0.1 Thermal calculations on images

```
from numba import njit
from numba.experimental import jitclass
from numba import int64
```

```
import numpy as np
from scipy import interpolate, special
import os
import time
import tempfile
import h5py, hickle
from multiprocessing import Pool
from pprint import pprint
```

```
# We extend the path instead of using `src.module` to be able to run generated files.
import sys
if 'src' not in sys.path: sys.path.append('src')
import wanglandau as wl
```

```
integer = int64
spec = [
    ('I0', integer[:]),
    ('I',  integer[:]),
    ('N',  integer),
    ('M',  integer),
#     ('Es', integer[:]),
    ('E',  integer),
    ('Ev', integer),
    ('dE', integer),
    ('dx', integer),
    ('i',  integer)
]

@jitclass(spec)
class StatisticalImage:
    def __init__(self, I0, I, M):
        if len(I0) != len(I):
            raise ValueError('Ground image I0 and current image I should have the same length.')
        if M < 0:
            raise ValueError('Maximum site value must be nonnegative.')
        self.I0 = I0
        self.I  = I
        self.N  = len(I0)
        self.M  = M
#         self.Es = self.energy_bins()
```

```
        self.E  = self.energy()
        self.Ev = self.E
        self.dE = 0
        self.dx = 0
        self.i  = 0
    def state(self):
        return self.I0.copy(), self.I.copy(), self.M
    def state_names(self):
        return 'I0', 'I', 'M'
    def copy(self):
        return StatisticalImage(*self.state())
    def energy_bins(self):
        E0 = 0
        Ef = np.sum(np.maximum(self.I0, self.M - self.I0))
        ΔE = 1
        return np.arange(E0, Ef + ΔE + 1, ΔE)
    def energy(self):
        return np.sum(np.abs(self.I - self.I0))
    def propose(self):
        i = np.random.randint(self.N)
        self.i = i
        x0 = self.I0[i]
        x = self.I[i]
        r = np.random.randint(2)
        if x == 0:
            dx = r
        elif x == self.M:
            dx = -r
        else:
            dx = 2*r - 1
        dE = np.abs(dx) if x0 == x else (dx if x0 < x else -dx)
        self.dx = dx
        self.dE = dE
        self.Ev = self.E + dE
    def accept(self):
        self.I[self.i] += self.dx
        self.E = self.Ev
```

0.1.1 Parallel Simulation

```
N = 16
Moff = 0
I0 = Moff * np.ones(N, dtype=int)
system_parameters = {
    'I0': I0,
```

```
    'I': I0.copy(),
    'M': 2**5 - 1
}
parallel_parameters = {
    'bins': 8,
    'overlap': 0.5,
    'steps': 1_000_000
}
wl_parameters = {
    'M': 1_000_000,
    'ε': 1e-10,
    'logf0': 1,
    'flatness': 0.1,
    'logging': False
}

system = StatisticalImage(**system_parameters) # Intermediate value
Es = system.energy_bins()
```

```
print('Parallel Wang-Landau simulation with')
for k, v in wl_parameters.items():
    print("\t", k, '\t', v)
print('on a {}.'.format(system.__class__.__name__))
```

```
Parallel Wang-Landau simulation with
     M   1000000
     ε   1e-10
     logf0   1
     flatness    0.1
     logging     False
on a StatisticalImage.
```

```
def parallel_wanglandau(subsystem): # Convenient form for `Pool.map`
    wl.urandom_reseed()
    state, Es = subsystem
    system = StatisticalImage(*state)
    print('(', end='', flush=True)
#     results = wl.wanglandau(system, Es, M = wlM, ε = 1e-8, logging=False)
    results = wl.wanglandau(system, Es, **wl_parameters)
    print(')', end='', flush=True)
    return results
```

```
print('Finding parallel bin systems ... ', end='', flush=True)
psystems = wl.parallel_systems(system, Es, **parallel_parameters)
print('done.')
```

```
Finding parallel bin systems ... done.
```

```
print('Running | ', end='', flush=True)
start_time = time.time()
with Pool() as pool:
    wlresults = pool.map(parallel_wanglandau, psystems)
print(' | done in', int(time.time() - start_time), 'seconds.')
```

```
Running | (
```

```
sEs, sS = wl.stitch_results(wlresults)
```

```
with tempfile.NamedTemporaryFile(mode='wb', prefix='wlresults-image-', suffix='.hdf5',
↪  dir='data', delete=False) as f:
    with h5py.File(f, 'w') as hkl:
        print('Writing results ... ', end='', flush=True)
        d = {}
        for k, v in zip(system.state_names(), system.state()):
            d.update({k: v})
        hickle.dump({
            'parameters': {
                'system': d,
                'wanglandau': {},
                'parallel': {}
            },
            'results': {
                'Es': sEs,
                'S': sS
            },
            'parallel_results': wlresults
        }, hkl)
        print('done: {}'.format(os.path.relpath(f.name)))
```

0.1.2 Results

```
import matplotlib.pyplot as plt
```

```
N, M = len(system_parameters['I0']), system_parameters['M']
```

```
for Es, S, H in wlresults:
    plt.plot(Es[:-1], S)
```

```
wlEs, S = sEs[:-1], sS
```

Fit a spline to interpolate and optionally clean up noise, giving WL g's up to a normalization constant.

```
gspl = interpolate.splrep(wlEs, S, s=0*np.sqrt(2))
wlgs = np.exp(interpolate.splev(wlEs, gspl) - min(S))
```

0.1.3 Exact solution

We only compute to halfway since g is symmetric and the other half's large numbers cause numerical instability.

```
def reflect(a, center=True):
    if center:
        return np.hstack([a[:-1], a[-1], a[-2::-1]])
    else:
        return np.hstack([a, a[::-1]])
```

The exact density of states for uniform values. This covers the all gray and all black/white cases. Everything else (normal images) are somewhere between. The gray is a slight approximation: the ground level is not degenerate, but we say it has degeneracy 2 like all the other sites. For the numbers of sites and values we are using, this is insignificant.

```
def bw_g(E, N, M, exact=True):
    return sum((-1)**k * special.comb(N, k, exact=exact) * special.comb(E + N - 1 - k*(M + 1), E
    ↪  - k*(M + 1), exact=exact)
        for k in range(int(E / M) + 1))
def exact_bw_gs(N, M):
    Es = np.arange(N*M + 1)
    gs = np.vectorize(bw_g)(np.arange(1 + N*M // 2), N, M, exact=False)
    return Es, reflect(gs, len(Es) % 2 == 1)
```

```
def gray_g(E, N, M, exact=True):
    return 2 * bw_g(E, N, M, exact=exact)
def exact_gray_gs(N, M):
    Es = np.arange(N*M + 1)
    gs = np.vectorize(gray_g)(np.arange(1 + N*M // 2), N, M, exact=False)
    return Es, reflect(gs, len(Es) % 2 == 1)
```

Expected results for black/white and gray.

```
bw_Es, bw_gs = exact_bw_gs(N=N, M=M)
gray_Es, gray_gs = exact_gray_gs(N=N, M=-1 + (M + 1) // 2)
```

Choose what to compare to.

```
Es, gs = bw_Es, bw_gs
```

Presumably all of the densities of states for different images fall in the region between the all-gray and all-black/white curves.

```
plt.plot(bw_Es / len(bw_Es), np.log(bw_gs), 'black', label='BW')
plt.plot(gray_Es / len(gray_Es), np.log(gray_gs), 'gray', label='Gray')
plt.plot(wlEs / len(wlEs), np.log(wlgs), label='WL')
plt.xlabel('E / MN')
plt.ylabel('ln g')
plt.title('N = {}, M = {}'.format(N, M))
plt.legend();
```

```
plt.plot(wlEs / len(wlEs), np.abs(wlgs - bw_gs) / bw_gs)
plt.title('Relative error')
plt.xlabel('E / MN')
plt.ylabel('ε(S)');
```

```
print('End of job.')
```