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>>> PTC - PyORBIT SIS18 Benchmark  
>>> An Iterative Journey
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>>> What's the SIS18 Benchmark?

- \* Created by Giuliano Franchetti to confirm space charge induced trapping of particles in a bunch during long term storage.

>>> Why should I care?

- \* We rely on PTC-PyORBIT for space charge simulations, so it's good to check that it is reliable.

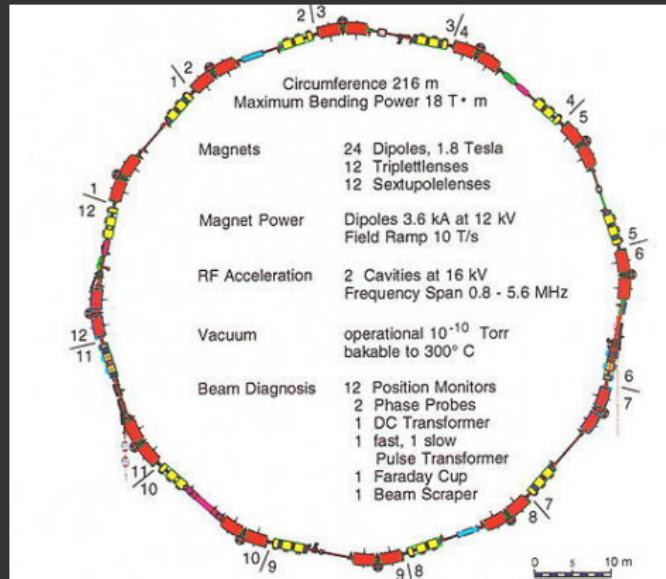
## >>> Motivation

- \* To verify that PTC-PyORBIT behaves as expected when considering physical effects that should be present for space charge studies.
- \* To understand the necessary ingredients for a space charge simulation with PTC-PyORBIT.
- \* To provide those that wish to use PTC-PyORBIT with nicely commented code examples.



## >>> The Accelerator

This benchmark uses the SIS18 (German SchwerIonenSynchrotron) lattice. In the benchmark lattice only one sextupole is present.



The reference lattice can be found [here](#), and the usable MAD-X lattice [here](#).



>>> Beam parameters (steps 1 - 6)

Parameter	Symbol	Value	Unit
Sextupole Strength (when used)	$K_2$	0.2	$m^{-2}$
Maximum Tuneshift	$\Delta Q_x$	0.1	-
Intensity	$N_p$	$2.95 \cdot 10^9$	particles
Horizontal Emittance	$\epsilon_x$	$4.91 \cdot 10^{-7}$	$m \ rad$
Vertical Emittance	$\epsilon_y$	$3.635 \cdot 10^{-7}$	$m \ rad$
Kinetik Energy	$E_k$	11.4	$\frac{MeV}{u}$
Gamma Transition	$\gamma_t$	5	-
Momentum Spread	$\frac{dp}{p}$	$2.5 \cdot 10^{-4}$	-
Lorentz Gamma	$\gamma$	1.012149995	-
Lorentz Beta	$\beta$	0.15448	-
Bunch Length	$\tau$	3472.7	ns
Synchrotron Tune	$Q_{synch}$	$\frac{1}{15000}$	-

Benchmark Information

Steps 1 - 6 Parameters

>>> Beam parameters (steps 7 - 9)

Parameter	Symbol	Value	Unit
Sextupole Strength (when used)	$K_2$	0.2	$m^{-2}$
Maximum Tuneshift	$\Delta Q_x$	0.1	-
Intensity	$N_p$	$1.95 \cdot 10^9$	particles
Horizontal Emittance	$\epsilon_x$	$4.91 \cdot 10^{-7}$	$m \ rad$
Vertical Emittance	$\epsilon_y$	$3.635 \cdot 10^{-7}$	$m \ rad$
Kinetik Energy	$E_k$	11.4	$\frac{MeV}{u}$
Gamma Transition	$\gamma_t$	5	-
Momentum Spread	$\frac{dp}{p}$	$2.5 \cdot 10^{-4}$	-
Lorentz Gamma	$\gamma$	1.012149995	-
Lorentz Beta	$\beta$	0.15448	-
Bunch Length	$\tau$	231.51	ns
Synchrotron Tune	$Q_{synch}$	$\frac{1}{1000}$	-

Benchmark Information  
Steps 7 - 9 Parameters



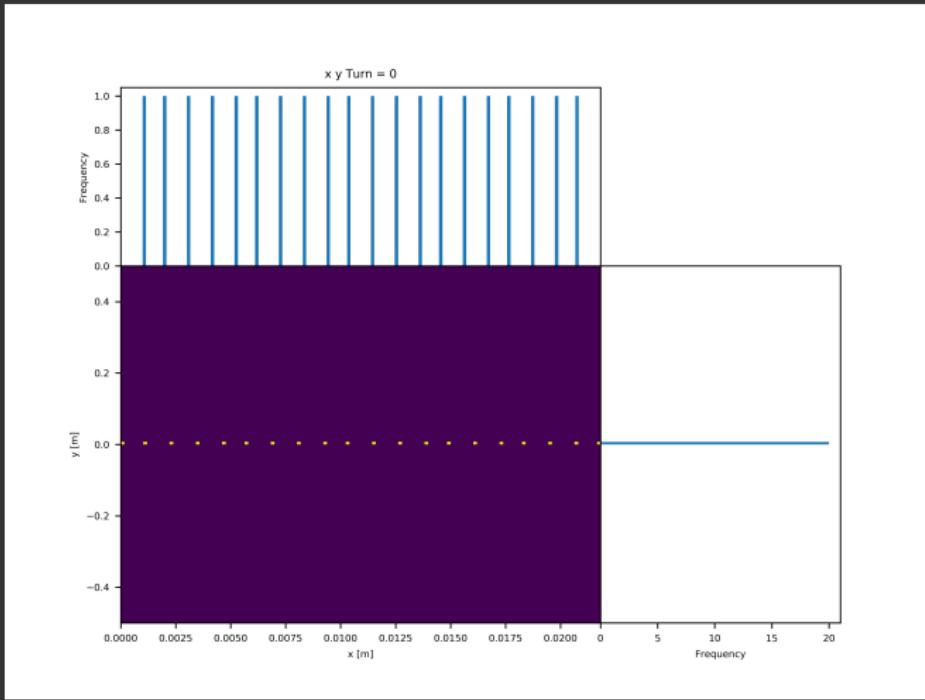
## >>> Particle Distributions: Poincaré

Poincaré distn:

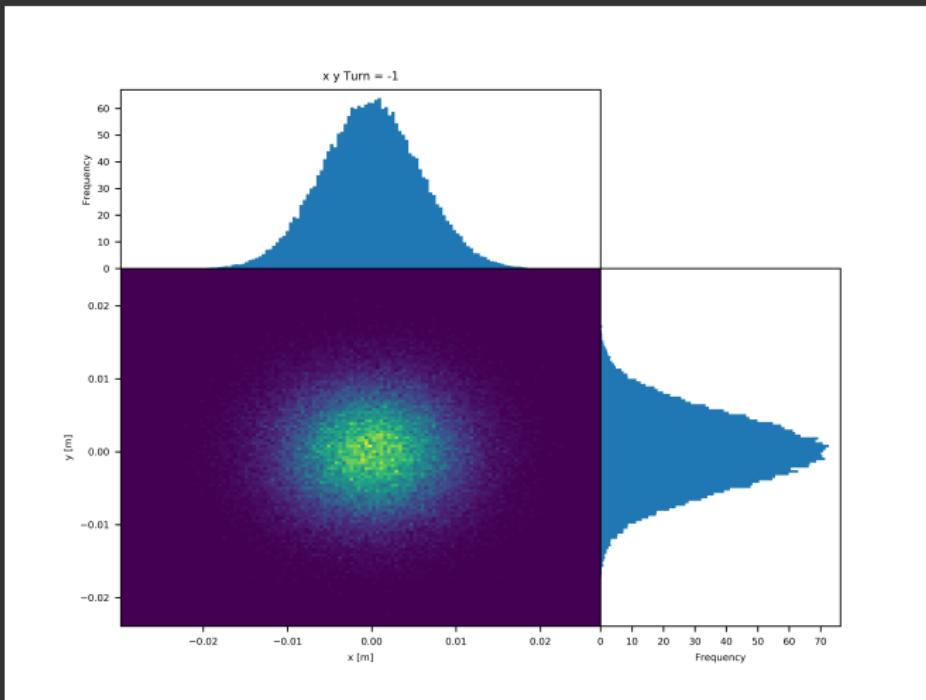
$$x_i = \frac{i\sigma_n}{N} \sqrt{\frac{\beta_{x0}\epsilon_x}{\beta\gamma}} \quad (1)$$

A simple approach - select a maximum amplitude in sigma  $\sigma_n$ , equally spaced (from 0) over the number of particles  $N$ , thus  $i$  is an integer iterator where we loop over  $i$  particles. Then the particle position is the square root of the beta function at the location of bunch creation and emittance (the local beam size)  $\beta_{x0}\epsilon_x$ , normalised with the Lorentz  $\beta\gamma$ .

## >>> Particle Distributions: Poincaré



### >>> Particle Distributions: 3D Gaussian



## >>> Linear Restoring Force

A linear restoring force was used to imitate the synchrotron motion:

The function is simple:

$$dE_i = dE_i + F \cdot z_i$$

$dE_i$  = energy offset of particle  $i$

$z_i$  =  $z$  co-ordinate of particle  $i$

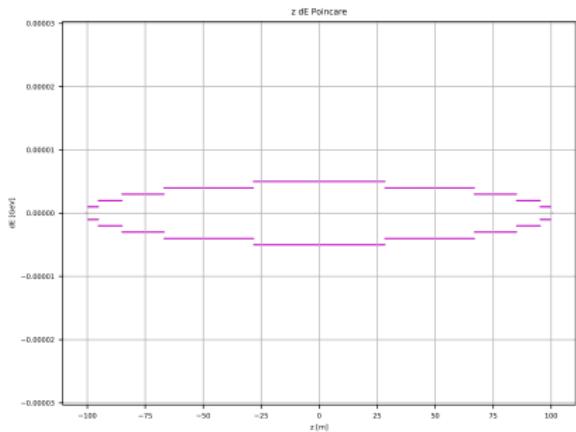
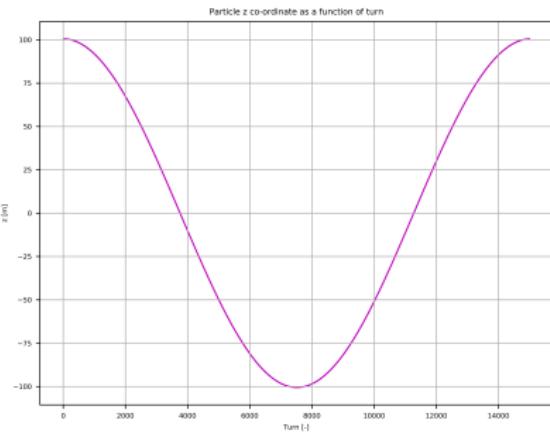
$F$  = restoring force

$F$  may be found analytically using an expression for the synchrotron tune, or empirically.

## >>> Linear Restoring Force

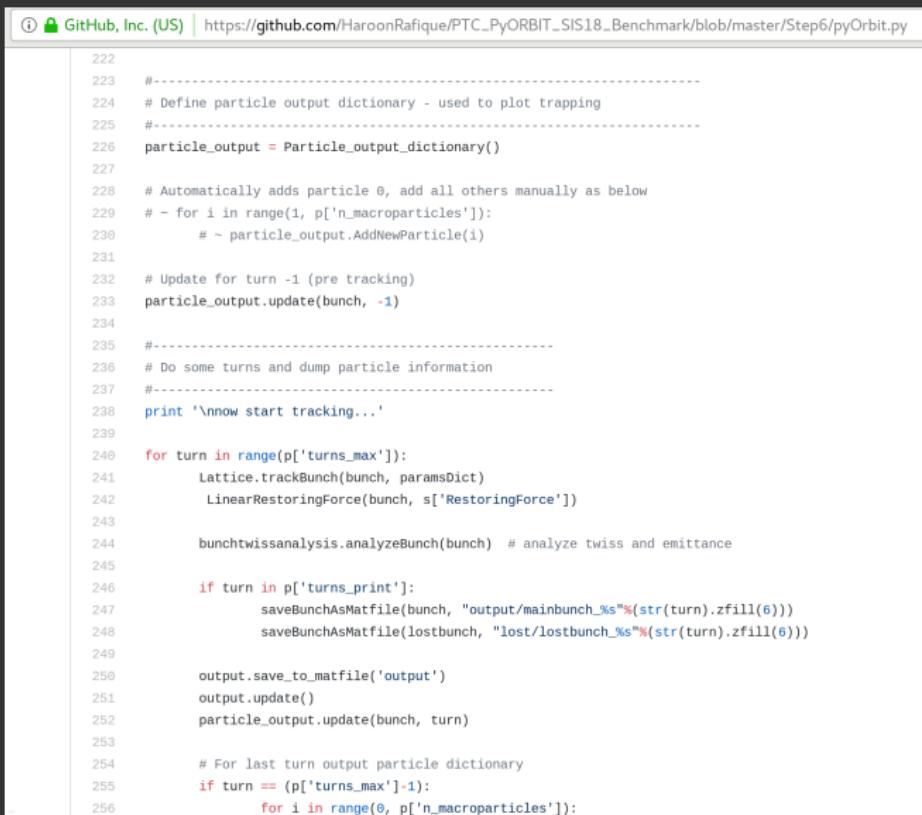
```
① GitHub, Inc. (US) | https://github.com/HaroonRafique/PTC_PyORBIT_SIS18_Benchmark/blob/master/Step6/lib/pyOrbit_LinearRestoringForce.py
 6     import math
 7     import orbit_mpi
 8
 9     from orbit_mpi import mpi_datatype, mpi_op
10     from bunch import Bunch
11
12     def LinearRestoringForce(bunch, force):
13
14         rank = 0
15         numprocs = 1
16
17         mpi_init = orbit_mpi.MPI_Initialized()
18         comm = orbit_mpi.mpi_comm.MPI_COMM_WORLD
19
20         if(mpi_init):
21             rank = orbit_mpi.MPI_Comm_rank(comm)
22             numprocs = orbit_mpi.MPI_Comm_size(comm)
23
24         nparts_arr_local = []
25         for i in range(numprocs):
26             nparts_arr_local.append(0)
27
28         nparts_arr_local[rank] = bunch.getSize()
29         data_type = mpi_datatype.MPI_INT
30         op = mpi_op.MPI_SUM
31
32         nparts_arr = orbit_mpi.MPI_Allreduce(nparts_arr_local,data_type,op,comm)
33
34         for i in range(bunch.getSize()):
35             en = bunch.dE(i)
36
37             en = en + bunch.z(i) * force
38
39             bunch.dE(i,en)
40
41     return
```

## >>> Linear Restoring Force



## >>> Particle Output Dictionary

The particle output dictionary was created to output individual particle data into a single file.

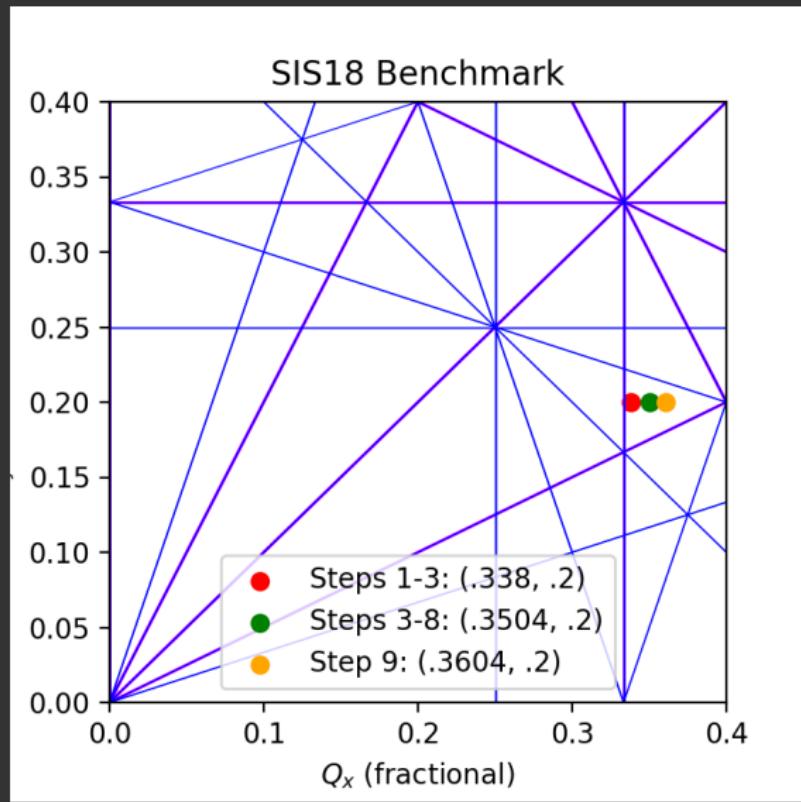


The screenshot shows a GitHub code editor with a script named 'pyOrbit.py'. The code is a Python script for generating particle output. It includes comments explaining the purpose of various sections and uses several modules like 'Lattice', 'LinearRestoringForce', and 'bunchnissanalysis'. The script defines a 'particle\_output' dictionary, adds particles, updates it with bunch information, performs tracking, and saves data to Matfiles at specific turns. The GitHub interface shows the file is 256 lines long and has been updated by HaroonRafique.

```
222 #-----
223 # Define particle output dictionary - used to plot trapping
224 #-----
225 #-----
226 particle_output = Particle_output_dictionary()
227
228 # Automatically adds particle 0, add all others manually as below
229 # - for i in range(1, p['n_macroparticles']):
230     # - particle_output.AddNewParticle(i)
231
232 # Update for turn -1 (pre tracking)
233 particle_output.update(bunch, -1)
234
235 #-----
236 # Do some turns and dump particle information
237 #-----
238 print '\nnnow start tracking...'
239
240 for turn in range(p['turns_max']):
241     Lattice.trackBunch(bunch, paramsDict)
242     LinearRestoringForce(bunch, s['RestoringForce'])
243
244     bunchnissanalysis.analyzeBunch(bunch) # analyze twiss and emittance
245
246     if turn in p['turns_print']:
247         saveBunchAsMatfile(bunch, "output/mainbunch_%s"%(str(turn).zfill(6)))
248         saveBunchAsMatfile(lostbunch, "lost/lostbunch_%s"%(str(turn).zfill(6)))
249
250     output.save_to_matfile('output')
251     output.update()
252     particle_output.update(bunch, turn)
253
254     # For last turn output particle dictionary
255     if turn == (p['turns_max']-1):
256         for i in range(0, p['n_macroparticles']):
```

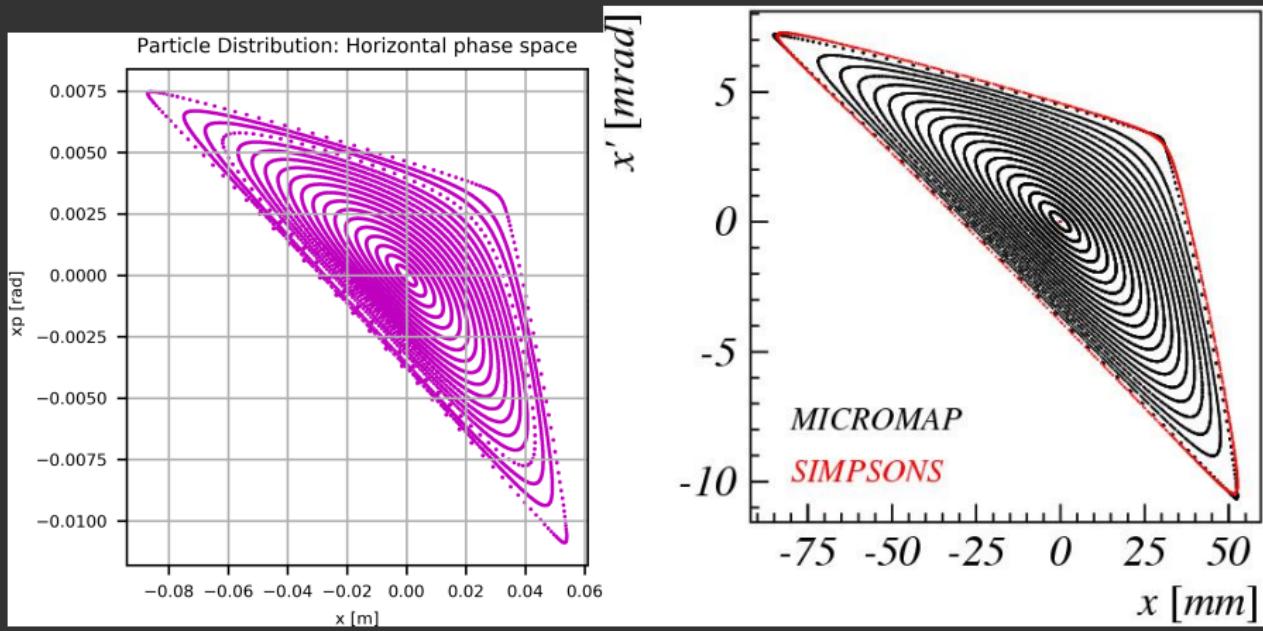
## SIS18 Benchmark

>>> Tune Diagram



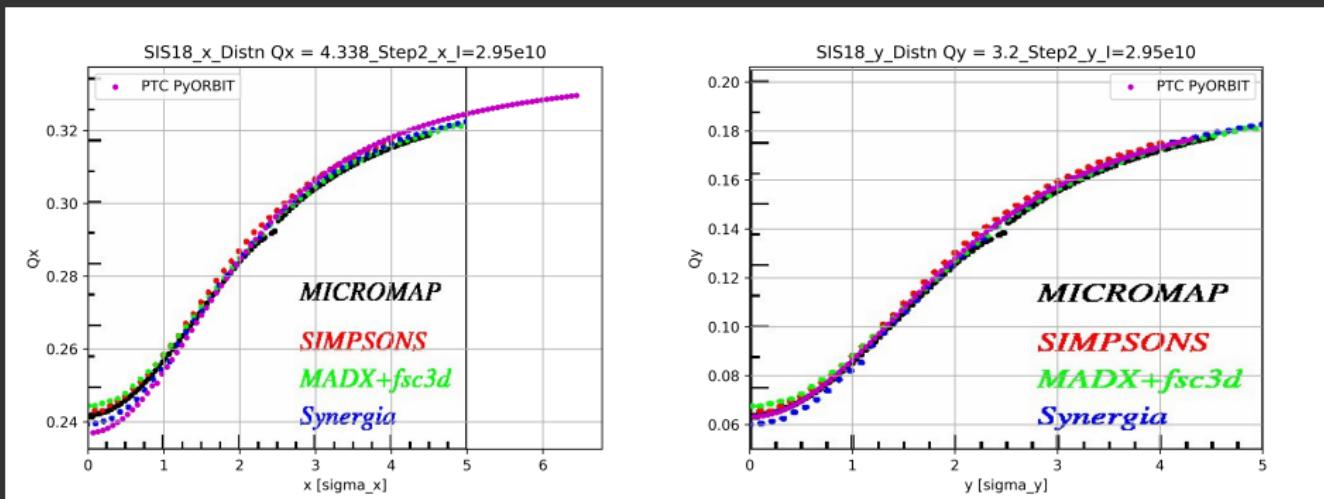
>>> Step1: Phase space  $(Q_x, Q_y) = (3.2, 4.338)$      $Q_s = \frac{1}{15000}$

This is our starting point. Sextupole on, no space charge - particles are stable up to  $6.42 \sigma_x$ .



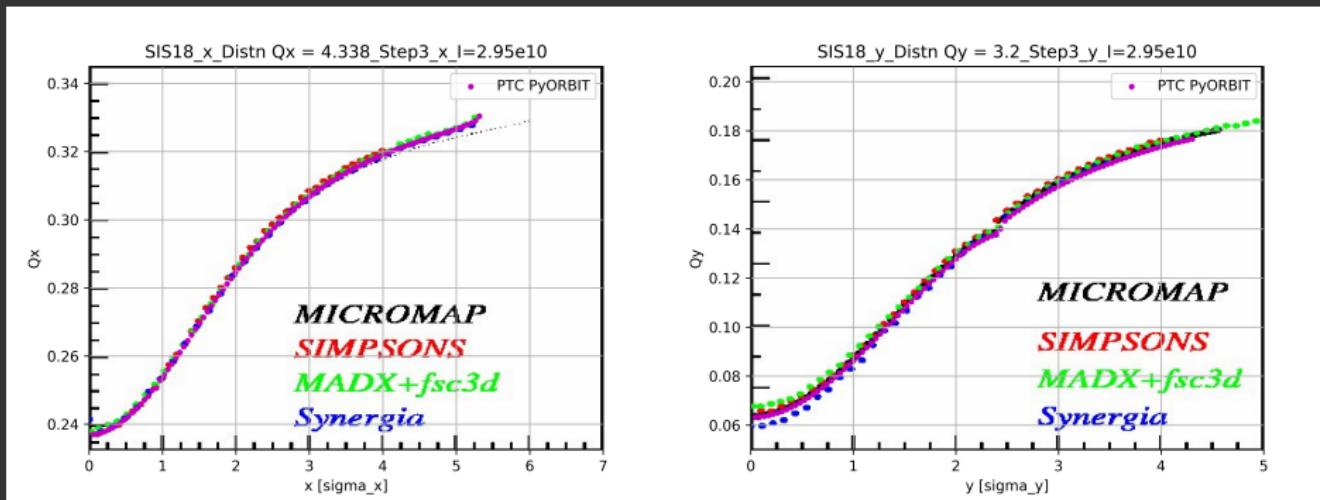
>>> Step2: Tunes with no sextupole  
 $(Q_x, Q_y) = (3.2, 4.338)$      $Q_s = \frac{1}{15000}$

This is the fractional tunespread as a function of particle amplitude  $x = \sqrt{\beta_x \gamma_x} x_0$  for the same lattice.



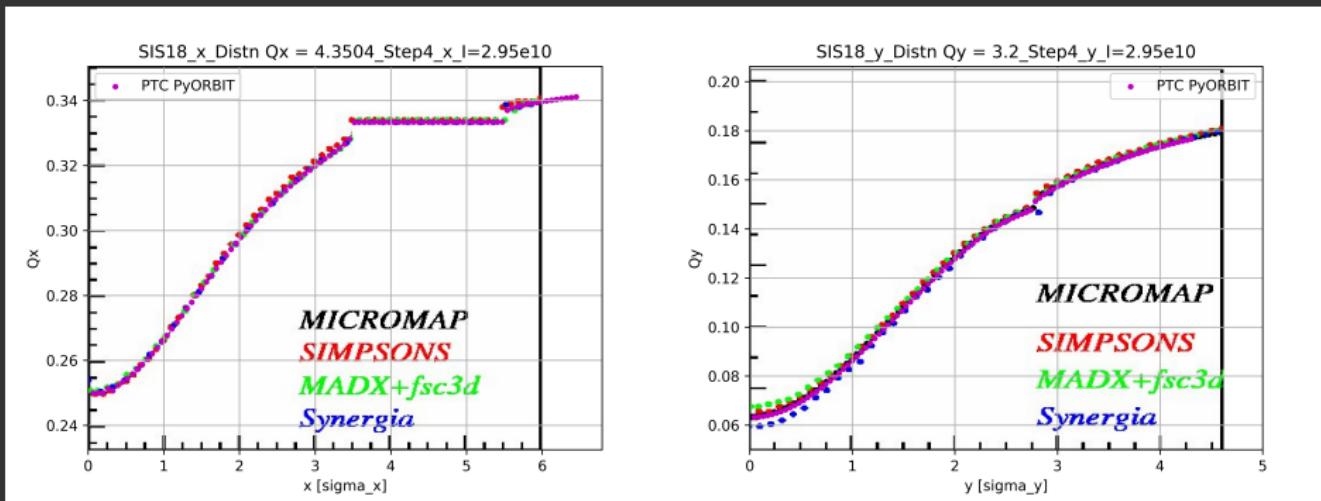
>>> Step3: Tunes with sextupole  
 $(Q_x, Q_y) = (3.2, 4.338)$      $Q_s = \frac{1}{15000}$

Now we power the sextupole, maintaining the tune. We note that the island is not visible here.

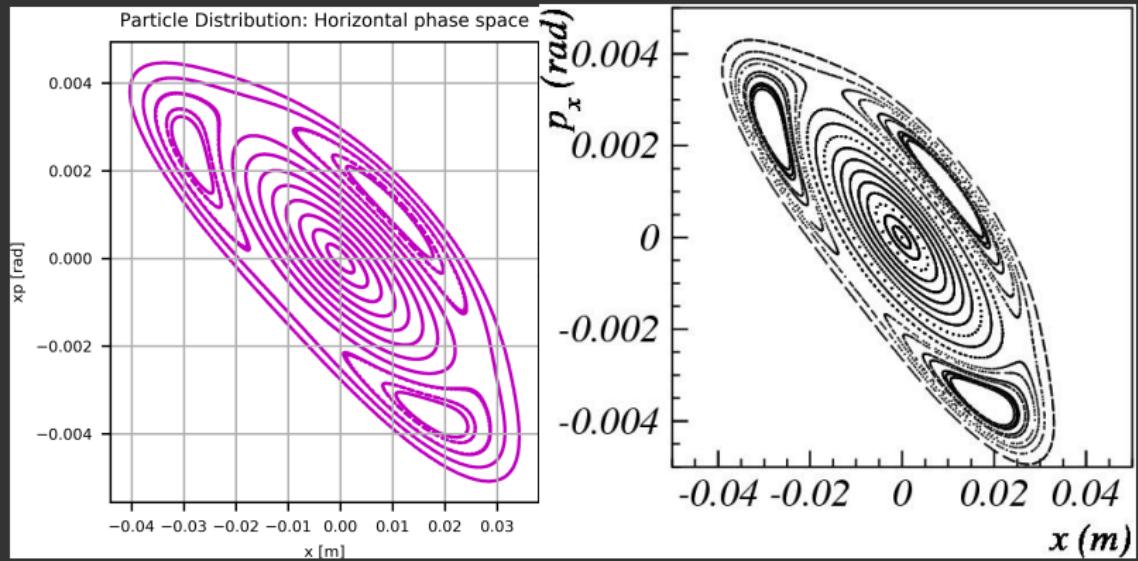


>>> Step4: Tunes with sextupole  
 $(Q_x, Q_y) = (3.2, 4.3504)$      $Q_s = \frac{1}{15000}$

We move further away from the third order resonance in order to view the islands.



>>> Step5: Phase space with sextupole  
 $(Q_x, Q_y) = (3.2, 4.3504)$      $Q_s = \frac{1}{15000}$

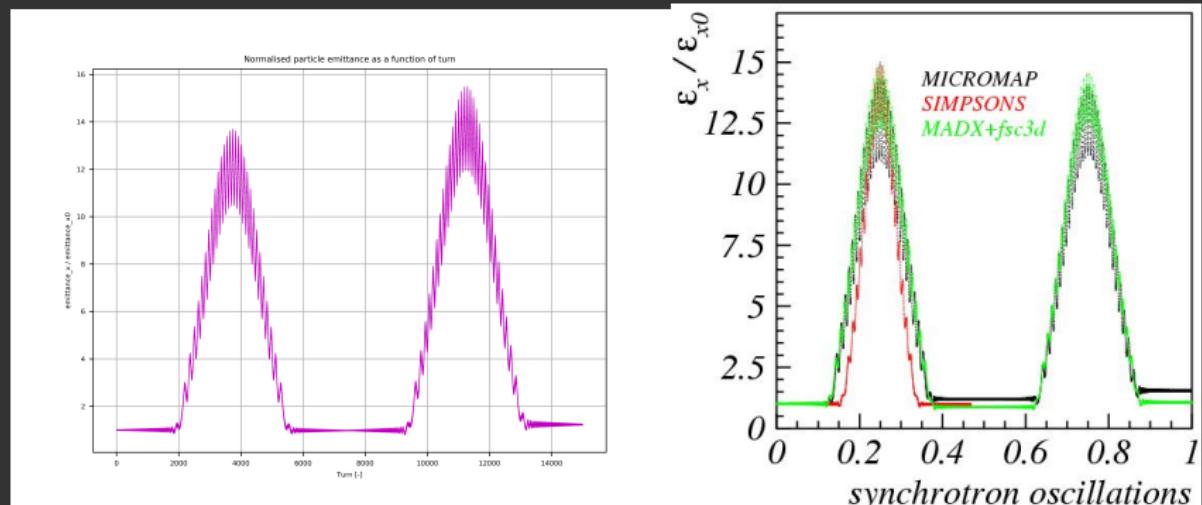


The amplitude of the island may be measured by the size of the flat region in the tune which crosses the resonance (previous slide). It is around  $2 \sigma_x$

>>> Step6: Slow trapping with sextupole

$$(Q_x, Q_y) = (3.2, 4.3504) \quad Q_s = \frac{1}{15000}$$

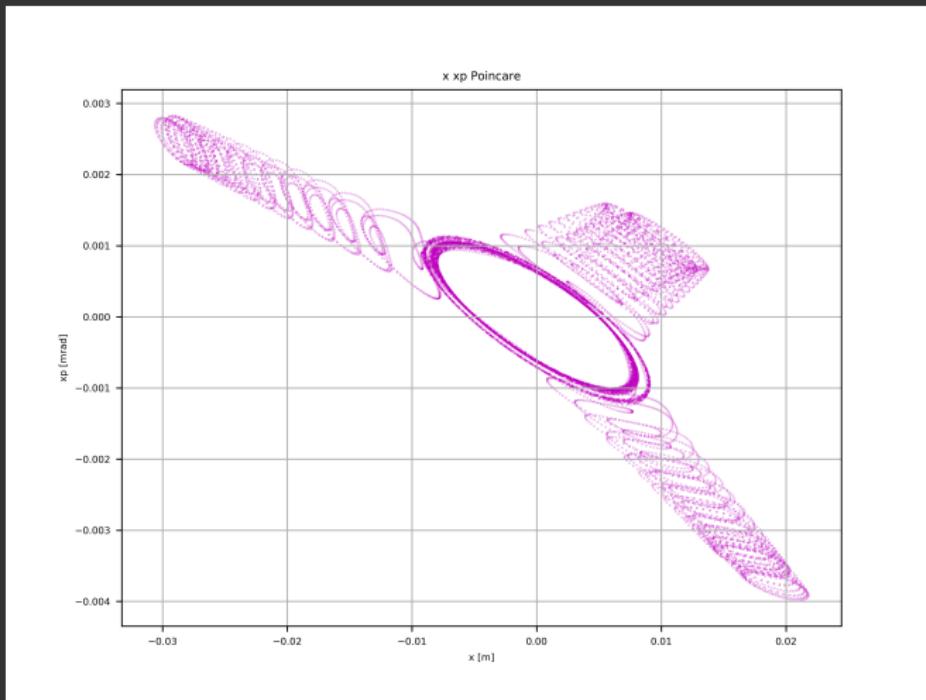
Trapping of a single particle for one synchrotron period - adiabatic crossing of the resonance.



Single particle emittance  $\epsilon_x = \beta_x p_x^2 + 2\alpha_x x p_x + \gamma_x x^2$  as a function of turn.

>>> Step6: Slow trapping with sextupole  
 $(Q_x, Q_y) = (3.2, 4.3504)$      $Q_s = \frac{1}{15000}$

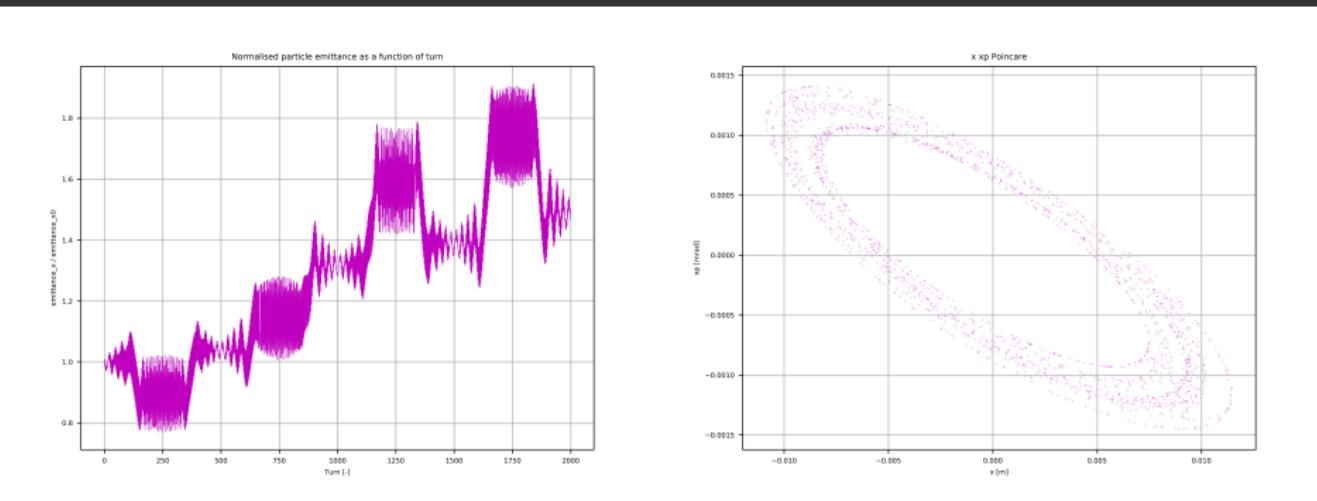
Pretty cool Poincaré section...



Single particle emittance  $\epsilon_x = \beta_x p_x^2 + 2\alpha_x x p_x + \gamma_x x^2$  as a  
function of turn

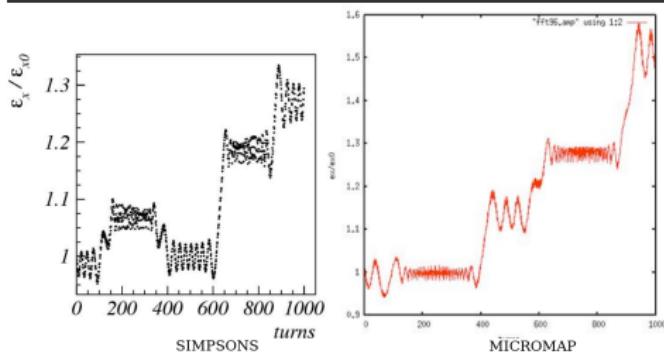
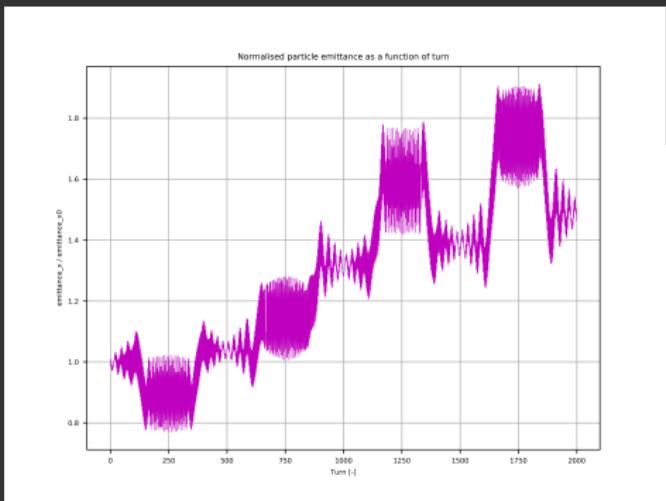
>>> Step7: Fast trapping with sextupole  
 $(Q_x, Q_y) = (3.2, 4.3504)$      $Q_s = \frac{1}{1000}$

Single particle scattering - crossing of the resonance too fast to be adiabatic.



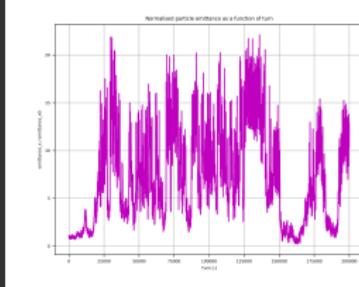
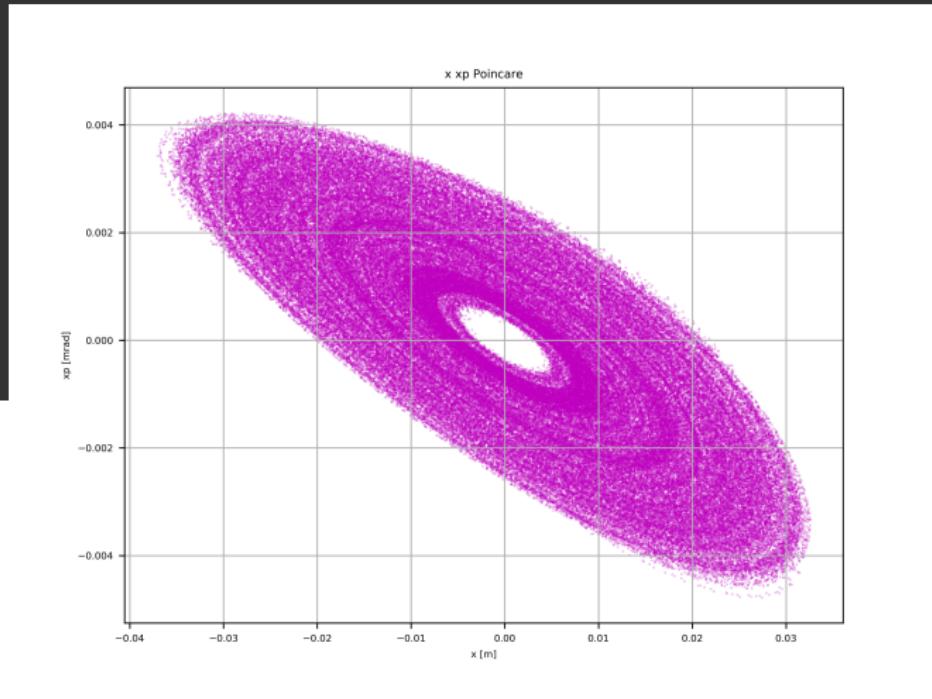
>>> Step7: Fast trapping with sextupole  
 $(Q_x, Q_y) = (3.2, 4.3504)$      $Q_s = \frac{1}{1000}$

Single particle scattering - crossing of the resonance too fast to be adiabatic.



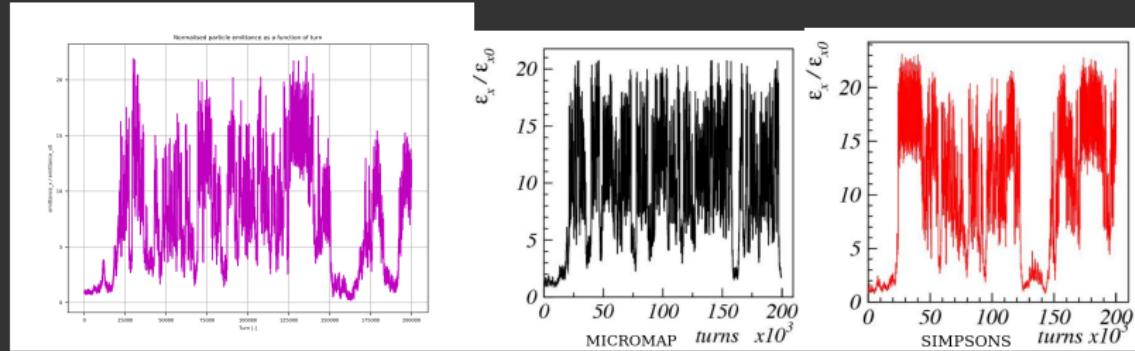
>>> Step8: Fast trapping with sextupole for many turns  
 $(Q_x, Q_y) = (3.2, 4.3504)$      $Q_s = \frac{1}{1000}$

Single particle scattering - over a longer number of synchrotron periods.



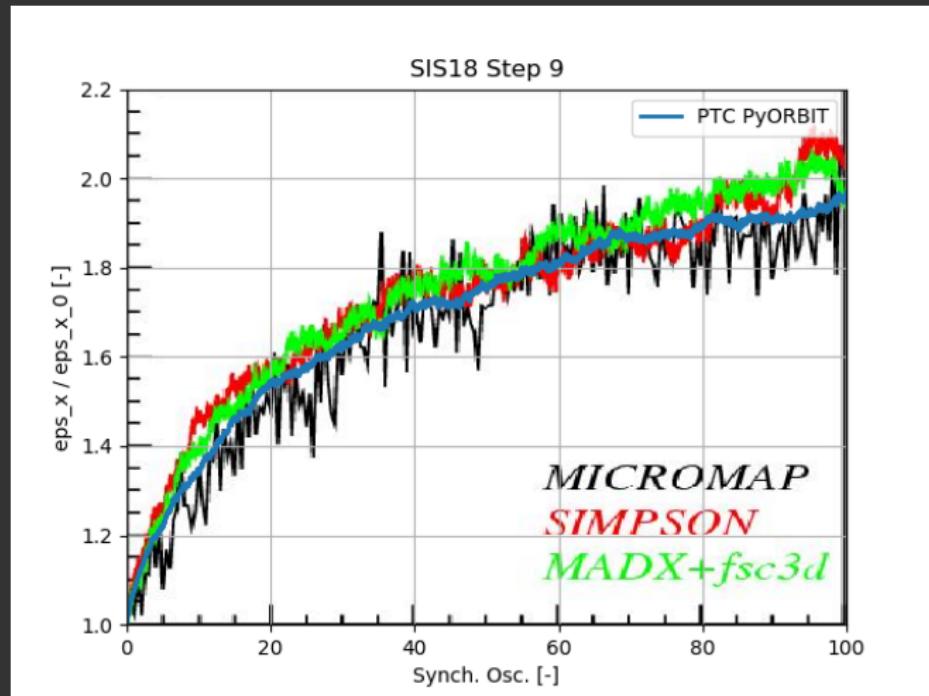
>>> Step8: Fast trapping with sextupole for many turns  
 $(Q_x, Q_y) = (3.2, 4.3504)$      $Q_s = \frac{1}{1000}$

Single particle scattering - over a longer number of synchrotron periods.



>>> Step9: Bunch emittance evolution

$$(Q_x, Q_y) = (3.2, 4.3604) \quad Q_s = \frac{1}{1000}$$



Full bunch of  $10^3$  particles - emittance growth over  $10^5$  turns.



>>> Github repository

Nearly everything required to run the benchmark in PTC-PyORBIT can be found at:

[https://github.com/HaroonRafique/PTC\\_PyORBIT\\_SIS18\\_Benchmark](https://github.com/HaroonRafique/PTC_PyORBIT_SIS18_Benchmark)  
Requirements:

- \* Run on LXPlus
- \* PyORBIT on AFS (LIU Project Space) - modification of environment variables?
- \* HTCondor scripts to run long simulations (provided)



>>> Don't Forget

Important things required for PTC-PyORBIT to function as expected:

- \* 'time.ptc' input file to make sure PTC is in the correct units
- \* initial parameters must be defined as expected and propagated consistently

## >>> Conclusions

- \* PTC-PyORBIT is in agreement with all other space charge codes that have completed this benchmark:

Person	Code	Laboratory
G. Franchetti	MICROMAP	GSI
S. Machida	SIMPSON	RAL
V. Kapin	MADX+ <a href="#">fsc3d</a>	ITEP
F. Schmidt	MADX+ORBITPTC	CERN
J. Amundson, E. Stern	Synergia	FERMILAB
J. Holmes	ORBIT	SNS

- \* It is imperative that the initial conditions are set up correctly.
- \* PTC-PyORBIT can be extended relatively easily to include new particle distributions, methods of handling output, etc.