum is f(x) = 100n obtained for x_i = 100 for all i.
        """
        c = self.GetIntegerChromosome(p, pop) # Get pth string as Numpy array
        val = np.sum(c)                       #   and sum it up.
        del c                                 # Delete "view" to internals.
        return float(val)                     # Always return a float.

n    = 10                # String length.
# (Command line arguments, integers, string length, and maximize it)
opt = MyPGA(sys.argv, PGA.DATATYPE_INTEGER, n, PGA.MAXIMIZE)
opt.SetRandomSeed(1)     # Set random seed for verification.
u_b =  100*np.ones(n)    # Define lower bound.
l_b = -100*np.ones(n)    # Define upper bound.
# Set the bounds.  Note, need to cast as C-combatible integers.
opt.SetIntegerInitRange(l_b.astype('intc'), u_b.astype('intc'))
opt.SetMaxGAIterValue(50)     # 50 generations for short output.
opt.SetUp()                   # Internal allocations, etc.
opt.Run(opt.maxint)           # Set the objective.
opt.Destroy()                 # Clean up PGAPack internals.
```

Running it yields the following output:

```
***Constructing PGA***
Iter #     Field      Value
10         Best       5.100000e+02
Iter #     Field      Value
20         Best       6.480000e+02
Iter #     Field      Value
30         Best       7.150000e+02
Iter #     Field      Value
40         Best       7.760000e+02
Iter #     Field      Value
50         Best       8.220000e+02
The Best Evaluation: 8.220000e+02.
The Best String:
#    0: [      27], [     100], [      86], [      98], [      97], [      99]
#    6: [      79], [      89], [      64], [      83]

***Destroying PGA context***
```

2.1.3 Example 3: MAXREAL

This is the same problem, but for real alleles. Here, note use of `SetMutationalType()` with the option of `PGA.MUTATION_RANGE`. This forces mutated allele values to remain within the initial range specified, useful for cases with constrained inputs. The default adds some (small) random amount, but over many

iterations, this can cause allele values to go significantly beyond the initial range.

```
"""
pypgapack/examples/example03.py  --  maxreal
"""
from pypgapack import PGA
import numpy as np
import sys

class MyPGA(PGA) :
    """
    Derive our own class from PGA.
    """
    def maxreal(self, p, pop) :
        """
        The maximum real sum problem.

        The alleles are doubles, and we solve
        .. math::
          \max f(x) &= \sum^N_{n=1} x_n \\
              s.t.  &= |x_i| \leq 100
        That maximum is  :math:`f_{\textrm{max}}(x) = 100n` obtained for
        :math:`x_i = 100, i = 1\ldots N`.
        """
        c = self.GetRealChromosome(p, pop) # Get pth string as Numpy array
        val = np.sum(c)                    #   and sum it up.
        del c                              # Delete "view" to internals.
        return val                         # Already a float.

n   = 10              # String length.
# (Command line arguments, doubles, string length, and maximize it)
opt = MyPGA(sys.argv, PGA.DATATYPE_REAL, n, PGA.MAXIMIZE)
opt.SetRandomSeed(1)  # Set random seed for verification.
u_b = 100*np.ones(n)  # Define lower bound.
l_b =-100*np.ones(n)  # Define upper bound.
# Set the bounds.  Default floats are handled without issue.
opt.SetRealInitRange(l_b, u_b)
# Force mutations to keep values in the initial range, a useful
#   feature for bound constraints.
opt.SetMutationType(PGA.MUTATION_RANGE)
opt.SetMaxGAIterValue(50) # 50 generations for short output.
opt.SetUp()               # Internal allocations, etc.
opt.Run(opt.maxreal)      # Set the objective.
opt.Destroy()             # Clean up PGAPack internals.
```

Running it yields the following output:

```
***Constructing PGA***
Iter #     Field      Value
10         Best       5.535524e+02
Iter #     Field      Value
20         Best       6.550377e+02
Iter #     Field      Value
30         Best       7.378405e+02
```

```
Iter #     Field      Value
40         Best       7.537097e+02
Iter #     Field      Value
50         Best       7.827324e+02
The Best Evaluation: 7.827324e+02.
The Best String:
#   0: [   66.95374], [   47.92554], [    91.9023], [   75.87186], [   96.31829]
#   5: [   85.24209], [   69.58556], [   90.59017], [   86.26947], [   72.07333]

***Destroying PGA context***
```

2.2 Examples of User Defined Operators

These examples explore one of the strengths of PGAPack, namely user-defined operators.

2.2.1 Example 4: User-defined String Initialization

We redo *Example 2: MAXINT* by initializing the strings with our own routine. Here, that's done by generating a permutation using Numpy.

```
"""
pypgapack/examples/example04.py  --  maxint with user initialization
"""
from pypgapack import PGA
import numpy as np
import sys
class MyPGA(PGA) :
    """
    Derive our own class from PGA.
    """
    def maxint(self, p, pop) :
        """
        The maximum integer sum problem.
        """
        c = self.GetIntegerChromosome(p, pop) # Get pth string as Numpy array
        val = np.sum(c)                       #   and sum it up.
        del c                                 # Delete "view" to internals.
        return float(val)                     # Always return a float.

    def init(self, p, pop) :
        """
        Random permutations using Numpy.
        """
        n    = self.GetStringLength()
        c    = self.GetIntegerChromosome(p, pop)
        perm = np.random.permutation(n)
        for i in range(0, n) :
            c[i] = perm[i]
        del c

```

```
n    = 10                # String length.
# (Command line arguments, integers, string length, and maximize it)
opt = MyPGA(sys.argv, PGA.DATATYPE_INTEGER, n, PGA.MAXIMIZE)
opt.SetRandomSeed(1)     # Set random seed for verification.
np.random.seed(1)        # Do the same with Numpy.
u_b =  100*np.ones(n)    # Define lower bound.
l_b = -100*np.ones(n)    # Define upper bound.
# Set the bounds.  Note, need to cast as C-compatible integers.
opt.SetIntegerInitRange(l_b.astype('intc'), u_b.astype('intc'))
opt.SetMaxGAIterValue(50)     # 50 generations for short output.
opt.SetUp()                   # Internal allocations, etc.
opt.Run(opt.maxint)           # Set the objective.
opt.Destroy()                 # Clean up PGAPack internals.
```

Running it yields the following output:

```
***Constructing PGA***
Iter #     Field      Value
10         Best       5.100000e+02
Iter #     Field      Value
20         Best       6.480000e+02
Iter #     Field      Value
30         Best       7.150000e+02
Iter #     Field      Value
40         Best       7.760000e+02
Iter #     Field      Value
50         Best       8.220000e+02
The Best Evaluation: 8.220000e+02.
The Best String:
#    0: [      27], [     100], [      86], [      98], [      97], [      99]
#    6: [      79], [      89], [      64], [      83]

***Destroying PGA context***
```

2.2.2 Example 5: User-defined Crossover Operator

This example solves "Oliver's 30-city Hamiltonian cycle Traveling Salesman Problem", as described in Poon and Carter, *Computer Ops Res.*, **22**, (1995). More importantly, it demonstrates use of a user-defined crossover operator, namely the "Tie-Breaking Crossover" (TBX1) of the same work.

The problem has 30 cities in a plane, and the goal is to minimize the distance traveled when visiting each city just once in a complete loop. The problem has 40 equivalent optima, each with a distance of 423.741 units. The coordinates of the cities are in the code, and are from Oliver's original paper by way of Steve Dower's site. See *Parallel Examples* for a parallel version that tries matching the cited results.

```
"""
pypgapack/examples/example05.py  --  traveling salesman
"""
from pypgapack import PGA
import numpy as np
import sys

```

```
8  class MyPGA(PGA) :
9      """
10     Derive our own class from PGA.
11     """
12     def tsm(self, p, pop) :
13         """
14         Oliver's 30 city traveling salesman problem.
15         """
16         c = self.GetIntegerChromosome(p, pop) # Get pth string as Numpy array
17         val = self.distance(c)
18         del c
19         return val
20
21     def distance(self, c) :
22         """
23         Compute the total distance for a set of cities.
24         """
25         # x and y coordinates by city
26         x = np.array([54.0,54.0,37.0,41.0, 2.0, 7.0,25.0,22.0,18.0, 4.0,\
27                       13.0,18.0,24.0,25.0,44.0,41.0,45.0,58.0,62.0,82.0,\
28                       91.0,83.0,71.0,64.0,68.0,83.0,87.0,74.0,71.0,58.0])
29         y = np.array([67.0,62.0,84.0,94.0,99.0,64.0,62.0,60.0,54.0,50.0,\
30                       40.0,40.0,42.0,38.0,35.0,26.0,21.0,35.0,32.0, 7.0,\
31                       38.0,46.0,44.0,60.0,58.0,69.0,76.0,78.0,71.0,69.0])
32         n   = self.GetStringLength()
33         val = 0.0
34         for i in range(0, n-1) :
35             val += np.sqrt( (x[c[i]]-x[c[i+1]])**2 + (y[c[i]]-y[c[i+1]])**2 )
36         val += np.sqrt( (x[c[0]]-x[c[n-1]])**2 + (y[c[0]]-y[c[n-1]])**2 )
37         assert(val > 423.70) # DEBUG.
38         return val
39
40     def tbx(self, p1, p2, pop1, c1, c2, pop2) :
41         """
42         Tie-breaking cross-over.  See Poon and Carter for details.
43         """
44         # Grab the city id's.
45         paren1 = self.GetIntegerChromosome(p1, pop1)
46         paren2 = self.GetIntegerChromosome(p2, pop1)
47         child1 = self.GetIntegerChromosome(c1, pop2)
48         child2 = self.GetIntegerChromosome(c2, pop2)
49         assert(np.sum(paren1)==435) # DEBUG
50         assert(np.sum(paren2)==435) # DEBUG
51
52         # Copy the parents to temporary vector for manipulation.
53         n = self.GetStringLength()
54         parent1 = np.zeros(n)
55         parent2 = np.zeros(n)
56         for i in range(0, n) :
57             parent1[i] = paren1[i]
58             parent2[i] = paren2[i]
59
60         # Code the parents using "position listing".
```

```
61          code1    = np.zeros(n)
62          code2    = np.zeros(n)
63          for i in range(0, n) :
64              code1[parent1[i]] = i + 1
65              code2[parent2[i]] = i + 1
66
67          # Randomly choose two cross-over points.
68          perm   = np.random.permutation(n)
69          point1 = np.min(perm[0:2])
70          point2 = np.max(perm[0:2])+1
71
72          # Exchange all alleles between the two points.  (It's unclear to me
73          #   whether these points should be inclusive or not; here, they are.)
74          temp = np.zeros(point2-point1)
75          for i in range(point1, point2) :
76              temp[i-point1] = parent1[i]
77              parent1[i]     = parent2[i]
78              parent2[i]     = temp[i-point1]
79
80          # Generate a cross-over map, a random ordering of the 0,1,...,n-1
81          crossovermap = np.random.permutation(n)
82
83          # Multiply each allele of the strung by n and add the map.
84          parent1 = parent1*n + crossovermap
85          parent2 = parent2*n + crossovermap
86
87          # Replace the lowest allele by 0, the next by 1, up to n-1.  Here,
88          #   we sort the parents first, and then for each element, find
89          #   where the increasing values are found in the original.  There
90          #   is probably a simpler set of functions built in somewhere.
91          sort1 = np.sort(parent1)
92          sort2 = np.sort(parent2)
93          for i in range(0, n) :
94              index = np.where(parent1 == sort1[i])
95              parent1[index[0][0]] = i
96              index = np.where(parent2 == sort2[i])
97              parent2[index[0][0]] = i
98
99          # Map the string back to elements.  These are the offspring.
100         tempchild1 = np.zeros(n)
101         tempchild2 = np.zeros(n)
102         for i in range(0, n) :
103             tempchild1[parent1[i]] = i
104             tempchild2[parent2[i]] = i
105         for i in range(0, n) :
106             child1[i] = tempchild1[i]
107             child2[i] = tempchild2[i]
108
109 # Number of cities.
110 n    = 30
111
112 opt = MyPGA(sys.argv, PGA.DATATYPE_INTEGER, n, PGA.MINIMIZE)
113
```

```
114 # One possible benchmark solution
115 reference = np.array([ 0, 2, 3, 4, 5, 6, 7, 8, 9,10, \
116                       11,12,13,14,15,16,17,18,19,20, \
117                       21,22,24,23,25,26,27,28,29, 1] )
118 print "Reference distance is: ", opt.distance(reference)
119
120 opt.SetRandomSeed(1)                    # Set seed for verification.
121 np.random.seed(1)                       # Do the same with Numpy.
122 opt.SetIntegerInitPermute(0, n-1)       # Start with random permutations.
123 opt.SetPopSize(400)                     # Large enough to see some success.
124 opt.SetMaxGAIterValue(100)              # Small number for output.
125 opt.SetCrossover(opt.tbx)               # Set a cross-over operation.
126 opt.SetMutation(PGA.MUTATION_PERMUTE)   # Mutate by permutation.
127 opt.SetUp()                             # Internal allocations, etc.
128 opt.Run(opt.tsm)                        # Set the objective and run.
129 opt.Destroy()                           # Clean up PGAPack internals.
```

Running it once yields the following output:

```
***Constructing PGA***
Reference distance is:  423.740563133
Iter #     Field      Value
10         Best       1.012977e+03
Iter #     Field      Value
20         Best       9.924890e+02
Iter #     Field      Value
30         Best       9.924890e+02
Iter #     Field      Value
40         Best       9.924890e+02
Iter #     Field      Value
50         Best       9.419172e+02
Iter #     Field      Value
60         Best       8.709637e+02
Iter #     Field      Value
70         Best       8.709637e+02
Iter #     Field      Value
80         Best       8.524659e+02
Iter #     Field      Value
90         Best       8.524659e+02
Iter #     Field      Value
100        Best       7.805903e+02
The Best Evaluation: 7.805903e+02.
The Best String:
#    0: [      29], [       3], [       0], [      12], [      13], [       9]
#    6: [       4], [       8], [      11], [      10], [       5], [       6]
#   12: [       7], [      17], [      14], [      18], [      16], [      15]
#   18: [      26], [      24], [      22], [      19], [      21], [      25]
#   24: [      27], [      23], [      20], [      28], [       2], [       1]

***Destroying PGA context***
```

2.2.3 Example 6: User-defined Mutation Operator

We redo *Example 2: MAXINT* using a custom mutation operator, largely following the PGAPack example.

```
"""
pypgapack/examples/example06.py  --  maxint with user mutation.
"""
from pypgapack import PGA
import numpy as np
import sys
class MyPGA(PGA) :
    """
    Derive our own class from PGA.
    """
    def maxint(self, p, pop) :
        """
        The maximum integer sum problem.
        """
        c = self.GetIntegerChromosome(p, pop) # Get pth string as Numpy array
        val = np.sum(c)                       #   and sum it up.
        del c                                 # Delete "view" to internals.
        return float(val)                     # Always return a float.

    def mutate(self, p, pop, pm) :
        """
        Mutate randomly within -n to n
        """
        n     = self.GetStringLength()
        c     = self.GetIntegerChromosome(p, pop)
        count = 0
        for i in range(0, n) :
            if self.RandomFlip(pm) :
                k = self.RandomInterval(1, 2*n)-n
                c[i] = k
                count += 1
        del c
        return count

n   = 10                # String length.
opt = MyPGA(sys.argv, PGA.DATATYPE_INTEGER, n, PGA.MAXIMIZE)
opt.SetRandomSeed(1)    # Set random seed for verification.
np.random.seed(1)       # Do the same with Numpy.
u_b =  100*np.ones(n)   # Define lower bound.
l_b = -100*np.ones(n)   # Define upper bound.
# Set the bounds.  Note, need to cast as C-combatible integers.
opt.SetIntegerInitRange(l_b.astype('intc'), u_b.astype('intc'))
opt.SetMaxGAIterValue(50)    # 50 generations for short output.
opt.SetMutation(opt.mutate)  # Set a custom mutation.
opt.SetUp()                  # Internal allocations, etc.
opt.Run(opt.maxint)          # Set the objective.
opt.Destroy()                # Clean up PGAPack internals.
```

Running it yields the following output:

```
***Constructing PGA***
Iter #     Field      Value
10         Best       5.270000e+02
Iter #     Field      Value
20         Best       7.150000e+02
Iter #     Field      Value
30         Best       7.280000e+02
Iter #     Field      Value
40         Best       8.030000e+02
Iter #     Field      Value
50         Best       8.220000e+02
The Best Evaluation: 8.360000e+02.
The Best String:
#    0: [      67], [      65], [      94], [      89], [      85], [      99]
#    6: [      80], [      98], [      87], [      72]

***Destroying PGA context***
```

2.2.4 Example 7: User-defined End of Generation Operator

This example is almost the same as *Example 1: MAXBIT* but it adds an end-of-generation operator. Here, we're using it to flip a random bit to 1. Of course, since we're maximizing the bit sum, this is "climbing the hill" to a better answer. This is a trivial example of such heuristics; in other situations, there are more complex, physically-motivated approaches. Another use of an end-of-generation operator would be for post-generation processing, such as plotting fitnesses, writing to file, etc.

```
"""
pypgapack/examples/example07.py  --  maxbit with end-of-generation hill climb
"""
from pypgapack import PGA
import sys
class MyPGA(PGA) :
    """
    Derive our own class from PGA.
    """
    def maxbit(self, p, pop) :
        """
        Maximum when all alleles are 1's, and that maximum is n.
        """
        val = 0
        # Size of the problem
        n = self.GetStringLength()
        for i in range(0, n) :
            # Check whether ith allele in string p is 1
            if self.GetBinaryAllele(p, pop, i) :
                val = val + 1
        # Remember that fitness evaluations must return a float
        return float(val)
    def climb(self):
        """
        Randomly set a bit to 1 in each string
```

```
        """
        popsize = self.GetPopSize()
        n = self.GetStringLength()
        for p in range(0, popsize) :
            i = self.RandomInterval(0, n - 1)
            self.SetBinaryAllele(p, PGA.NEWPOP, i, 1)

# (Command line arguments, 1's and 0's, string length, and maximize it)
opt = MyPGA(sys.argv, PGA.DATATYPE_BINARY, 100, PGA.MAXIMIZE)
opt.SetRandomSeed(1)        # Set random seed for verification.
opt.SetMaxGAIterValue(50)   # 50 generations (default 1000) for short output.
opt.SetEndOfGen(opt.climb) # Set a hill climbing heuristic
opt.SetUp()                 # Internal allocations, etc.
opt.Run(opt.maxbit)         # Set the objective.
opt.Destroy()               # Clean up PGAPack internals
```

Running it yields the following output:

```
***Constructing PGA***
Iter #     Field      Value
10         Best       7.000000e+01
Iter #     Field      Value
20         Best       7.600000e+01
Iter #     Field      Value
30         Best       8.300000e+01
Iter #     Field      Value
40         Best       8.800000e+01
Iter #     Field      Value
50         Best       9.100000e+01
The Best Evaluation: 9.300000e+01.
The Best String:
[ 11111111111111111110111111111111111110111110110111110111111110111 ]
[ 11111111111111111111111101111111111 ]
***Destroying PGA context***
```

Notice that for the same settings, the heuristic improved the best solution a little bit. Of course, flipping just one bit out of 100 shouldn't be expected to work miracles.

2.3 Parallel Examples

The following examples illustrate use of `pypgapack` in a parallel setting using the `mpi4py` package.

2.3.1 Example 8: Parallel MAXBIT

We adapt our favorite *Example 1: MAXBIT* using MPI. We up the string length and population to bring out timing differences.

```
"""
pypgapack/examples/example08.py  --  parallel maxbit
"""
```

```
from pypgapack import PGA
from mpi4py import MPI
import sys
class MyPGA(PGA) :
    """
    Derive our own class from PGA.
    """
    def maxbit(self, p, pop) :
        """
        Maximum when all alleles are 1's, and that maximum is n.
        """
        val = 0
        # Size of the problem
        n = self.GetStringLength()
        for i in range(0, n) :
            # Check whether ith allele in string p is 1
            if self.GetBinaryAllele(p, pop, i) :
                val = val + 1
        # Remember that fitness evaluations must return a float
        return float(val)

comm = MPI.COMM_WORLD           # Get the communicator.
rank = comm.Get_rank()          # Get my rank.
if rank == 0 :                  # Just to show it works, have node 0
    seed = 1                    #   set seed=1 and n=500 and have all
    n    = 500
else :                          #   other nodes set seed=n=0.  Then,
    seed = 0                    #   broadcast them to all nodes with
    n    = 0
seed = comm.bcast(seed, root=0) #   node 0 as the root process,
n    = comm.bcast(n,    root=0) #   and verify by printing.
print " node=", rank, " seed=", seed, " n=", n

if rank == 0 :
    t_start = MPI.Wtime()  # Start the clock.
# (Command line arguments, 1's and 0's, string length, and maximize it)
opt  = MyPGA(sys.argv, PGA.DATATYPE_BINARY, n, PGA.MAXIMIZE)
opt.SetPopSize(1000)
opt.SetRandomSeed(seed)         # Set random seed for verification.
opt.SetMaxGAIterValue(50)       # 50 generations for short output.
opt.SetUp()                     # Internal allocations, etc.
opt.Run(opt.maxbit)             # Set the objective.
opt.Destroy()                   # Clean up PGAPack internals
if rank == 0 :
    t_end = MPI.Wtime()
    print "Elapsed time = ", t_end-t_start, " seconds."
MPI.Finalize() # Should be called automatically, but good practice.
```

Running it using `mpirun -np 1 python example08.py` yields the following output:

```
 node= 0  seed= 1  n= 500
***Constructing PGA***
Iter #     Field      Value
```

```
10          Best       2.930000e+02
Iter #      Field      Value
20          Best       3.030000e+02
Iter #      Field      Value
30          Best       3.080000e+02
Iter #      Field      Value
40          Best       3.240000e+02
Iter #      Field      Value
50          Best       3.280000e+02
The Best Evaluation: 3.290000e+02.
The Best String:
[ 1001010110110110111110111111011001010111010111101100011011100011 ]
[ 1111111101101111111110100101100011101110011111110110110011110001 ]
[ 1010100111101111111111101110101100111111110101100110011111111001 ]
[ 0101011111111100011111111111011011101100111101111011101110010110000 ]
[ 1011110100011110010110010110011011111001010001001011000101101111 ]
[ 0111101101111100111101011100111111010110111111100101110010110111 ]
[ 1011000101110001111100011111110111110100101110000010010011100010 ]
[ 1011101100101110111001110000011111111111101101111111 ]
***Destroying PGA context***
Elapsed time =  2.06364393234  seconds.
```

Running it using `mpirun -np 2 python example08.py` yields the following output:

```
 node= 0  seed= 1  n= 500
***Constructing PGA***
 node= 1  seed= 1  n= 500
***Constructing PGA***
Iter #      Field      Value
10          Best       2.930000e+02
Iter #      Field      Value
20          Best       3.030000e+02
Iter #      Field      Value
30          Best       3.080000e+02
Iter #      Field      Value
40          Best       3.240000e+02
Iter #      Field      Value
50          Best       3.280000e+02
***Destroying PGA context***
The Best Evaluation: 3.290000e+02.
The Best String:
[ 1001010110110110111110111111011001010111010111101100011011100011 ]
[ 1111111101101111111110100101100011101110011111110110110011110001 ]
[ 1010100111101111111111101110101100111111110101100110011111111001 ]
[ 0101011111111100011111111111011011101100111101111011101110010110000 ]
[ 1011110100011110010110010110011011111001010001001011000101101111 ]
[ 0111101101111100111101011100111111010110111111100101110010110111 ]
[ 1011000101110001111100011111110111110100101110000010010011100010 ]
[ 1011101100101110111001110000011111111111101101111111 ]
***Destroying PGA context***
Elapsed time =  1.14050102234  seconds.
```

This was using a dual core laptop with probably more browser windows open than needed, but the results

look good. Overall, PGAPack focuses parallelism on the object function evaluation. Hence, if your objective function is expensive to evaluate, you can expect relatively good scaling up to the number of strings replaced every generation (the default is 1/10 of the total).

2.3.2 Example 9: Parallel Traveling Salesman

We adapt the Traveling Salesman Problem to parallel and try matching the results Poon and Carter found for the TBX1 cross-over operator. The GA parameters used by Poon and Carter are not entirely clear. They cite results for a population of 21 over 300 iterations with a cross-over to mutation probability ratio of 0.8/0.2. The exact nature of the "swap" mutation is cited from another work I don't have at hand, and the selection appears to be standard elitist, i.e all but the best is replaced. I set a swap operator that always swaps a user-set number of pairs. Using 3 pairs (i.e. 10%) seems to get close to the cited results. Also, because PGAPack doesn't like odd population sizes, I use 22.

```
"""
pypgapack/examples/example09.py  --  traveling salesman in parallel
"""
from pypgapack import PGA
from mpi4py import MPI
import numpy as np
import sys

class MyPGA(PGA) :
    """
    Derive our own class from PGA.
    """
    def tsm(self, p, pop) :
        """
        Oliver's 30 city traveling salesman problem.
        """
        c = self.GetIntegerChromosome(p, pop) # Get pth string as Numpy array
        val = self.distance(c)
        del c
        return val

    def distance(self, c) :
        """
        Compute the total distance for a set of cities.
        """
        # x and y coordinates by city
        x = np.array([54.0,54.0,37.0,41.0, 2.0, 7.0,25.0,22.0,18.0, 4.0,\
                      13.0,18.0,24.0,25.0,44.0,41.0,45.0,58.0,62.0,82.0,\
                      91.0,83.0,71.0,64.0,68.0,83.0,87.0,74.0,71.0,58.0])
        y = np.array([67.0,62.0,84.0,94.0,99.0,64.0,62.0,60.0,54.0,50.0,\
                      40.0,40.0,42.0,38.0,35.0,26.0,21.0,35.0,32.0, 7.0,\
                      38.0,46.0,44.0,60.0,58.0,69.0,76.0,78.0,71.0,69.0])
        n   = self.GetStringLength()
        val = 0.0
        for i in range(0, n-1) :
            val += np.sqrt( (x[c[i]]-x[c[i+1]])**2 + (y[c[i]]-y[c[i+1]])**2 )
        val += np.sqrt( (x[c[0]]-x[c[n-1]])**2 + (y[c[0]]-y[c[n-1]])**2 )
        assert(val > 423.70) # Debug.
```

```
        return val

    def tbx(self, p1, p2, pop1, c1, c2, pop2) :
        """
        Tie-breaking cross-over.  See Poon and Carter for details.
        """
        # Grab the city id's.
        paren1 = self.GetIntegerChromosome(p1, pop1)
        paren2 = self.GetIntegerChromosome(p2, pop1)
        child1 = self.GetIntegerChromosome(c1, pop2)
        child2 = self.GetIntegerChromosome(c2, pop2)
        assert(np.sum(paren1)==435) # DEBUG
        assert(np.sum(paren2)==435) # DEBUG

        # String length.
        n = self.GetStringLength()
        parent1 = np.zeros(n)
        parent2 = np.zeros(n)

        for i in range(0, n) :
            parent1[i] = paren1[i]
            parent2[i] = paren2[i]

        # Code the parents using "position listing".
        code1   = np.zeros(n)
        code2   = np.zeros(n)
        for i in range(0, n) :
            code1[parent1[i]] = i + 1
            code2[parent2[i]] = i + 1

        # Randomly choose two cross-over points.
        perm   = np.random.permutation(n)
        point1 = np.min(perm[0:2])
        point2 = np.max(perm[0:2])+1

        # Exchange all alleles between the two points.
        temp = np.zeros(point2-point1)
        for i in range(point1, point2) :
            temp[i-point1] = parent1[i]
            parent1[i]     = parent2[i]
            parent2[i]     = temp[i-point1]

        # Generate a cross-over map, a random ordering of the 0,1,...,n-1
        crossovermap = np.random.permutation(n)

        # Multiply each allele of the strung by n and add the map.
        parent1 = parent1*n + crossovermap
        parent2 = parent2*n + crossovermap

        # Replace the lowest allele by 0, the next by 1, up to n-1.
        sort1 = np.sort(parent1)
        sort2 = np.sort(parent2)
        for i in range(0, n) :
```

```
92              index = np.where(parent1 == sort1[i])
93              parent1[index[0][0]] = i
94              index = np.where(parent2 == sort2[i])
95              parent2[index[0][0]] = i
96
97          tmpc1 = np.zeros(n)
98          tmpc2 = np.zeros(n)
99          # Map the string back to elements.  These are the offspring.
100         for i in range(0, n) :
101             tmpc1[parent1[i]] = i
102             tmpc2[parent2[i]] = i
103         for i in range(0, n) :
104             child1[i] = tmpc1[i]
105             child2[i] = tmpc2[i]
106
107     def swap(self, p, pop, pm) :
108         """
109         Random swap of allele pairs.  Note, nswap must be set!
110         """
111         n     = self.GetStringLength()
112         c     = self.GetIntegerChromosome(p, pop)
113         index = np.random.permutation(n)
114         for i in range(0, self.nswap) :
115             i1 = index[2*i  ]
116             i2 = index[2*i+1]
117             tmp1      = c[i1]
118             tmp2      = c[i2]
119             c[i1]     = tmp2
120             c[i2]     = tmp1
121         del c
122         return 0
123
124     def init(self, p, pop) :
125         """
126         Random initial states.  We do this so that we can enforce the same
127         initial guesses for all runs to compare against the Poon and Carter.
128         """
129         n = self.GetStringLength()
130         c = self.GetIntegerChromosome(p, pop)
131         np.random.seed(p)
132         perm = np.random.permutation(n)
133         for i in range(0, n) :
134             c[i] = perm[i]
135         del c
136
137 comm = MPI.COMM_WORLD           # Get the communicator.
138 rank = comm.Get_rank()          # Get my rank.
139 t_start = MPI.Wtime()           # Start the clock.
140 n        = 30                   # Number of cities.
141 numrun   = 25                   # Number of runs to average.
142 besteval = np.zeros(numrun)
143 for i in range(0, numrun) :
144     opt = MyPGA(sys.argv, PGA.DATATYPE_INTEGER, n, PGA.MINIMIZE)
```

```
145     opt.SetInitString(opt.init) # Set an initialization operator.
146     opt.SetCrossoverProb(0.8)   # (Default is 85%)
147     opt.SetPopSize(22)          # 22 rather than 21
148     opt.SetNumReplaceValue(21)  # Keep the best 22-21 = 1 string = elitist.
149     opt.SetMaxGAIterValue(300)  # 300 generations, like the reference.
150     opt.SetCrossover(opt.tbx)   # Set a cross-over operation.
151     opt.SetMutation(opt.swap)   # Set a mutate operation.
152     opt.nswap = 3               # Number of pairs to swap in mutation.
153     opt.SetUp()                 # Internal allocations, etc.
154     opt.Run(opt.tsm)            # Set the objective and run.
155     if rank == 0 :
156         best        = opt.GetBestIndex(PGA.OLDPOP)
157         besteval[i] = opt.GetEvaluation(best, PGA.OLDPOP)
158     opt.Destroy()              # Clean up PGAPack internals.
159 
160 if rank == 0 :
161     print "  MEAN: ", np.mean(besteval) # Print out the mean
162     print " SIGMA: ", np.std(besteval)  # and standard deviation.
163     print "   MIN: ", np.min(besteval)  # Print out the mean
164     print "   MAX: ", np.max(besteval)  # Print out the mean
165     t_end = MPI.Wtime()
166     print "Elapsed time = ", t_end-t_start, " seconds."
167 
168 MPI.Finalize() # Should be called automatically, but good practice.
```

Running it using `mpirun -np 1 python example09.py` yields the following output (last several lines):

```
300        Best       9.143837e+02
The Best Evaluation: 9.143837e+02.
The Best String:
#    0: [       7], [      29], [       5], [       2], [      28], [      22]
#    6: [       8], [      10], [      18], [      19], [       3], [      15]
#   12: [      14], [      27], [      16], [       4], [      24], [      20]
#   18: [      17], [      12], [      26], [      25], [       0], [       9]
#   24: [       6], [      13], [      23], [      11], [      21], [       1]

***Destroying PGA context***
  MEAN:  844.962788368
 SIGMA:  97.2494217692
   MIN:  622.625633555
   MAX:  997.810825333
Elapsed time =  67.2148938179  seconds.
```

Running it using `mpirun -np 6 python example09.py` yields the following output (last several lines):

```
Iter #     Field      Value
300        Best       8.004605e+02
The Best Evaluation: 8.004605e+02.
The Best String:
#    0: [       7], [      18], [      17], [      14], [      25], [       5]
#    6: [      19], [      24], [      16], [       2], [       3], [      22]
```

```
#   12: [      21], [       1], [      23], [      26], [       8], [      28]
#   18: [       9], [      12], [      11], [      15], [      20], [      10]
#   24: [      27], [      29], [      13], [       0], [       4], [       6]

***Destroying PGA context***
  MEAN:  832.288188184
 SIGMA:  79.9132513415
   MIN:  597.338150861
   MAX:  942.115139223
Elapsed time =  45.1982860565  seconds.
```

This was using a 6-core machine, but the speedup was barely 25%—why? Well, evaluating the distance of 30 cities is cheap compared to the sorting occuring on the master process for the cross-over operation. As noted earlier, PGAPack's parallelism is definitely meant for expensive evaluations relative to everything else. Here, we see *some* speedup, just not very good. Also compare the mean and standard deviation to the cited 829.00 and 72.69. The parallel results are much closer, which may be a fluke, but it may be worth investigating.

2.3.3 Example 10: Optimizing a Slab Reactor

In this problem, the goal is to optimize a "slab reactor". While the physics is out of our scope, the essential idea is as follows. We have 8 slabs, each 20 cm thick, and each of a different "fuel". These slabs are situated together in some pattern. The pattern is surrounded on either side by water, and beyond the right side water is vacuum while on the left side is an effective mirror, i.e. what goes in must return. This definition makes it so the sequence [0,1,...,7] is different from [7,...,1,0]. What we want to do is to maximize the multiplication factor "keff" (which is related to how long the reactor could produce energy before needing more fuel) and make the power distribution as flat as possible (which in real reactors is related to several important safety margins). The latter is quantified by the "peaking factor", defined as the maximum assembly power divided by the average power of all assemblies, and we seek to minimize this. The objective function is a weighted sum of these objectives.

The crossover used is the heuristic tie-breaking crossover (HTBX) described in Carter, *Advances in Nuclear Science and Technology*., **25**, (1997). The basic idea is similar to the TBX operator of the TSP examples but includes extra problem information, in this case the "reactivity" of the fuel as quantified by "k-infinity". Basically, a more "reactive" fuel produces more neutrons with time, or, in other words, contributes more to the energy production of the reactor. Because of the way the materials are ordered in this example, they are already sorted by reactivity. Hence, TBX and HTBX are identical in implementation for this case.

We run the problem over 100 generations and with a population of size 50. We employ a rather elitist strategy, replacing 40 strings each generation. We also disallow identical strings.

```
"""
pypgapack/examples/example10.py  --  optimize a reflected 8 slab reactor
"""

from pypgapack import PGA
from mpi4py import MPI
import numpy as np
import sys
import matplotlib.pyplot as plt
```

```
from scipy import factorial
from matplotlib import rc
rc('text', usetex=True)
rc('font', family='serif')

class MyPGA(PGA) :
    """
    Derive our own class from PGA.
    """
    def f(self, p, pop) :
        """
        Minimize peaking and maximize keff using weighted objective.
        """
        pattern = self.GetIntegerChromosome(p, pop)
        keff, peak = self.flux(pattern)
        del pattern
        delta = 0
        if peak > 1.8 :
            delta = peak - 1.8
        self.evals += 1
        return 1.0*keff - 10.0*delta

    def flux(self, pattern, plot=0) :
        """
        Solve for the flux via simple mesh-centered finite differences.

        Returns keff and the maximum-to-average assembly-averaged
        fission density ratio.
        """
        # Coarse mesh boundaries.  (Specific to n=8!!)
        coarse = np.array([0.,20.,40.,60.,80.,100.,120.,140.,160.,180.,200.])
        # Fine meshes per coarse mesh.
        fine = 20
        dx = 20.0 / fine
        # Material map (simplifies coefficients)
        mat = np.zeros(fine * (len(coarse) - 1))
        mat[0:fine] = 10 # reflector
        j = fine
        for i in range(0, self.number_slabs) :
            mat[j:(j + fine)] = pattern[i] # one of the slabs
            j += fine
        mat[j:(j + fine)] = 10 # reflector
        # System size
        n = fine * (len(coarse) - 1)
        # Materials
        D, R, F, S = self.materials()
        # Coefficient Matrix (just diagonals) and Vectors
        AD = np.zeros((n, 2))
        AL = np.zeros((n, 2))
        AU = np.zeros((n, 2))
        nufission = np.zeros((n, 2))
        scatter12 = np.zeros(n)
        for g in range(0, 2) :
```

```
            # reflective left boundary
            b = 1
            AU[0, g] = -2.0 * D[g, mat[0]] * D[g, mat[0]] / \
                       (D[g, mat[0]] * dx + D[g, mat[1]] * dx)
            AD[0, g] = -AU[0, g] + 2.0 * D[g, mat[0]] * (1.0 - b) / \
                       (4.0 * D[g, mat[0]] * (1.0 + b) + dx * (1.0 - b)) + \
                       dx * R[g, mat[0]]
            nufission[0, g] = dx * F[g, mat[0]]
            scatter12[0] = dx * S[mat[0]]
            # vacuum right boundary
            b = 0
            AL[n - 2, g] = -2.0 * D[g, mat[n - 1]] * D[g, mat[n - 2]] / \
                           (D[g, mat[n - 1]] * dx + D[g, mat[n - 2]] * dx)
            AD[n - 1, g] = -AL[n - 2, g] + \
                        2.0 * D[g, mat[n - 1]] * (1.0 - b) / \
                       (4.0 * D[g, mat[n - 1]] * (1.0 + b) + \
                       dx * (1.0 - b)) + \
                       dx * R[g, mat[n - 1]]
            nufission[n - 1, g] = dx * F[g, mat[n - 1]]
            scatter12[n - 1] = dx * S[mat[n - 1]]
            # internal cells
            for i in range(1, n - 1) :
                AL[i - 1, g] = -2.0 * D[g, mat[i]] * D[g, mat[i - 1 ]] / \
                               (D[g, mat[i]] * dx + D[g, mat[i - 1]] * dx)
                AU[i, g] = -2.0 * D[g, mat[i]] * D[g, mat[i + 1 ]] / \
                           (D[g, mat[i]] * dx + D[g, mat[i + 1]] * dx)

                AD[i, g] = -(AL[i - 1, g] + AU[i, g]) + dx * R[g, mat[i]]
                nufission[i, g] = dx * F[g, mat[i]]
                scatter12[i] = dx * S[mat[i]]
        # Initiate the fluxes.
        phi1 = np.zeros(n)
        phi2 = np.zeros(n)
        # Use a guess for fission density based on fission cross section.
        fission_density = nufission[:, 1]
        # and normalize it.
        fission_density = fission_density / np.sqrt(np.sum(fission_density**2))
        fission_density0 = np.zeros(n)
        # Initialize the downscatter source.
        scatter_source = np.zeros(n)
        # Initial eigenvalue guess.
        keff = 1
        keff0 = 0
        # Set errors.
        errorfd = 1.0
        errork = 1.0
        it = 0
        while (errorfd > 1e-5 and errork > 1e-5 and it < 200) :
            # Solve fast group.
            self.tridiag(AU[:, 0], AL[:, 0], AD[:, 0], \
                         fission_density / keff, phi1)
            # Compute down scatter source.
            scatter_source = phi1 * scatter12
```

```
116             # Solve thermal group.
117             self.tridiag(AU[:, 1], AL[:, 1], AD[:, 1], scatter_source, phi2)
118             # Keep old values.
119             fission_density0[:] = fission_density
120             keff0 = keff
121             # Update density and eigenvalue
122             fission_density[:] = phi1 * nufission[:, 0] + \
123                                  phi2 * nufission[:, 1]
124             keff = keff0 * np.sum(fission_density) / \
125                            np.sum(fission_density0)
126             # Update errors.  Use Linf norm on density.
127             errorfd = np.max(np.abs(fission_density - fission_density0))
128             errork = np.abs(keff - keff0)
129             it += 1
130         # Now we average the fission density over each fueled coarse mesh.
131         slab_fission_density = np.zeros(self.number_slabs)
132         j = fine
133         for i in range(0, self.number_slabs) :
134             slab_fission_density[i] = \
135                 np.mean(fission_density[j:(j + fine) - 1])
136             j += fine
137         mean_fission_density = np.mean(fission_density)
138         peaking = slab_fission_density / mean_fission_density
139         max_peaking = np.max(peaking)
140         return keff, max_peaking
141 
142     def tridiag(self, U, L, D, f, y):
143         """
144         Tridiagonal solver.
145 
146         This assumes vectors U, L, and D are of the same length.  The right
147         hand side is f and the solved unknowns are returned in y.
148         """
149         N = len(D)
150         w = np.zeros(N)
151         v = np.zeros(N)
152         z = np.zeros(N)
153         w[0] = D[0]
154         v[0] = U[0] / w[0]
155         z[0] = f[0] / w[0]
156         for i in range(1, N) :
157             w[i] = D[i] - L[i - 1] * v[i - 1]
158             v[i] = U[i] / w[i]
159             z[i] = (f[i] - L[i - 1] * z[i - 1]) / w[i]
160         y[N - 1] = z[N - 1]
161         for i in range(N - 2, -1, -1) :
162             y[i] = z[i] - v[i] * y[i + 1]
163 
164     def materials(self):
165         """
166         10 fuels with one 1 reflector by row.  Represents burnup of 0, 5,
167         10, 15, 20, 25, 30, 35, 40, and 45 MWd/kg for one assembly type.
168         """
```

```
169         D = np.array([                                                  \
170             [1.4402e+00, 1.4429e+00, 1.4453e+00, 1.4467e+00, 1.4476e+00,\
171              1.4483e+00, 1.4489e+00, 1.4496e+00, 1.4507e+00, 1.4525e+00,\
172              1.3200e+00],                                               \
173             [3.7939e-01, 3.7516e-01, 3.7233e-01, 3.7045e-01, 3.6913e-01,\
174              3.6818e-01, 3.6749e-01, 3.6699e-01, 3.6649e-01, 3.6615e-01,\
175              2.6720e-01]])
176         R = np.array([                                                  \
177             [2.5800e-02, 2.5751e-02, 2.5755e-02, 2.5840e-02, 2.5958e-02,\
178              2.6090e-02, 2.6226e-02, 2.6362e-02, 2.6559e-02, 2.6802e-02,\
179              2.5700e-02],                                               \
180             [1.1817e-01, 1.2301e-01, 1.2306e-01, 1.2277e-01, 1.2223e-01,\
181              1.2136e-01, 1.2017e-01, 1.1871e-01, 1.1622e-01, 1.1275e-01,\
182              5.1500e-02]])
183         F = np.array([                                                  \
184             [7.9653e-03, 7.6255e-03, 7.2724e-03, 6.9344e-03, 6.6169e-03,\
185              6.3189e-03, 6.0453e-03, 5.7908e-03, 5.4413e-03, 5.0379e-03,\
186              0.00000000],                                               \
187             [1.6359e-01, 1.7301e-01, 1.7681e-01, 1.7634e-01, 1.7355e-01,\
188              1.6931e-01, 1.6424e-01, 1.5869e-01, 1.5005e-01, 1.3894e-01,\
189              0.00000000]])
190         S = np.array(                                                   \
191             [1.5204e-02, 1.5152e-02, 1.4958e-02, 1.4828e-02, 1.4744e-02,\
192              1.4691e-02, 1.4652e-02, 1.4626e-02, 1.4601e-02, 1.4587e-02,\
193               2.3100e-02])
194         return D, R, F, S
195
196     def kinf(self, pattern) :
197         """
198         Compute kinf for each slab.
199         """
200         D, R, F, S = self.materials()
201         k = np.zeros(len(pattern))
202         for i in range(0, len(pattern)) :
203             k[i] = (F[0, i] + F[1, i] * S[i] / R[1, i]) / R[0, i]
204         return k
205
206     def htbx(self, p1, p2, pop1, c1, c2, pop2) :
207         """
208         Heuristic tie-breaking cross-over.  See Carter for details. Actually,
209         for this problem, HTBX is equivalent to TBX, since the materials are
210         already ordered by reactivity.
211         """
212         # Grab the city id's.
213         paren1 = self.GetIntegerChromosome(p1, pop1)
214         paren2 = self.GetIntegerChromosome(p2, pop1)
215         child1 = self.GetIntegerChromosome(c1, pop2)
216         child2 = self.GetIntegerChromosome(c2, pop2)
217         # Copy the parents to temporary vector for manipulation.
218         n = self.GetStringLength()
219         parent1 = np.zeros(n)
220         parent2 = np.zeros(n)
221         for i in range(0, n) :
```

```
            parent1[i] = paren1[i]
            parent2[i] = paren2[i]
        # Code the parents using "position listing".
        code1 = np.zeros(n)
        code2 = np.zeros(n)
        for i in range(0, n) :
            code1[parent1[i]] = i + 1
            code2[parent2[i]] = i + 1
        # Randomly choose two cross-over points.
        perm = np.random.permutation(n)
        point1 = np.min(perm[0:2])
        point2 = np.max(perm[0:2]) + 1
        # Exchange all alleles between the two points.
        temp = np.zeros(point2 - point1)
        for i in range(point1, point2) :
            temp[i - point1] = parent1[i]
            parent1[i] = parent2[i]
            parent2[i] = temp[i - point1]
        # Generate a cross-over map, a random ordering of the 0,1,...,n-1
        crossovermap = np.random.permutation(n)
        # Multiply each allele of the strung by n and add the map.
        parent1 = parent1 * n + crossovermap
        parent2 = parent2 * n + crossovermap
        # Replace the lowest allele by 0, the next by 1, up to n-1.  Here,
        #   we sort the parents first, and then for each element, find
        #   where the increasing values are found in the original.  There
        #   is probably a simpler set of functions built in somewhere.
        sort1 = np.sort(parent1)
        sort2 = np.sort(parent2)
        for i in range(0, n) :
            index = np.where(parent1 == sort1[i])
            parent1[index[0][0]] = i
            index = np.where(parent2 == sort2[i])
            parent2[index[0][0]] = i
        # Map the string back to elements.  These are the offspring.
        tempchild1 = np.zeros(n)
        tempchild2 = np.zeros(n)
        for i in range(0, n) :
            tempchild1[parent1[i]] = i
            tempchild2[parent2[i]] = i
        for i in range(0, n) :
            child1[i] = tempchild1[i]
            child2[i] = tempchild2[i]

    def eog(self):
        """
        Log some data for each generation.
        """
        best                = self.GetBestIndex(PGA.OLDPOP)
        bestpattern         = self.GetIntegerChromosome(best, PGA.OLDPOP)
        iter                = self.GetGAIterValue()
        keff, peak          = opt.flux(bestpattern)
        self.besteval[iter-1] = self.GetEvaluation(best, PGA.OLDPOP)
```

```
275         self.bestkeff[iter-1] = keff
276         self.bestpeak[iter-1] = peak
277 
278 comm = MPI.COMM_WORLD           # Get the communicator.
279 rank = comm.Get_rank()          # Get my rank.
280 size = comm.Get_size()
281 t_start = MPI.Wtime()           # Start the clock.
282 n = 8                           # Number of fueled slabs (1-10)
283 opt = MyPGA(sys.argv, PGA.DATATYPE_INTEGER, n, PGA.MAXIMIZE)
284 opt.SetRandomSeed(1)                      # Set seed for verification.
285 np.random.seed(1)                         # Do the same with Numpy.
286 opt.SetIntegerInitPermute(0, n - 1)       # Start with random permutations.
287 opt.SetPopSize(50)                        # Large enough to see some success.
288 opt.SetNumReplaceValue(40)                # Keep the best half.
289 opt.SetMaxGAIterValue(50)                 # Small number for output.
290 opt.SetCrossover(opt.htbx)                # Set a cross-over operation.
291 opt.SetEndOfGen(opt.eog)                  # End of generation info
292 opt.SetMutation(PGA.MUTATION_PERMUTE)     # Mutate by permutation.
293 opt.SetNoDuplicatesFlag(PGA.TRUE)         # Keep no duplicate patterns.
294 opt.SetUp()                               # Internal allocations, etc.
295 opt.number_slabs = n                      # Must be set to string size.
296 opt.evals = 0
297 opt.besteval = np.zeros(51)
298 opt.bestkeff = np.zeros(51)
299 opt.bestpeak = np.zeros(51)
300 opt.Run(opt.f)
301 
302 if rank > 0:
303     evals = np.array([opt.evals],dtype='i')
304     comm.Send([evals,MPI.INT], dest=0, tag=13)
305 else :
306     evals = np.array([1],dtype='i')
307     for i in range(1, size) :
308         comm.Recv([evals,MPI.INT], source=i, tag=13)
309         opt.evals += evals[0]
310 
311 if rank == 0 :
312     best = opt.GetBestIndex(PGA.OLDPOP) # Get the best string
313     bestpattern = opt.GetIntegerChromosome(best, PGA.OLDPOP)
314     keff, peak = opt.flux(bestpattern)  #   and its keff and peak
315     print " best keff = ", keff, " and peak = ", peak
316     t_end = MPI.Wtime()
317     print "Elapsed time = ", t_end-t_start, " seconds."
318     print "# Evaluations = ", opt.evals
319     plt.plot( np.arange(0,51), opt.besteval, 'b',                          \
320               np.array([0,51]), np.array([1.06611227 , 1.06611227]), 'g--',\
321               lw=2) # Plot the objective as a function of generations
322                     #   against the reference solution.
323     plt.title('Convergence of Objective')
324     plt.xlabel(' generation')
325     plt.ylabel(' objective ')
326     plt.legend(('HTBX', 'solution'),loc=4, shadow=True)
327     plt.grid(True)
```

```
328     plt.show()
329
330 MPI.Finalize()
331 opt.Destroy()
```

Running it using `mpirun -np 1 python example10.py` yields the following output:

```
***Constructing PGA***
Iter #     Field      Value
10         Best       1.060239e+00
Iter #     Field      Value
20         Best       1.065283e+00
Iter #     Field      Value
30         Best       1.065808e+00
Iter #     Field      Value
40         Best       1.065808e+00
Iter #     Field      Value
50         Best       1.065808e+00
The Best Evaluation: 1.065808e+00.
The Best String:
#    0: [       4], [       2], [       1], [       5], [       3], [       0]
#    6: [       7], [       6]

 best keff =  1.06416554234  and peak =  2.00346511469
Elapsed time =  179.573677063  seconds.
# Evaluations =  1990
***Destroying PGA context***
```

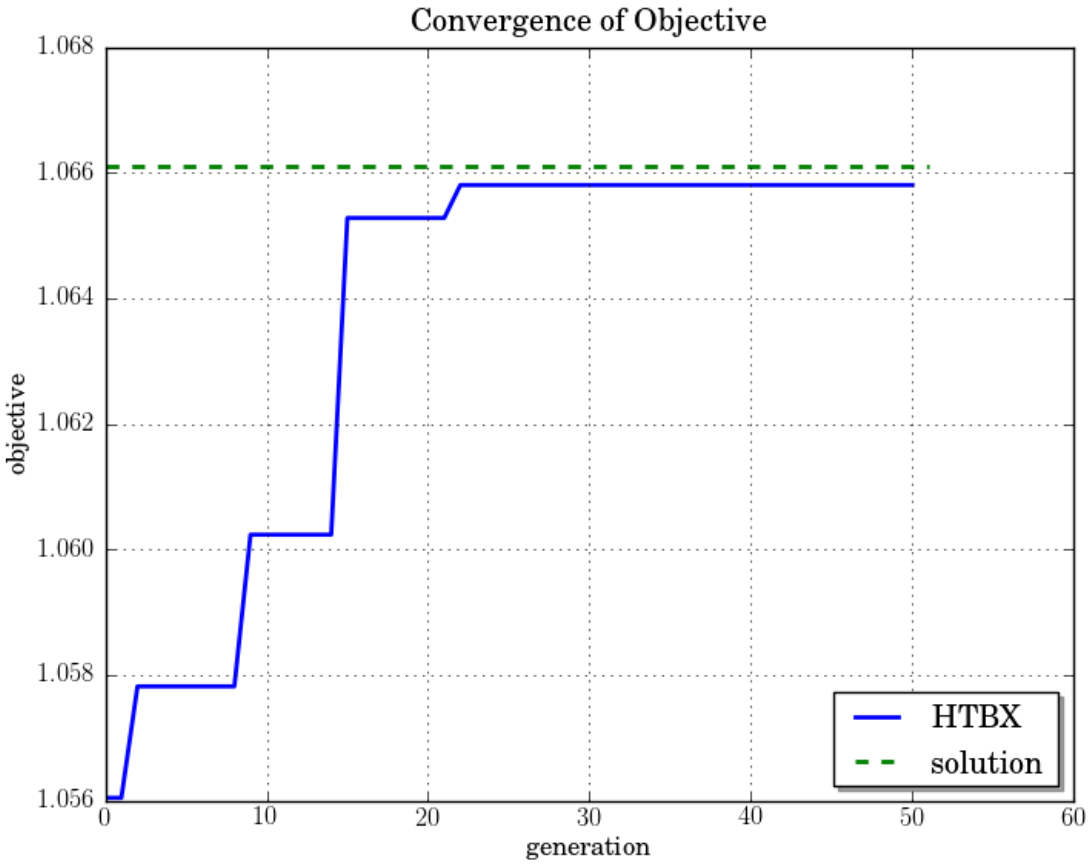
Running it using `mpirun -np 2 python example10.py` yields the following output:

```
***Constructing PGA***
***Constructing PGA***
Iter #     Field      Value
10         Best       1.060239e+00
Iter #     Field      Value
20         Best       1.065283e+00
Iter #     Field      Value
30         Best       1.065808e+00
Iter #     Field      Value
40         Best       1.065808e+00
Iter #     Field      Value
50         Best       1.065808e+00
The Best Evaluation: 1.065808e+00.
The Best String:
#    0: [       4], [       2], [       1], [       5], [       3], [       0]
#    6: [       7], [       6]

 best keff =  1.06416554234  and peak =  2.00346511469
Elapsed time =  108.014204025  seconds.
# Evaluations =  1990
***Destroying PGA context***
***Destroying PGA context***
```

Running it using `mpirun -np 5 python example10.py` yields the following output:

```
***Constructing PGA***
***Constructing PGA***
***Constructing PGA***
***Constructing PGA***
***Constructing PGA***
Iter #     Field      Value
10         Best       1.060239e+00
Iter #     Field      Value
20         Best       1.065283e+00
Iter #     Field      Value
30         Best       1.065808e+00
Iter #     Field      Value
40         Best       1.065808e+00
Iter #     Field      Value
50         Best       1.065808e+00
The Best Evaluation: 1.065808e+00.
The Best String:
#    0: [       4], [       2], [       1], [       5], [       3], [       0]
#    6: [       7], [       6]

 best keff =  1.06416554234  and peak =  2.00346511469
Elapsed time =  51.1656098366  seconds.
# Evaluations =  1990
***Destroying PGA context***
***Destroying PGA context***
***Destroying PGA context***
***Destroying PGA context***
***Destroying PGA context***
```

The reference solution for this objective is 1.06611227 , which was found by directly solving each of the ~80000 possible solutions. As we can observe, the GA does quite well. The parallel performance is also quite good, which makes sense as this problem has a comparitively expensive evaluation function. Note that for `np` above 2, PGApack uses the master process for communications, and hence `np=5` has just 4 compute processes.


Convergence of Objective
objective
generation
1.068
1.066
1.064
1.062
1.060
1.058
1.056
0
10
20
30
40
50
60
HTBX
solution

CHAPTER

THREE

METHODS

To be completed.

CHAPTER

FOUR

API REFERENCE

4.1 Introduction

This section provides a reference for all functions defined in the `pypgapack` module that are *extensions* of the basic PGAPack API. All PGAPack functions are contained in the `pypgapack.PGA` class. The PGAPack library is typically used as follows:

```
double evaluate(PGAContext *ctx, int p, int pop);
PGAContext *ctx;
ctx = PGACreate(&argc, argv, PGA_DATATYPE_BINARY, 100, PGA_MAXIMIZE);
PGASetUp(ctx);
PGARun(ctx, evaluate);
PGADestroy(ctx);
```

The `ctx` object is created explicitly by the user and then passed as the first argument to all subsequent function calls, with function names taking the form `PGAxxx`. For `pypgapack`, `ctx` is a *private* member of `PGA` created during construction, and all `PGA` members drop the *PGA* prefix and the initial `ctx` argument. So, for example,

```
ctx = PGACreate(&argc, argv, PGA_DATATYPE_BINARY, 100, PGA_MAXIMIZE);
PGASetUp(ctx);
```

in C/C++ becomes

```
obj = pypgapack.PGA(sys.argv, PGA.DATATYPE_BINARY, 100, PGA.MAXIMIZE)
obj.SetUp()
```

in Python. For all functions included in PGAPack, the user is directed to the pgapack documentation. What follows is a description of the few new methods added for `pypgapack` that make life in Python a bit easier.

4.2 pypgapack API

The easiest way to see what `pypgapack` offers is to do the following:

```
>>> import pypgapack as pga
>>> dir(pga)
['PGA', 'PGA_swigregister', '__builtins__', '__doc__', '__file__',
```

```
 '__name__', '__package__', '_newclass', '_object', '_pypgapack',
 '_swig_getattr', '_swig_property', '_swig_repr', '_swig_setattr',
 '_swig_setattr_nondynamic']
```

This command works with any Python module. Our interest is in the `PGA` class. We do the same for this:

```
>>> dir(pga.PGA)
['BinaryBuildDatatype', 'BinaryCopyString', 'BinaryCreateString',
 'BinaryDuplicate', 'BinaryHammingDistance', 'BinaryInitString',
 'BinaryMutation', 'BinaryOneptCrossover', 'BinaryPrint',
 'BinaryPrintString', 'BinaryTwoptCrossover', 'BinaryUniformCrossover',...
```

and find a really long list of class members, most of which are directly from PGAPack. In the following, we document only those not included in PGAPack, as use of the PGAPack functionality is covered above (i.e. drop the `ctx` argument and `PGA` prefix).

class `PGA`

PGA wrapper class.

`__init__`(*argv*, *datatype*, *n*, *direction*)

Construct the PGA context. This essentially wraps the PGACreate function, so see the PGAPack documentation.

Parameters

- **argv** – system argument
- **datatype** – allele dataype; can be `PGA.DATATYPE_XXX`, where `XXX` is `BINARY`, `INTEGER`, and so on.
- **n** – size of the unknown, i.e. number of alleles of type `datatype`
- **direction** – either `PGA.MAXIMIZE` or `PGA.MINIMIZE`

`GetIntegerChromosome`(*p*, *pop*)

Get direct access to the *p*-th integer chromosome string in population *pop* .

Parameters

- **p** – string index
- **pop** – population index

Returns string as numpy array of integers

`GetRealChromosome`(*p*, *pop*)

Get direct access to the *p*-th double chromosome string in population *pop*.

Parameters

- **p** – string index
- **pop** – population index

Returns string as numpy array of floats

`SetInitString`(*f*)

Set a function for initializing strings. The function `f` provided **must** have the signature `f(p,`

`pop)`, but should almost certainly be an inerited class member with the signature `f(self, p, pop)`. See PGAPack documentation for more about user functions.

Parameters **f** – Python function

See Also:

Example 4: User-defined String Initialization for an example on string initialization.

`SetCrossover`(*f*)

Set a function for the crossover operation. The function `f` provided **must** have the signature `f(a,b,c,d,e,f)`, but should almost certainly be an inerited class member with the signature `f(self,a,b,c,d,e,f)`. See PGAPack documentation for more about user functions.

Parameters **f** – Python function

See Also:

Example 5: User-defined Crossover Operator for an example on setting the crossover operator.

`SetMutation`(*f*)

Set a function for the mutation operator. The function `f` provided **must** have the signature `f(p, pop, prob)`, but should almost certainly be an inerited class member with the signature `f(self, p, pop, prob)`. See PGAPack documentation for more about user functions.

Parameters **f** – Python function

See Also:

Example 6: User-defined Mutation Operator for an example on setting the mutation operator.

`SetEndOfGen`(*f*)

Set a function for an operator to be performed at the end of each generation. The function `f` provided **must** have the signature `f(pop)`, but should almost certainly be an inerited class member with the signature `f(self, pop)`. Such an operator can be used to implement hill-climbing heuristics. See PGAPack documentation for more about user functions.

Parameters **f** – Python function

See Also:

Example 7: User-defined End of Generation Operator for an example on setting the an end of generation operator.

CHAPTER

FIVE

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CHAPTER

SIX

INDICES AND TABLES

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