

Intra-Beam scattering studies for CLIC damping rings

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Thanks to : M. Martini, Y. Papaphilippou

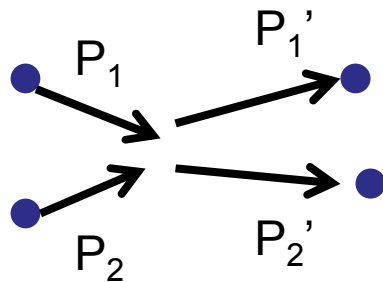
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CONTENTS

- Introduction
- Status of IBS calculations
- MOCAC code
- Conclusions

Intra-Beam Scattering

IBS is the effect due to multiple Coulomb scattering between charged particles in the beam



- Only Particles at the same position interact
- Only the momenta of the particles change

All the scattering events produce a variation of the particle distribution of the beam:

$$\rho(\mathbf{r}, \mathbf{p}, t), \quad N_b = \int \rho(\mathbf{r}, \mathbf{p}, t) d^3 \mathbf{p} d^3 \mathbf{r}$$

Considering only IBS, the evolution of the distribution function is given by:

$$\frac{\partial \rho}{\partial t}(\mathbf{r}, \mathbf{p}, t) \stackrel{IBS}{=} \int \left(\frac{\partial N}{\partial t}(\mathbf{r}, t, \mathbf{p}' \rightarrow \mathbf{p}) - \frac{\partial N}{\partial t}(\mathbf{r}, t, \mathbf{p} \rightarrow \mathbf{p}') \right) d^3 \mathbf{p}'$$

Also mean values of functions of the particle coordinates change:

$$\epsilon_k = \beta_k r_k'^2 + 2\alpha_k r_k r_k' + \gamma_k r_k^2$$

$$\epsilon_{k,RMS}(t) = \frac{\langle \epsilon_k \rangle_t}{2} = \frac{1}{2} \int \epsilon_k(r_k, r_k') \rho(\mathbf{r}, \mathbf{r}', t) d^3 \mathbf{r}' d^3 \mathbf{r}$$

It would be possible to follow the evolution of the beam parameters if the evolution of the distribution $\rho(\mathbf{r}, \mathbf{p}, t)$ was known.

The effect of the IBS on the beam emittance is evaluated by means of the growth rates:

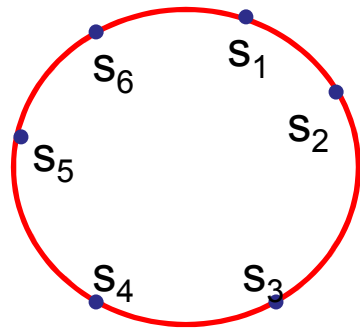
$$\frac{1}{\tau_{k,IBS}} = \frac{1}{\epsilon_{k,RMS}} \frac{d \langle \epsilon_k \rangle}{dt} \quad k = x, z, s$$

Growth rates are calculated at specific points on the lattice.

The Bjorken-Mtingwa formulation gives, for the growth rates:

$$\frac{1}{\tau_k} = \frac{r_0^2 c N(\log)}{8\pi\gamma^4 \beta^3 \epsilon_x \epsilon_z \epsilon_s} \int_0^\infty \frac{\lambda^{1/2} d\lambda}{|L + \lambda I|^{1/2}} \left\{ \text{Tr} L^{(k)} \text{Tr}(L + \lambda I)^{-1} - 3 \text{Tr} L^{(k)} (L + \lambda I)^{-1} \right\}$$

Growth rates are calculated at different points of the lattice and then averaged over the ring:



$$\frac{1}{\tau} = \sum_{k=1}^N \frac{S_{k+1} - S_k}{S} \frac{1}{\tau_k} \quad S = \sum_{k=1}^N S_k$$

Considering only the contribution of IBS, the evolution of the emittance is given by :

$$\frac{d\epsilon_{k,RMS}}{dt} = \frac{\epsilon_{k,RMS}}{\tau_{k,IBS}}$$

Taking into consideration also radiation damping and quantum excitation, the emittance evolves according to the equation:

$$\frac{d\epsilon_{k,RMS}}{dt} = -\frac{2}{\tau_k}(\epsilon_{k,RMS} - \epsilon_{k,0,RMS}) + \frac{2\epsilon_{k,RMS}}{\tau_{k,IBS}}$$

For the CLIC damping rings the parameters are:

- Damping time = 1.5 , 1.5 , 0.8 ms
- IBS Growth time = 1.9, 6, 3.9 ms
- Horizontal emittance (No IBS) = 89 nm
- Horizontal emittance (IBS) = 461 nm
- Vertical emittance (No IBS) = 3.0 nm
- Vertical emittance (IBS) = 4.4 nm
- Longitudinal emittance (No IBS) = 2256 eV s
- Longitudinal emittance (IBS) = 3460 eV s

In classical IBS theory $\tau_{k,IBS}$ is calculated assuming:

Bjorken-Mtingwa:

- Coulomb Scattering transition rate from QED
- Gaussian Distribution

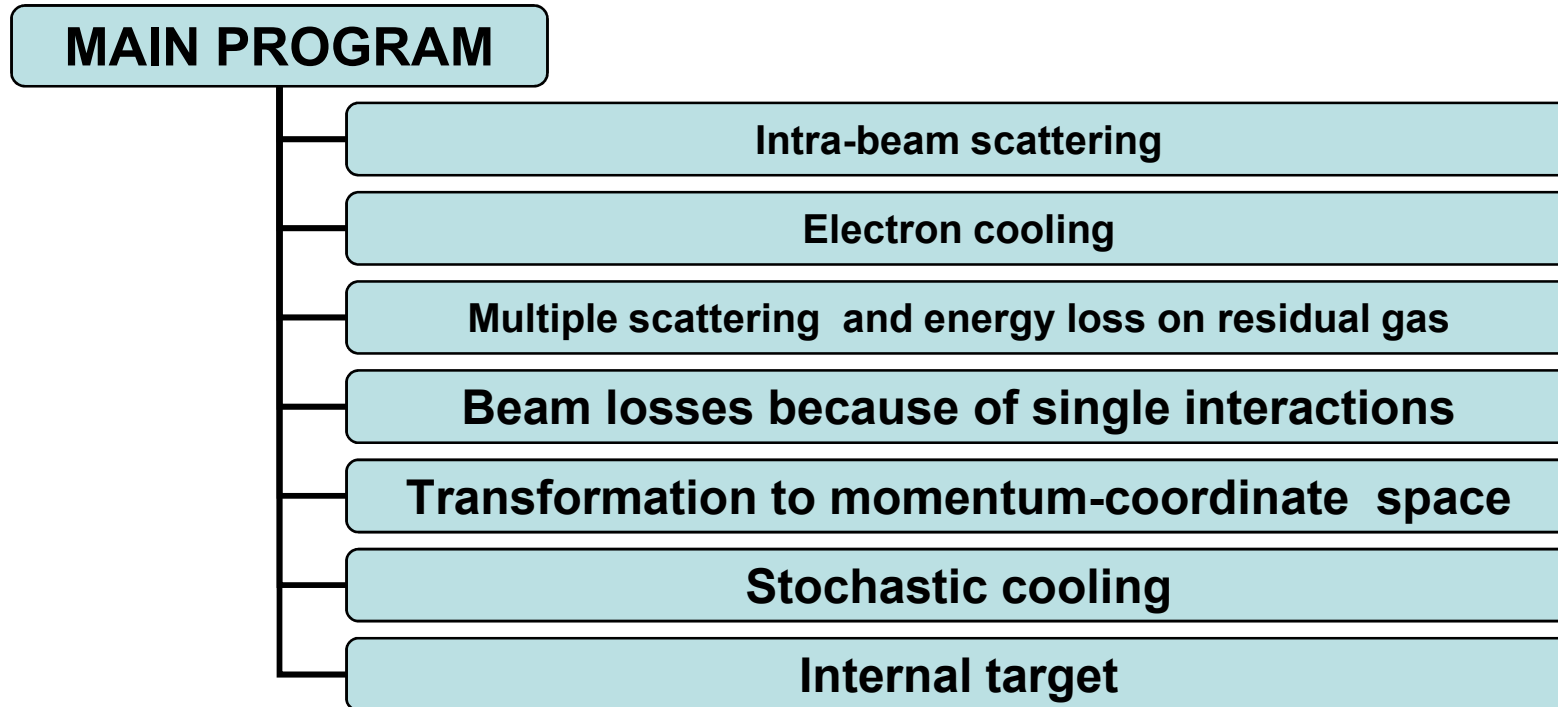
Piwinsky-Martini:

- Rutherford cross section
- Gaussian Distribution

- How do $\tau_{k,IBS}$ change if the distribution is not Gaussian ?
- Is it possible to follow the evolution of the particle distribution in the DR ?

MOCAC Code

Developed by P. R. Zenkevich and A. E. Bolshakov, *ITEP, Moscow, Russia*



- Not dedicated to IBS study.
- Rather complicated structure
- Not modified since 2006

It needs to be elaborated

MOCAC Code: test

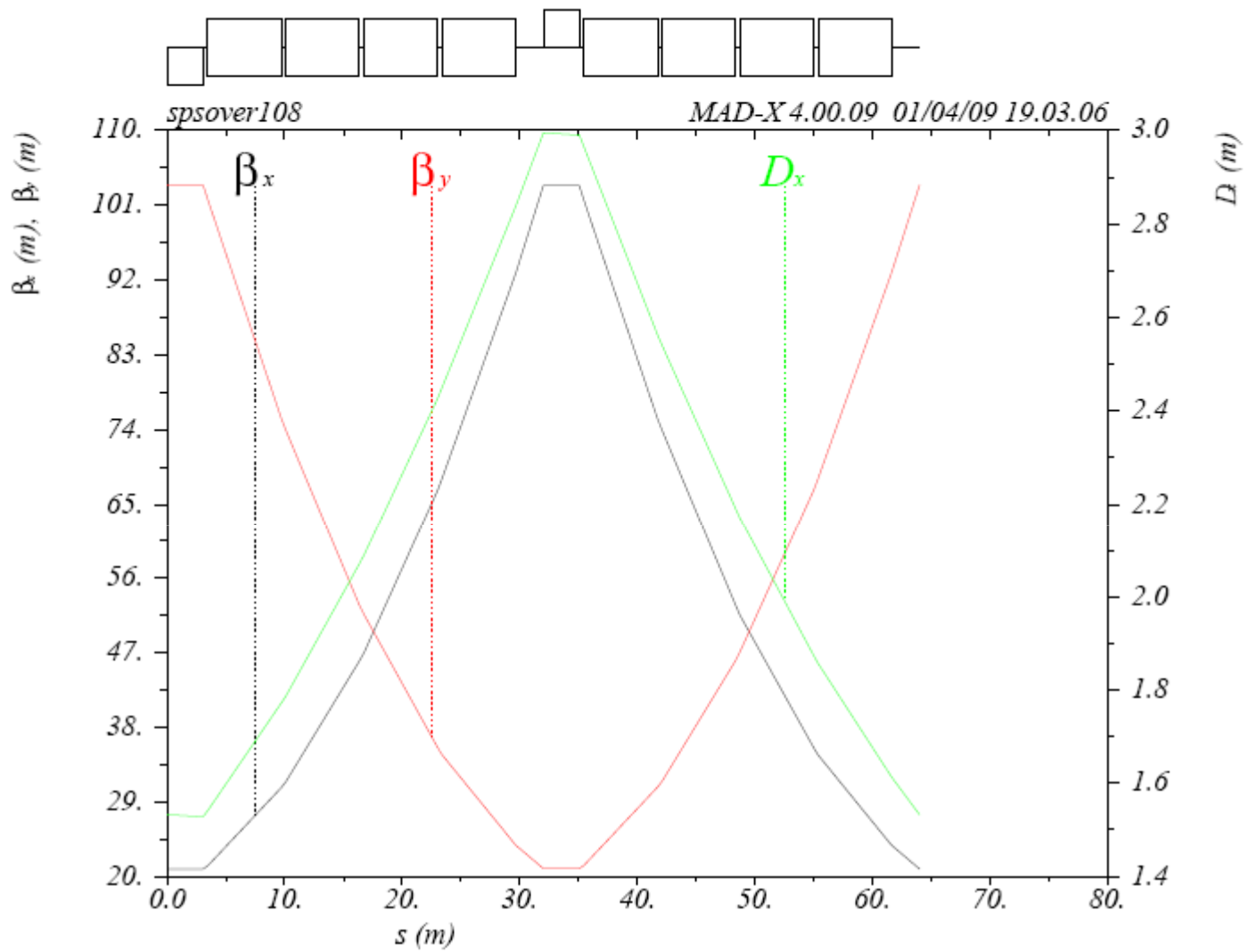
Due to the complicated structure of the code it was not easy to fit the parameters of the program for the test.

29-janv-09

SPS 1/108 lattice: $p=26$ GeV/c, $N=1E14$ (1 bunch), $eHn=eVn=0.5$ um, $lbunch=1$ ns, $dpp=2.5E-3$

	1/TL [s-1]	1/TX [s-1]	1/TY [s-1]	Cb(Log)
MADX: BJORKEN-MTINGWA Coulomb log computed	2,91E-03	1,22E-01	1,51E-02	15,742
Mathematica: BJORKEN-MTINGWA Coulomb log computed	3,05E-03	1,30E-01	1,65E-02	15,747
Mathematica: BJORKEN-MTINGWA Coulomb log=20	3,87E-03	1,65E-01	2,09E-02	20,000
Mathematica: PIWINSKI	3,81E-03	1,62E-01	2,28E-02	
MOCAC	-2,70E+00	2,88E+00	5,06E-01	

A more accurate comparison will be possible after cleaning of the code.



MOCAC Code: IBS routine

MOCAC 'tracks' particles through their invariants.

Transversal invariants:

$$\epsilon_x(i) = \beta_x \left(x'_i - D' \frac{\Delta p_i}{p} \right)^2 + 2\alpha_x \left(x'_i - D' \frac{\Delta p_i}{p} \right) \left(x_i - D \frac{\Delta p_i}{p} \right) + \gamma_x \left(x_i - D \frac{\Delta p_i}{p} \right)^2$$

$$\epsilon_z(i) = \beta_z z'_i{}^2 + 2\alpha_z z_i z'_i + \gamma_z z_i^2$$

Longitudinal invariant:

$$\epsilon_s(i) = \left(\frac{\Delta p_i}{p} \right)^2 + \frac{(2\pi)^2 \nu_s^2}{\left(\alpha - \frac{1}{\gamma^2} \right)^2 C^2} \Delta s_i^2 \quad i = 1, \dots, Num.Part.$$

Phases of the particles are not tracked.

MOCAC Code: IBS routine

The initial distribution of the beam is a randomly generated Gaussian*:

Only the invariants of the particles are generated according to the probability density distribution:

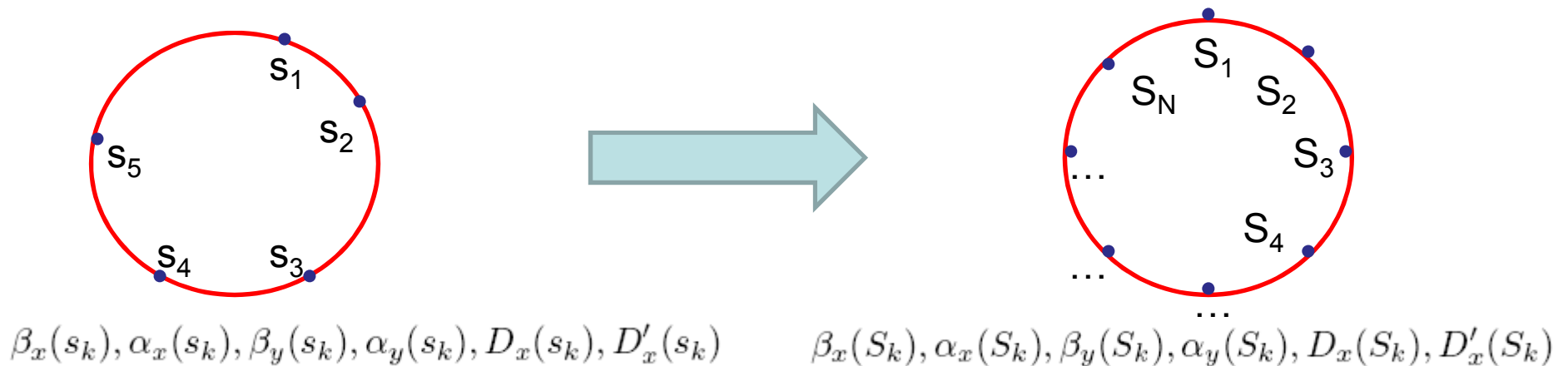
$$p(\epsilon_k) = \frac{1}{\langle \epsilon_k \rangle} e^{-\frac{\epsilon_k}{\langle \epsilon_k \rangle}} \quad \langle \epsilon_k \rangle = 2\epsilon_{k,RMS}$$

The initial emittances of the beam are given in the input file.

Taking uniformly distributed phases assures a Gaussian distribution of the particles.

MOCAC Code: IBS routine

- Input parameters are read from an input file (ASCII)
- Optical functions of the lattice are read from a MAD 8 output file
- A cubic interpolation of the optical functions is performed to calculate the values of the functions at N equally spaced points of the lattice (total length of the lattice is kept constant)



MOCAC Code: IBS routine

The total length of the simulation is $N_{INJ} \times T_{INJ}$

T_{INJ} is the injection period (not used for IBS)

Each period T_{INJ} is divided in N_{IBS} sub-periods of length $T_{IBS} = T_{INJ} / N_{IBS}$

In each sub-period T_{IBS} the scattering routine is run at each of the equally spaced N points of the ring.

The scattering routine is called at intervals $\Delta T = T_{IBS} / N$

After each call of the scattering routine the invariants of every particle are changed according to the momentum change of the particles.

MOCAC Code: Scattering routine

The coordinates and momenta of the particles are generated.

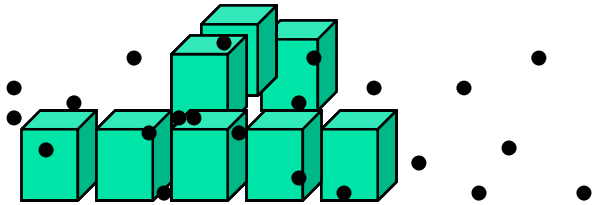
ϕ_x ϕ_z ϕ_s are randomly generated uniformly in $[0, 2\pi]$ for each particle.

$$\begin{aligned}\frac{\Delta p_i}{p} &= \sqrt{\epsilon_s(i)} \cos(\phi_s) & \Delta s_i &= \frac{(\alpha - \frac{1}{\gamma^2})C}{2\pi\nu_s} \sqrt{\epsilon_s(i)} \sin(\phi_s) \\ x'_i &= -\frac{\sqrt{\epsilon_x(i)}}{\sqrt{\beta_x}} [\alpha_x \cos(\phi_x) + \sin(\phi_x)] + D'_x \frac{\Delta p_i}{p} & x_i &= \sqrt{\epsilon_x(i)\beta_x} \cos(\phi_x) + D_x \frac{\Delta p_i}{p} \\ z'_i &= -\frac{\sqrt{\epsilon_z(i)}}{\sqrt{\beta_z}} [\alpha_z \cos(\phi_z) + \sin(\phi_z)] & z_i &= \sqrt{\epsilon_z(i)\beta_z} \cos(\phi_z)\end{aligned}$$

The scattering routine implements the Binary Collision Model
(*T. TAKIZUKA and H. ABE, Journal of Computational Physics 25, 1977*)

MOCAC Code: Scattering routine (BCM)

Particles are grouped in geometric cells.

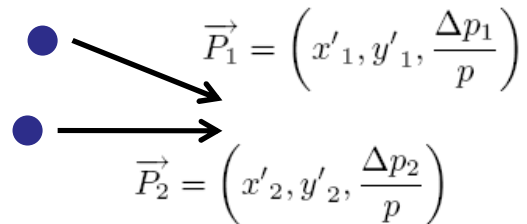


The density inside each cell is considered uniform $\rho = N_p/V$
Particles in the same cell are considered at the same point in space and then can scatter.

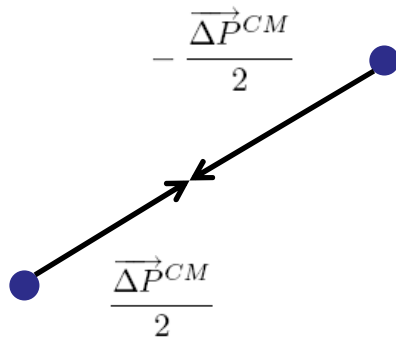
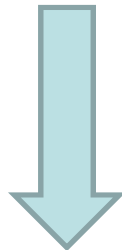
The particles inside each cell are randomly coupled and a 'collision' between the particles of each couple is simulated.

MOCAC Code: Collision simulation

The collision is simulated in the center of mass frame, which is assumed to be the beam rest frame, for any couple of particles.



$$\overline{\Delta \vec{P}} = \left(x'_2 - x'_1, y'_2 - y'_1, \frac{\Delta p_2}{p} - \frac{\Delta p_1}{p}\right)$$



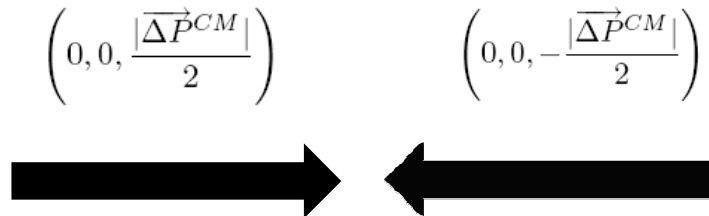
$$\overline{\Delta \vec{P}}^{CM} = \left(x'_2 - x'_1, y'_2 - y'_1, \frac{1}{\gamma} \left(\frac{\Delta p_2}{p} - \frac{\Delta p_1}{p}\right)\right)$$

MOCAC Code: Collision simulation

Operating the following rotation of the beam frame:

$$R = \begin{pmatrix} \frac{\Delta P_x^{CM} \Delta P_s^{CM}}{|\Delta \vec{P}^{CM}| |\Delta \vec{P}_\perp^{CM}|} & \frac{\Delta P_z^{CM} \Delta P_s^{CM}}{|\Delta \vec{P}^{CM}| |\Delta \vec{P}_\perp^{CM}|} & -\frac{|\Delta \vec{P}_\perp^{CM}|}{|\Delta \vec{P}^{CM}|} \\ -\frac{\Delta P_z^{CM}}{|\Delta \vec{P}_\perp^{CM}|} & \frac{\Delta P_x^{CM}}{|\Delta \vec{P}_\perp^{CM}|} & 0 \\ \frac{\Delta P_x^{CM}}{|\Delta \vec{P}^{CM}|} & \frac{\Delta P_z^{CM}}{|\Delta \vec{P}^{CM}|} & \frac{\Delta P_s^{CM}}{|\Delta \vec{P}^{CM}|} \end{pmatrix}$$

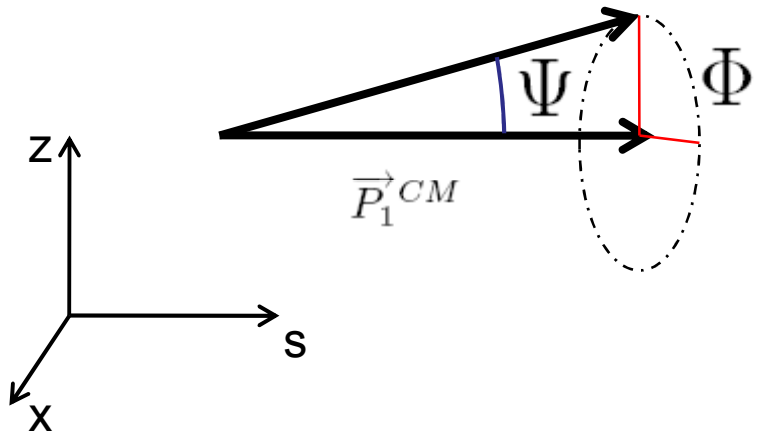
we are lead to the simplest case:



MOCAC Code: Collision simulation

The momentum change can be now expressed by the angles Ψ and Φ

$$\vec{P}_1^{CM} = \vec{P}_1^{CM} + \overline{\Delta P}_1^{CM}$$



Φ is uniformly distributed in $[0, 2\pi]$

$$\overline{\Delta P}_1^{CM} = \left(\frac{|\overline{\Delta P}^{CM}|}{2} \cos(\Phi) \sin(\Psi), \frac{|\overline{\Delta P}^{CM}|}{2} \sin(\Phi) \sin(\Psi), \frac{|\overline{\Delta P}^{CM}|}{2} (\cos(\Psi) - 1) \right)$$

MOCAC Code: Collision simulation

The scattering angle is determined by:

$$\Psi = 2r_i \sqrt{\frac{8\pi c\rho L_c \Delta T}{\gamma^5 \beta^3 |\overrightarrow{\Delta P}^{CM}|^3}} \quad L_c = \log \left(\frac{\Delta z \beta^2 \gamma^2 |\overrightarrow{\Delta P}^{CM}|^2}{2r_i} \right)$$

There is a discrepancy between this formula and *P. Zenkevich, NIM A, 577 (2007)*:

$$\sin \left(\frac{\Psi}{2} \right) = r_i \sqrt{\frac{4\pi c\rho L_c \Delta T}{\gamma^5 \beta^3 N |\overrightarrow{\Delta P}^{CM}|^{3/2}}}$$

To be clarified by Zenkevich.

MOCAC Code: Collision simulation

Averaging the momentum change over all the particles we should get:

$$\vec{F} = \frac{d \langle \Delta \vec{P}_1 \rangle}{dt} = -\frac{2\pi c r_i^2}{\gamma^5 \beta^3} \left\langle L_c \frac{\vec{P}_1 - \vec{P}_2}{|\vec{P}_1 - \vec{P}_2|^3} \right\rangle$$

\vec{F} is the friction force in the Fokker-Planck equation:

$$\frac{\partial \rho}{\partial t} = -\frac{\partial}{\partial P_m} (F_m \rho) + \frac{1}{2} \frac{\partial^2}{\partial P_m \partial P_{m'}} (D_{m,m'} \rho)$$

MOCAC Code: Collision simulation

By transforming back $\overrightarrow{\Delta P_1}^{CM}$ to the laboratory frame it is possible to compute:

$$\overrightarrow{P_1'}^{LAB} = \overrightarrow{P_1}^{LAB} + \overrightarrow{\Delta P_1}^{LAB} \quad \overrightarrow{P_2'}^{LAB} = \overrightarrow{P_2}^{LAB} + \overrightarrow{\Delta P_2}^{LAB}$$

Using the momentum conservation: $\overrightarrow{\Delta P_2}^{LAB} = -\overrightarrow{\Delta P_1}^{LAB}$

The new invariants for the 2 particles are then calculated.

MOCAC Code: Possible changes

- Complete the cleaning of the code and run new tests (in progress)
- Read the optical function from a MADX output file (short time)
- Tracking the particles using the phase advances (not an issue)
- Replace the scattering dynamics with a dynamics compatible with the classical IBS theory (been examined, feasible-medium time)
- Implement the damping effect in the code
(under investigation by Zenkevich and Bolshakov – will be needed after IBS investigations completed)

Conclusions

MOCAC contains good ideas to investigate the evolution of the beam distribution in the Damping Ring

- It is possible to include radiation damping
- It is possible to use different scattering dynamics
- It is compatible with MAD X

But it needs to be made compatible with the classical IBS calculations

THANKS.

The End