Collaboration in intra-beam scattering activities for the CERN CLIC Damping Rings Dr P. Zenkevich and Dr A. Bolshakov

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## Credits

- "Upgrade of the MOCAC Code and Creation of Library Modules", A. Bolshakov, P. Zenkevich, ITEP, Moscow.
- "Kinetics Effects in Multiple Intra-beam Scattering" (IBS), P.R.Zenkevich", A.E.Bolshakov", O. Boine-Frenkenheim" - "ITEP, Moscow, Russia, "GSI, Darmstadt, Germany.
- "Last Advances in Analysis of Intra-Beam Scattering in the Hadron Storage Rings", P. Zenkevich, ITEP, Moscow, 2006.
- "A new algorithm for the kinetic analysis of intra-beam scattering in storage rings", P.R.Zenkevich\*, A.E.Bolshakov\*, O. Boine-Frenkenheim\*\*, 2005, -\*ITEP, Moscow, Russia, \*\*GSI, Darmstadt, Germany.
- "Hybrid Monte Carlo Methods for Fluid and Plasma Dynamics", R. Caflisch, Mathematics Department, Materials Science & Engineering Department, UCLA, 2007.
- "A Binary Collision Model for Plasma Simulation with a Particle Code", T. Takizuka\*, H. Abe\*\*, 1977, - \*Japan Atomic Energy Research Institute, Tokai, Ibaraki, Japan, \*\* Department of Electronics, Kyoto University, Japan.

## Proposed workplan

### • The aim of the collaboration

Investigate the issue of equilibrium beam distribution (emittance) and relaxation time in the CLIC damping rings (DR) in the presence of strong intra-beam scattering (IBS) effect, radiation damping and quantum excitation.

### • What is expected to be achieved?

- 1. Detailed description of the kinetic analysis of intra-beam scattering.
- 2. Full understanding of numerical code for IBS *MOCAC* (P. Zenkevich and A. Bolshakov) and algorithms used.
- 3. Implementation of *MOCAC* at CERN with application to the CLIC DR.
- 4. Benchmarking of *MOCAC* against Gaussian models (Bjorken-Mtingwa, Piwinski) on simplified DR lattices (single cell, cf. Y. Papaphilippou).
- 5. Upgrading of *MOCAC* code to include the radiation damping and quantum excitation effects.

# **MOnte-CArlo Code (MOCAC) 1**

### Structure of MOCAC



- Module 1: Intra-beam scattering
  - The method can be regarded as "collective map" in momentum space; input data is a set of 6-D vectors characterizing the ensemble.
- Module 5: Transformation to momentum-coordinate space
  - Transform input set of particles invariants in output set of particle 6-D vectors in momentum coordinate space.

# MOnte-CArlo Code (MOCAC) 2

- MOCAC IBS module is based on "binary collision model" (BCM): the idea is to change the real IBS by a set of artificial "scattering" events built in such a way that the mean invariants rates are same as due to real IBS process.
- *Algorithm steps (change of kernel calculation by successive application of the "binary collision" map in 3D momentum space):*
- 1. Choose phases randomly for a given set of macro-particles invariants.
- 2. Calculate the momenta and coordinates of macro-particles.
- 3. Work out the macro-particles distribution on the space cells and link test particles to a "partner" in the cell.
- 4. Apply binary collision map on each macro-particle pair (test & "partner") using "local ensemble" for each cell: i.e. calculate the collision angles for given "collision events" and update momentum components for the test particle and its partner.
- 5. Calculate the new set of macro-particles invariants.

## MOnte-CArlo Code (MOCAC) 3

- This operation can be considered as "collective map"; each particle is "test" and "field" particle simultaneously.
- The map is repeated through the time interval  $\Delta t$ .
- The lattice is described as set of discrete points related to different longitudinal coordinates. Each point is characterized by its set of optical parameters.
- This algorithm is many-dimensional integration by Monte-Carlo method and needs high number of macro-particles (10<sup>5</sup>-10<sup>6</sup>) to diminish a noise.

• Compute the scattering angle  $\Theta_{sc}^{i,j}$  for two colliding macro-particles using the expression ( $\rho_0$  = particle density, u =momentum, log  $\Lambda$  = Coulomb logarithm).

$$\sin\left(\frac{\Theta_{sc}^{i,j}}{2}\right) = \sqrt{\frac{2A\rho_0 \log \Lambda^{i,j} \Delta t}{N_{mp} \left|\vec{u}^i - \vec{u}^j\right|^{3/2}}}$$

- Compute the momentum change of each interacting particle (test & partner).
- Azimuthal angle  $\Phi_{sc}^{ij}$  is defined by random choice on interval [0,  $2\pi$ ].
- This map was included in *MOCAC*. This idea was suggested earlier by T. Takizuka & H. Abe (1977). At each step random choice of partner are assumed.



Coordinates in the laboratory frame and the relative momentum u at the time t.





Change in the relative momentum frame after collision at the time  $t+\Delta t$ .

## **Collision Map 2**

Integration step [s]	$\Delta t \approx T_{\rm IBS} / 100 N_{\rm per}$
Number of macro-particles	$N_{\rm mp} \ge N_{\rm long} N_{\rm tr}$
Number of points per period	$N_{\rm per} \times N_{\rm mag}/5$
Size of transverse cell [m]	$\Delta_{tr} \leq \min(\sigma_{x,y}/5, D\sigma_p)$
Size of longitudinal cell [m]	$\Delta_{long} \geq \sigma_s / 10$
Maximal collision angle [rad]	$\chi_{\rm max} < 0.5$

List of code parameters

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## **Collision Map 3**

- T. Takizuka & H. Abe, J. Comp. Phys. 25 (1977).
- T & A binary collision model is equivalent to the collision term in Landau-Fokker-Planck equation
  - The scattering angle  $\Theta$  is chosen randomly from a Gaussian random variable  $\delta \equiv \tan(\Theta/2)$ .
  - $\delta$  has mean 0 and variance  $\langle \delta \rangle^2 = (n_{\rm L} e^4 \log \Lambda / 8\pi \varepsilon_0 m^2 u^3) \Delta t$ .
  - Parameters: log A = Coulomb logarithm u = relative velocity.
- Simulation
  - Every particle collides once in each time interval: scattering angle depends on  $\Delta t$ .
  - Implemented in ICEPIC by Birdsall, Cohen and Procassini.