

- **Physics model**
- **Code implementation**
- **Typical usage**
- **Performance**
- **Documentation**
- **License**
- **Future plans**

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- The Beam-beam force
	- **depends on the bunch parameters**
	- E is highly non-linear
	- **Couples all bunches together**

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- Many different oscillation mode
- Interplay with other effects

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- Linear transport (6D, with chromaticity, linear bucket)
- **Head-on collision (HFMM 4D, FPPS 4D, soft-Gaussian 4-6D)**
- **Long range collision (soft-Gaussian 4D)**
- Noise source (white / colored)
- Collimator ("*in or out* " model)
- \blacksquare Impedance (equidistant slices and wake tables)
- **Linear detuning (e.g. due to octupole)**
- **Transverse feedback (perfect, ADT-like)**
- Synchrotron radiation (damping, quantum excitation)

Impact on CERN machines Colliding beam stability

instability of colliding beams in the LHC, damped by the transverse feedback in agreement with COMBI simulations

Horizonta

 Simulation of coupled bunch coupled beam instability of two trains of 36 bunches colliding long-range

Impact on CERN machines Beam transfer function

- Beam transfer function measurements are used as a diagnostic tool to quantify Landau damping
- Can simulate complex features that are not accessible with the analytical models (e.g. chromaticity, distorted particle distributions)

Impact on CERN machines Emittance effects

- Noise on colliding beams generates an important emittance growth
	- \rightarrow specification for HL-LHC crab cavities
	- \rightarrow specification for PC ripple, ground motion tolerance, etc (HL-LHC, FCC-hh)
- Can simulate complex features that are not accessible with the analytical models (e.g. 2nd order effects, tune effects, chromaticity, …)
- \blacksquare Identify other observables to test the models
- **Solver noise is critical for this** type of study

Impact on CERN machines Emittance effects

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 0.28

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- Need to solve Poisson equation in 2D (open boundary) twice per interaction per turn (i.e. efficiently): several methods exist
- Compute first order moments of the distribution and compute the field based on a Gaussian distribution : −*r* 2
- Fastest and most precise solver, yet it is not accurate especially for non-Gaussian beams
- Used to compute long-range beam-beam interactions since the dependency on the details of the beam distribution is less critical
- Sufficient for most coherent stability studies

Head-on beam-beam interaction Fast Multipole Method (4D)

Group far particles and approximate the kick with a multipolar expansion :

- Smart bookkeeping allows to do reduce the complexity from $O(N^2)$ to $O(N)$
- Accurate (≠precise) for any type of distribution
- Still quite slow for large particle density, especially when the beams are slightly separated
- \blacksquare In a parallel implementation, the position and charge of all particles have to be exchanged between processes, could be slow in clusters without shared memory

Head-on beam-beam interaction Hybrid FMM (4D)

- Distribute the charge of the core on a rectangular mesh, to reduce the computing time and the amount of data to be transfered
	- Grid properties are set with preprocessor directives (need to recompile!)
- E Initially designed for space charge, the HFMM was also used in BEAMX (single 6D beam-beam interaction)
- Working horse of COMBI, since it was heavily benchmarked against analytical formulas, other codes (BeamBeam3D) and observations at RHIC and LHC

- The implementation of the HFMM is robust and fast for most application, but is difficult to maintain and would require major refactoring to be parallelisable (and probably wouldn't be efficient anymore!)
- The quad tree algorithm is very efficient, but really noisy
	- Shows a saturation of the noise level at large number of macroparticle $($ >10 $⁶)$ that is not compatible with present needs</sup>
- We (A. Florio) looked into possibilities to improve :
	- Parallel version of the FMM, with a fixed grid (F2M2)
	- **Fast Polar Poisson Solver (FPPS)**

- Distribute the charge on a polar mesh, solve Poisson on the angular coordinate using the FFT and finite differences on the radial coordinate
	- **Removes the need for copies as well as the artifacts** introduced with a FFT solver on a 2D Cartesian mesh
	- Stretch the radial coordinate to simulate open-boundary condition
	- **Fast, accurate and more precise** than the HFMM
	- **Trivial implementation of the** second level of parallelisation with OpenMP
	- **Singularity at the center is not** trivial to handle

Head-on beam-beam interaction Soft-Gaussian 6D model

Based on K. Hirata's BBC

 \rightarrow Compute the weak-strong beam-beam kick for all particles, taking the other beam's longitudinal slices' first order moments to model the strong beam (soft-Gaussian)

24 Note : A fully self-consistent 6D solver is implemented in BEAMX (HFMM) and BeamBeam3D (FFT) \rightarrow higher computing requirements

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Impedance

A la HEADTAIL :

- Slice the bunch longitudinally (equidistant)
- Compute the charge and average positions of each slice
- Apply the kicks to trailing particles and trailing bunches based on the wake function (wake table or resonator implemented in the development branch)

→ Requires *long term* (i.e. few turns) data storage and communication between bunches \rightarrow see later

- Only single kick per turn possible
- **-** Benchmarked with HEADTAIL multibunch (N. Mounet) (TMCI, coupled bunch instability rise times, octupole thresholds)

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Synchrotron radiation

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- Full implementation (Based on S. White's):
	- **For each particle at each arc (or** each turn), compute the number of photons emitted based on Poisson distribution
	- Compute each photon energy based on probability density (analytical + lookup table)
	- Apply longitudinal and transverse kicks
- Gaussian noise model (default)
	- When several photons are emitted per arc (or per turn), the energy loss per particle becomes Gaussian distributed

 \rightarrow Apply averaged damping on all particles, with a single particle Gaussian white noise (longitudinal and transverse) (based on either the radiation integrals or the equilibrium emittance and damping time)

- Some old (1987) routines are written in F77 (HFMM, BBC), the newest features are implemented as C++ class (FPPS*)
	- **Most features are implemented in F90 or C**
- The wrapping and the first level of parallelisation (MPI) $($ \rightarrow combi.c, master.c, slave.c) is written in C
- Compiled and tested with gcc and icc only
- A second level of parallelisation based on OpenMP is implemented in all functions (only where it offers a gain)
	- Requires an implementation of MPI with the level of thread support MPI_THREAD_FUNNELED (note: usually MPI_THREAD_SINGLE still works, but no guarantees)
- The FPPS relies on FFTW3
- * Pythonized version available in PyPIC

Master and slave initialisation

- At initialisation the process with ID 0 becomes the masters (combi.c \rightarrow master.c)
	- Reads input files (checks their integrity)
	- Makes a mapping MPI ID to bunches, cancels unused processes
	- Sends initialisation data to slaves and waits for return values of all slaves
	- Do one test turn (check integrity, compute collision pattern, both the slave and the master allocate the memory that they will need during the execution)
- Other process are slaves (combi.c \rightarrow slave.c)
	- Read input files
	- Allocate memory for the beam (array of double representing 7 coordinates (6D phase space + charge) of each particle
	- **Enter the 'while(1)' loop:**
		- Wait for an instruction from the master (action code, …)
		- Execute (slave.c \rightarrow Fortran / C / C++ functions)
		- Send completion message

$*$ in

- Name of other input files
- Select type of output and set output file names
- **Machine and beam parameters**
- \blacksquare * fill
	- Define the bunch configuration (i.e. filling scheme in the LHC)
- $\overline{}$ $\overline{}$ $\overline{}$ $\overline{}$ $\overline{}$ $\overline{}$
	- Define the list of actions representing the machine (equivalent to the sequence in MAD, but for two interacting beams)

First level of parallelization

- Usually, the number of bunches is larger than available CPU per node → MPI required
	- 1 **master** process
	- **1 slave** process per bunch
- Both beams go through the same sequence of action, but in opposite direction
	- a, It is the responsibility of the user (with the help of few helper output at initialisation) to ensure the consistency between the filling scheme and the action sequence

Comunication

Comunication

Comunication

Some effects (now only the impedance \rightarrow ADT features ?) requires memory of the particle distribution of the different bunches over a some turns

 \rightarrow A PassageRecord instance is created evey time a bunch reaches an Impedance Action which contains :

- **The absolute time of the interaction**
- First order moments of the particle distributions (either per bunch or per longitudinal slice)
- The PassageRecord is send to the master and forwarded to the slaves when needed
	- Each slave keeps in memory its own deque with PassageRecord of all other bunghes
	- Maximum length of the deque is set by the 'wake length' in input

Typical setup

* The total number of bunches in the FCChh would be out of reach currently

First level of parallelization Performance

- \blacksquare Total number of actions to be performed $\sim N^{\,2}_{\,bunch}$
- Number of actions per bunch ~ *Nbunch*

First level of parallelization Drawbacks

- No gain for actions requiring heavy processing
- Potentially large waste of resources (waiting processes)
	- **Figure 1** Flexibility in the action sequence is a requirement driven by the different needs

- No change on the first level of parallelization
- **Parallelize loops using OpenMP**

- ✔ *Simple implementation*
- ✔ *Many optimisation features available in OpenMP*
- ✗ *Resources are wasted because of the busy-waiting of idling MPI processes*
	- 42 *cluster (use with care, even on your machine...)* ✗ *Control of priorities at the OS scheduler level is available in COMBI (enable RTS) but it requires privileges that are usually not granted on a shared*

- **Requires major modification of the first level of** parallelization :
	- Third type of process : **helper**
- **Slaves** and **helpers** have access to shared memory (using POSIX memory mapping)

- ✔ *No idling MPI process*
- ✗ *Complicated implementation*
- ✗ *Work sharing algorithm have to be implemented*

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Performance Study cases

Low waste (High perfromance

with first level of parallelization) :

High waste (low perfromance with • first level of parallelization) :

 \sim ~95% of the code

seems parallelizable

Performance COMBhy – high waste config.

- 23 bunches \rightarrow 47 MPI processes running on 48 CPUs, only 2 of which are working at a time
	- \rightarrow Waste of resources still present

Performance COMBhy – high waste config.

- The waste of resource can be mitigated by introducing system calls that to lower the priority of idling MPI process with respect to processing ones in the OS scheduler
- The gain is marginal in absence of oversubsribtion but otherwise significant
- Changing processes priorities requires capabilities that are not always granted

Performance

- The second level of paralelisation offers a major speedup wrt to the single level only, both in the high and low waste configurations
- The shared memory version does not allow for a major speedup in the high waste configuration

Solver performance

- The noise introduced by the field solver is critical, especially when studying slow emittance effects
- **Estimated by executing it on a** random distribution for which an analytical formula exists
	- Average of several seeds

$$
\Delta_{\text{solver}} = \frac{\langle |k_{\text{solver}} - k_{\text{th}}| \rangle}{\langle |k_{\text{th}}| \rangle}
$$

- The best performing is the soft-Gaussian as it is fast and its noise approaches the theoretical minimum 1/sqrt(N)
- The FPPS offers a major speed up for a given noise amplitude wrt to the **HFMM**

FPPS parallelisation performance

Excellent speedup using OpenMP on the most of the loops (not within fftw)

- Many studies are run on local machines (usually 4 to 12 cores), which performances allow for decent studies
- **For most studies the code performs well, but requires** proper parallel infrastructure, as well as manpower for setting up, testing and optimisation on each architecture

 \rightarrow Preferred resources : Multiple nodes with fast connection, regular memory and possibly several cores per node

- **EPFL infrastructures** (accessible on request by L. Rivkin's Ph D. students) offer such capabilities, most results obtained with COMBI where based on their facilities
- **Need to investigate potential of current CERN** infrastructures

Future plans

- **Current implementation has great potential**
	- Maintain and use the code to produce results (including implementation of 'minor' actions for interplay studies)
- Emittance studies are limited by the amount of numerical noise due to the field solver with a finite number of macro-particles (10 $^{\circ}$ \rightarrow 10 $^{\circ}$)
	- GPU acceleration is not suited for strong-strong simulations due to the communication requirements (imposed by the physics)
		- \rightarrow Need for computing resources (large number of CPUs per node)
- Long term stability studies (either due to slow rise times, or slow distorsions of the particle distribution) requires several turns (~10 $\mathrm{^6}$ or more) and several parameter scans

 \rightarrow Need for computing resources (large number of CPUs per node, possibly large number of nodes for multibunch studies)

Effort on the 6D beam-beam effects just (re-)started

 \rightarrow Effectively increases the computational need by a factor \sim 50 (number of slices per bunch)

- Testing and parallelisation of the 6D soft-Gaussian solver
- Ľ Implementation of a self-consistent 6D solver ?

Documentation

- Source are available on svn : <https://svnweb.cern.ch/cern/wsvn/COMBI>
	- **doc/action codes**
	- doc/input conventions
	- doc/goottoknow
- <https://cds.cern.ch/record/1987672/files/CERN-THESIS-2014-246.pdf> appendix A
- <https://twiki.cern.ch/twiki/bin/view/ABPComputing/COMBI>

better than nothing

- CERN copyright

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