

Space charge simulation update

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KURRI-FFAG collaboration meeting

Code development

- Space charge module of Simpsons (space charge simulation code for synchrotron)

plus

- Tracking module of S-code (FFAG tracking code).

Simpsons-FFAG?

Space charge module

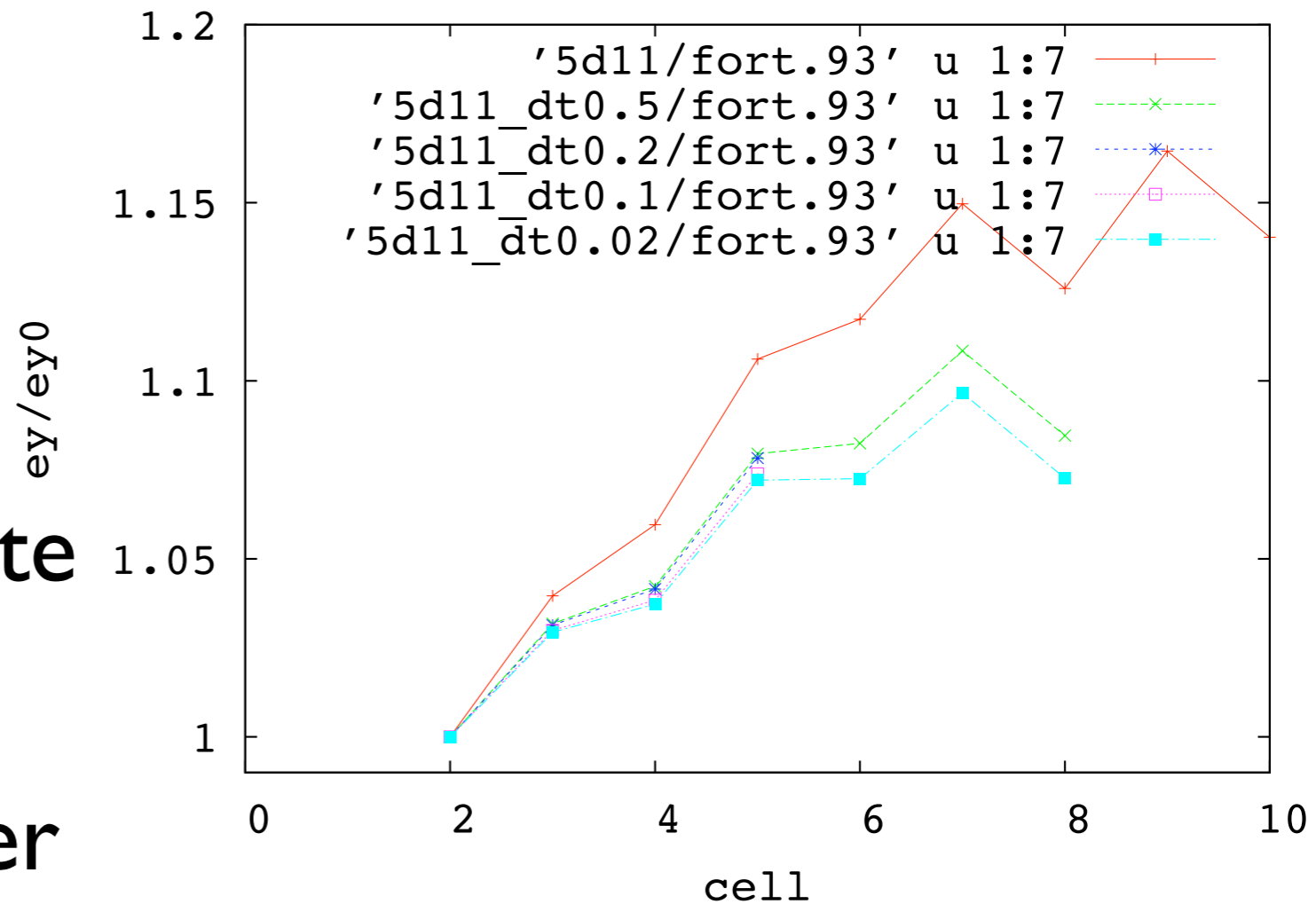
- n_s (=20) slices in longitudinal direction.
- Within a slice, transverse space charge is calculated using *r-theta* grids (cells) assuming no longitudinal variation.
- Density difference between cells in adjacent slice is used for longitudinal space charge.
- Model between 2.5-D and 3-D.

Optimization of tracking parameters (I)

- Number of space charge kicks per cell.
- Number of macro particles.
- Number of grids (cells in *r-z-theta*) for single bunch.

Optimization of tracking parameters (2)

- Number of space charge kicks per cell depends on space charge strength.
- Optimize it at moderate space charge strength.
- For simulation of higher current, increase # of kicks proportional to the strength.



red: 12 kicks per cell
green: 24 kicks
blue: 60 kicks
pink: 120 kicks
light blue: 600 kicks

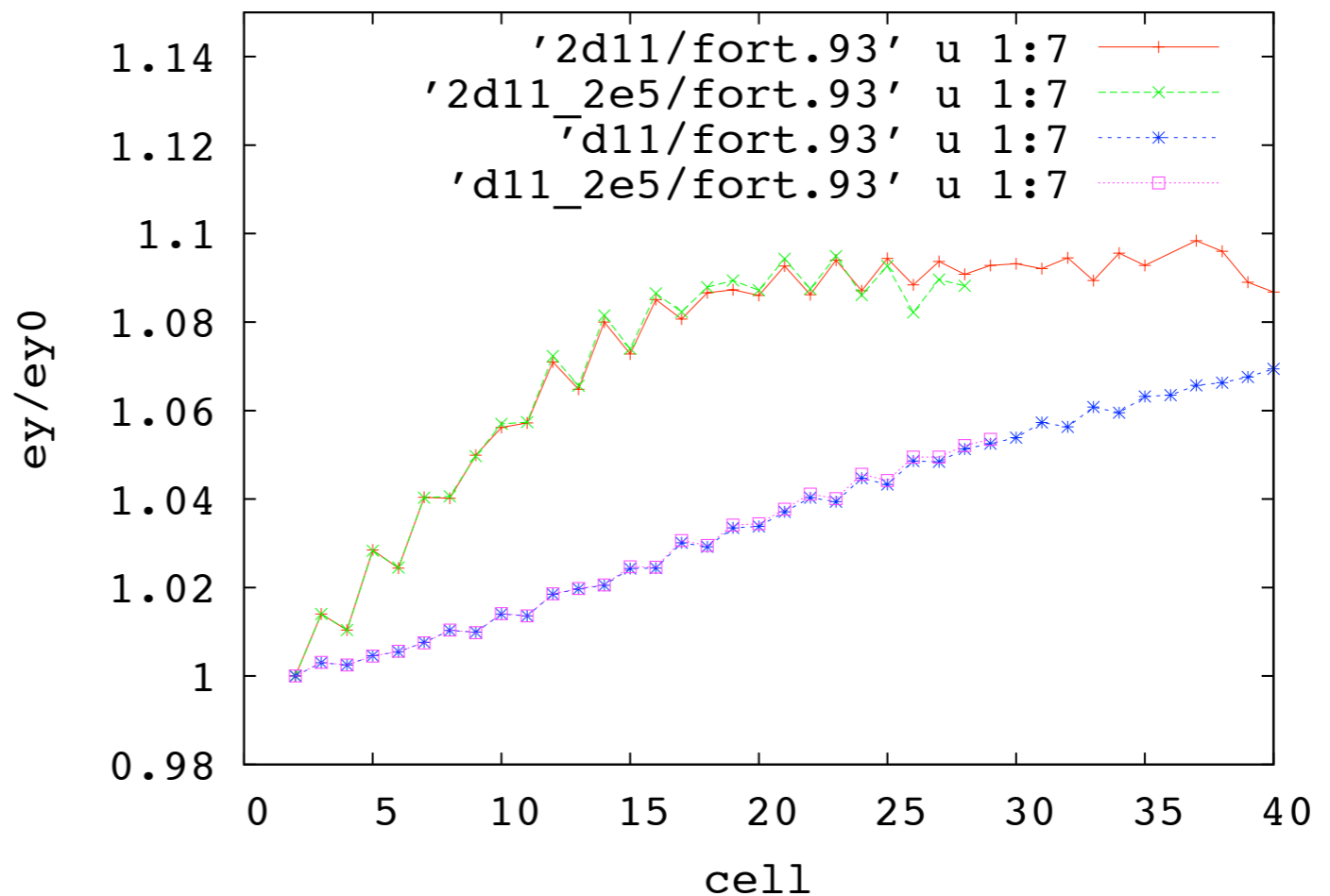
Optimization of tracking parameters (3)

- Number of macro particles depends on number of grids (cells).
- It does not depends on space charge strength.
- Once we have reasonable number of particles per cell, keep the same parameters.

Optimization of tracking parameters (4)

- Rule of thumb is 10 macro particles per cell.
 - if there are grids 20 in r and z, 16 in theta,
 - $20 (r) \times 20 (z) \times 16 (\text{theta}) \times 10 = 64,000$

Macro particles of 20,000 and 200,000 give almost identical results for two intensity cases.



Optimization of tracking parameters (5)

- Choice of number of cell is arbitrary at the moment.
- 20 (r) x 20 (z) x 16 (theta) seems to be the minimum from previous experience.

Space charge tune shift

- Gaussian distribution in 3-D (cut at 2.5 sigma), the maximum tune shift is

$$\Delta Q_y = - \frac{r_p n_t}{\pi (4 \epsilon_{rms,un}) (1 + \sqrt{\epsilon_x / \epsilon_y}) \beta^2 \gamma^3} \frac{1}{B_f}$$

- $\epsilon_{y, rms, unnor} = 1.3 \times 10^{-6}$, $n_t = 0.6 \times 10^{11}$, $B_f = 0.42$ gives
 - Effectively, $\epsilon_x \sim 10 \epsilon_y$ because of dispersion function.
 - $\Delta Q_y = -0.142$
- Tune in the model lattice is (1.816, 2.292).

Observation expected (I)

- Without error fields, only systematic resonances affect a beam.
- $Q_y=2.0$ is not a systematic integer resonance.
- However, because of 8-fold symmetry, $Q_y=2.0$ is quarter resonance ($q_y=0.250$) in each cell.
- Cell tune without tune shift is $(0.227, 0.286)$.
- $q_y=0.250$ is second order (?) parametric resonance.
(Okamoto and Yokoya, NIM A482 p.51 2002).

Observation expected (2)

- Tune shift vs intensity

intensity in a ring	cell $dq_y (q_y)$	ring $dQ_y (Q_y)$
0.6×10^{11}	-0.018 (0.268)	-0.142 (2.150)
1.8×10^{11}	-0.053 (0.233)	-0.426 (1.866)
2.4×10^{11}	-0.071 (0.215)	-0.568 (1.724)
3.6×10^{11}	-0.107 (0.179)	-0.852 (1.440)



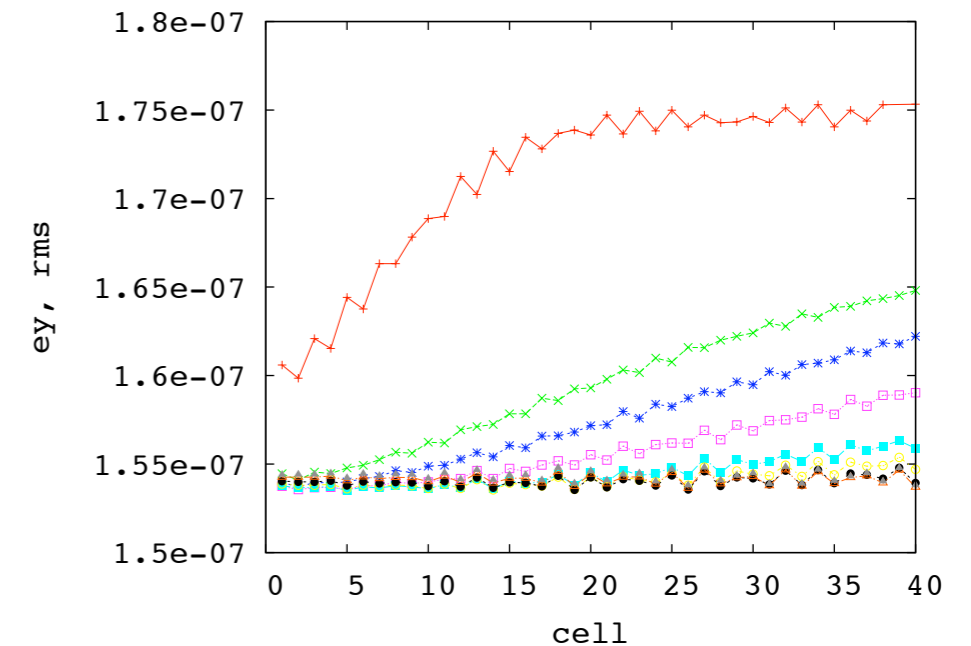
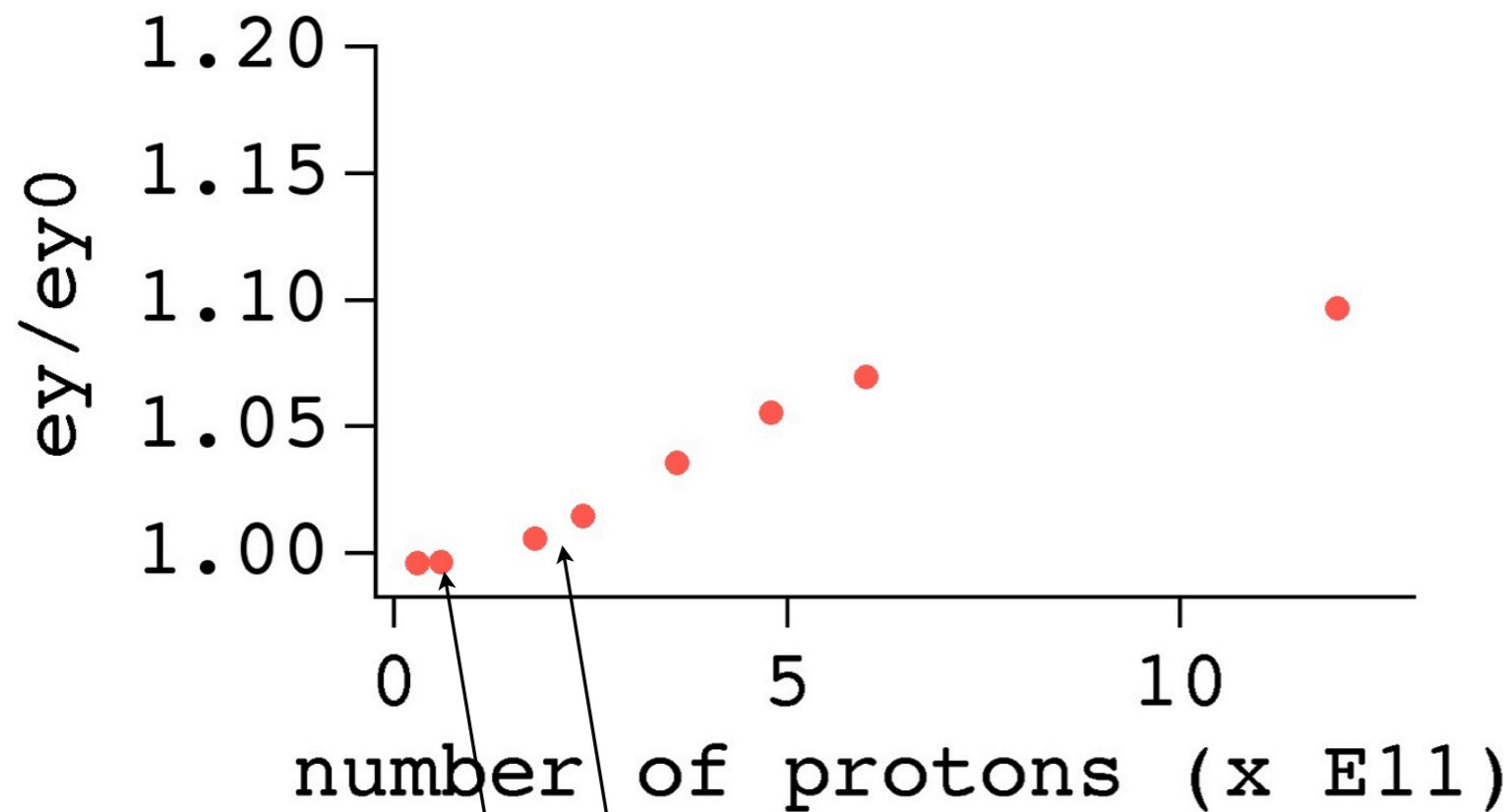
when maximum tune shift hits $q_y=0.250$



when coherent tune shift hits $q_y=0.250$.

Simulation results

- Emittance growth (5th turn) vs intensity.



when coherent tune shift hits $q_y=0.250$.

when maximum tune shift hits $q_y=0.250$

Summary

- New space charge simulation code for FFAG shows reasonable results.
- In the ideal lattice without errors, quarter resonance in each cell is the primary source of emittance growth.
- Next step
 - Multi-turn injection.
 - COD makes $Q_y=2.0$ another source of growth.
 - Comparison with growth from scattering.