
Weekly Report

June 21-June 27

Christina

Monte-Carlo Simulation

Some improvements:

1. Preprocess function: added `rotate_pads=True` for the simulated data only
 - a. The real & simulated data are now on the same coordinates
2. Commented out the “process shifted z by 38mm due to trigger delay” command in the preprocess function
 - a. Rather nice fitting for event_305
 - b. However, some events who do go through trigger delays now experience a shift in z-direction

Gradient Descent - a local optimization method

Gradient Descent Principles

1. We choose a starting point within a function , and then take steps towards the direction with negative (approximate) gradient
2. Conjugate gradient: used to approximate the answer \mathbf{x} to the equation $Q\mathbf{x} + \mathbf{b} = \mathbf{g}$, where Q is a symmetrical matrix.
 - a. Let $\{d_0, \dots, d_{n-1}\}$ be a set of nonzero Q -orthogonal vectors, the sequence $\mathbf{x}_{k+1} = \mathbf{x}_k + a_k * d_k$ (where a is step size) converges to a unique solution $Q\mathbf{x} + \mathbf{b} = 0$ that minimizes the function.

Chi-square function

1. The function we used to evaluate the chi-square value is `minimizer.run_tracks`
 - a. Takes in our initial guess (in the form of a 2D array), the experimental data points (x y z) and the experimental hit pattern (pad amplitudes indexed by pad number).
 - b. Returns an array consisting of the three chi-square values (position, energy, vertex components)

Scipy function construction

1. We define a function $f(y)$ that can be accepted by `Scipy.optimize` functions

```
def f(y): #accept y as a vector consisting of six parameters
```

```
    ctr = np.zeros([1,6])
```

```
    ctr[0] = y #transform y to a 2D array acceptable by run_tracks function
```

```
    chi_result = minimizer.run_tracks(ctr, exp_pos, exp_hits)
```

```
    return sum(chi_result[0]) #We want to minimize the sum
```

Scipy function construction

2. We define the jacobian with a numerical derivative function `scipy.optimize.approx_fprime` so we can implement it into our optimization function to get a more accurate result
3. We then create a callback function (callbackF) which prints out the intermediate result of each iteration when we call the optimization function
4. Run the optimization function
 - a. `results = scipy.optimize.minimize(f #object function, ctr0 #initial guess, method="CG", jac=jacobian, callback=callbackF, options={'eps':1e-13} #step size)`

Differential evolution - a global optimization method

Differential Evolution Principles

1. The method iteratively try to improve a existing candidate solution -> stochastic method, the best solution is not guaranteed to be found
2. Starts with a “population” of candidate solutions, points with random positions in the search space.
3. For each candidate, pick two distinct candidates (a, b) from the population
4. Compute the candidate’s potentially new position (aka trial candidate):
 - a. We want to “mutate” the best candidate c_0
 - b. $C' = c_0 + \text{mutation} * (a - b)$
 - c. For ‘best1bin’ strategy, we take a random number in $[0,1)$ and if it is less than the recombination constant (explained later) the parameters is loaded from c' (else it’s loaded from c_0).
 - d. If the trial candidate c' is better than the original candidate c_0 , it replaces c_0 .

Scipy function parameters

1. Bounds: (min, max) pairs for each element in our input vector. `bounds = [(-1,1), (-1, 1), (0, 1), (0,5), (-2 * pi, 2 * pi), (-2 * pi, 2 * pi)]`
2. Maxitr: # of iterations allowed
3. Popsiz: population size. Default is 15; larger size improves chance of convergence
4. Mutation: a float or a range - larger mutation value increases the search radius but will slow down convergence. If mutation is given a range, the constant is randomly changed per iteration to increase speed.
5. Recombination constant: float in [0,1]. Increased RC will allow more mutants to become trial candidates, but will decrease population stability

Results

Event #305

1. Conjugate Gradient $\chi = 24.713$

```
results = scipy.optimize.minimize(f, ctr0, method="CG", jac=jacobian, callback=callbackF, options={'gtol': 30.0, 'eps': 1e-10})
cg = results.x
```

```
302 0.000360-0.000141 0.734463 2.614517-3.002859 1.951079 24.713554
303 0.000360-0.000141 0.734463 2.614517-3.002859 1.951079 24.713554
304 0.000360-0.000141 0.734463 2.614517-3.002859 1.951079 24.713554
305 0.000360-0.000141 0.734463 2.614517-3.002859 1.951079 24.713554
306 0.000360-0.000141 0.734463 2.614517-3.002859 1.951079 24.713554
307 0.000360-0.000141 0.734463 2.614517-3.002859 1.951079 24.713554
308 0.000360-0.000141 0.734463 2.614517-3.002859 1.951079 24.713554
309 0.000360-0.000141 0.734462 2.614516-3.002859 1.951079 24.713481
310 0.000360-0.000141 0.734462 2.614516-3.002858 1.951079 24.713390
311 0.000360-0.000141 0.734462 2.614515-3.002858 1.951079 24.713326
312 0.000360-0.000141 0.734462 2.614515-3.002858 1.951079 24.713326
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317 0.000360-0.000141 0.734462 2.614515-3.002858 1.951079 24.713326
318 0.000360-0.000141 0.734462 2.614515-3.002858 1.951079 24.713326
319 0.000360-0.000141 0.734462 2.614515-3.002858 1.951079 24.713326
320 0.000360-0.000141 0.734462 2.614515-3.002858 1.951079 24.713326
```

2. Monte-Carlo

```
uvw, (cu, cv) = mcfitter.preprocess(xyzs[:,0:5], center=(cx, cy), rotate_pads=False) # get calibrated set of data
uvw_values = uvw.values #transform pd file to arrays
print(cu,cv)
mcres, minChis, all_params, good_param_idx = mcfitter.process_event(uvw, cu, cv, return_details=True)
|
print(mcre)
```

```
-43.91625428724363 118.39379347393248
```

```
{'x0': 0.004871498439906168, 'y0': -0.011977943314897467, 'z0': 0.7347184090083536, 'enu0': 2.8186679462143145, 'azi0': -2.920989043120735, 'pol0': 1.9263176741477495, 'posChi2': 3.97700712247824, 'enChi2': 7.1981754856024605, 'vertChi2': 3.344052462098103, 'lin_scatt_ang': 1.2420273973461806, 'lin_beam_int': 720.0566082022592, 'lin_chi2': 36.34021928440234, 'rad_curv': 117.8081237563669, 'brho': 0.20911788771969209, 'curv_en': 2.0944258199546852, 'curv_ctr_x': -43.91625428724363, 'curv_ctr_y': 118.39379347393248}
```

$$\text{Chi} = 3.977 + 7.198 + 3.344 = 14.519$$

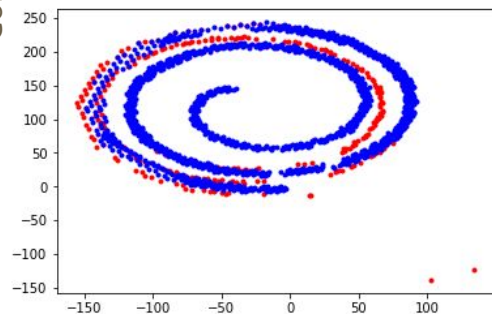
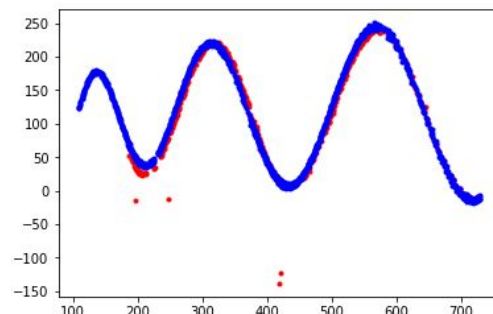
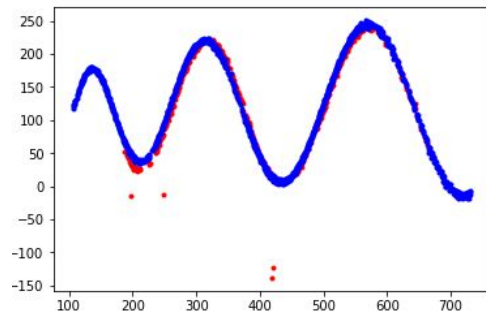
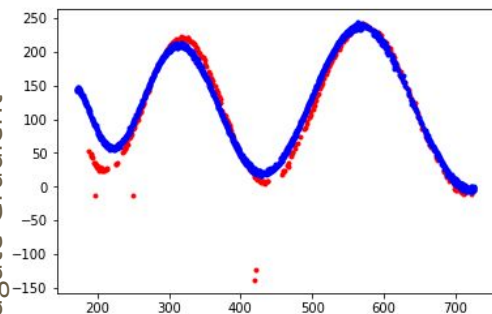
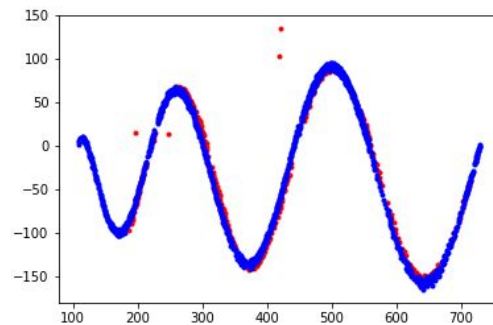
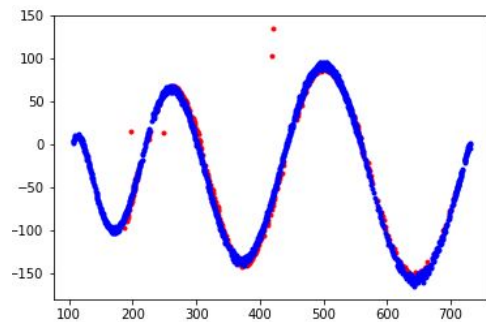
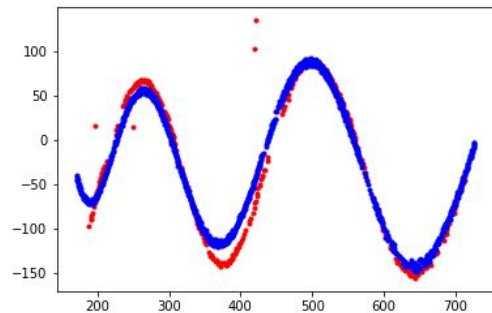
3. Differential evolution

chi=14.208

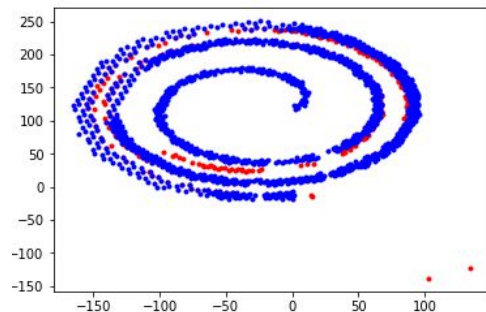
```
i = 1
def callbackF(x, convergence=10):
    global i
    print('{0:4d}{1: 3.6f}{2: 3.6f}{3: 3.6f}{4: 3.6f}{5: 3.6f}{6: 3.6f}{7: 3.6f}'.format(i,x[0],x[1],x[2],x[3],x[4],x[5],
    i += 1
bounds = [(-1,1), (-1, 1), (0, 1), (0,5), (-2 * pi, 2 * pi), (-2 * pi, 2 * pi)]
results = scipy.optimize.differential_evolution(f, bounds, callback=callbackF, maxiter=10000, strategy='best1bin',\
    recombination=0.8, popsize=15, mutation=(0.5, 1.0))

'''    The differential evolution strategy should be one of:
        'best1bin' - shorter time but usually needs to run multiple times to get the best result
        'best1exp' - shorter time but usually needs to run multiple times to get the best result
        'randlexp' - takes a long time but converges
        'randtobest1exp' - takes a long time but converges
        'currenttobest1exp' - takes a long time but converges
        'best2exp'
        'rand2exp'
        'randtobest1bin'
        'currenttobest1bin'
        'best2bin'
        'rand2bin'
        'rand1bin'
    The default is 'best1bin'.
'''

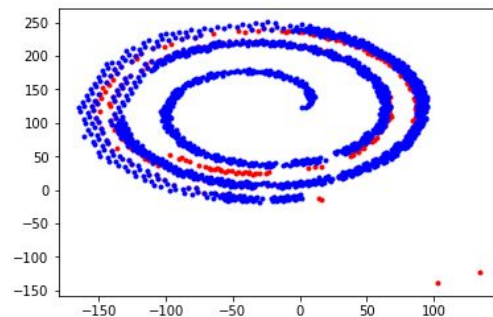
94 0.003659-0.012275 0.733378 2.809433 0.211693-1.926856 14.577127
95 0.003659-0.012275 0.733378 2.809433 0.211693-1.926856 14.577127
96 0.003659-0.012275 0.733378 2.809433 0.211693-1.926856 14.577127
97 0.002564-0.011759 0.733055 2.809754 0.207638-1.926092 14.260398
98 0.002564-0.011759 0.733055 2.809754 0.207638-1.926092 14.260398
99 0.002564-0.011759 0.733055 2.809754 0.207638-1.926092 14.260398
100 0.002564-0.011759 0.733055 2.809754 0.207638-1.926092 14.260398
101 0.002564-0.011759 0.733055 2.809754 0.207638-1.926092 14.260398
102 0.002564-0.011759 0.733055 2.809754 0.207638-1.926092 14.260398
103 0.002564-0.011759 0.733055 2.809754 0.207638-1.926092 14.260398
104 0.002306-0.011989 0.732534 2.808862 0.204325-1.926166 14.207878
105 0.002306-0.011989 0.732534 2.808862 0.204325-1.926166 14.207878
106 0.002306-0.011989 0.732534 2.808862 0.204325-1.926166 14.207878
```

Monte Carlo



Differential Evolution



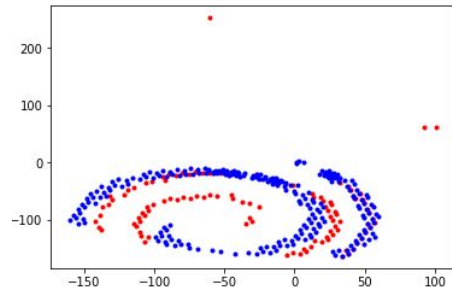
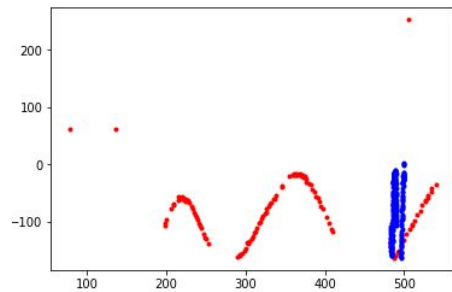
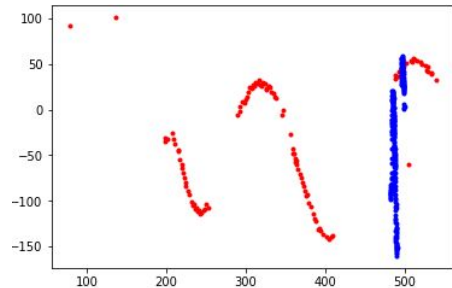
Event #456

Monte Carlo: $\chi = 12.365 + 7.088 + 0.033 = 19.486$

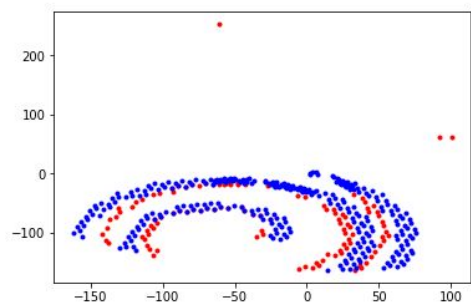
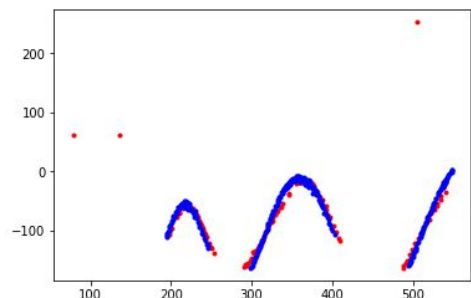
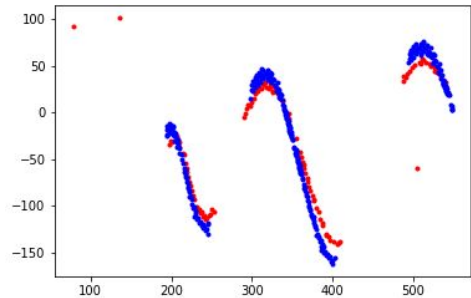
Gradient: 110.8 (fail)

Differential evolution: with population size=30,
recombination constant = 0.6, $\chi = 27.32$

Conjugate Gradient



Monte Carlo



Differential Evolution

