Experimental and numerical studies on the proposed application of hollow electron beam collimation for the LHC at CERN

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Abstract

This thesis work was carried out in the framework of U.S. LHC Accelerator Research Program (US-LARP), a collaboration between the European Organization for Nuclear Research (CERN) and the U.S. Department of Energy. The first half of the work was completed at Fermilab (USA), the location of the Tevatron, a proton-antiproton collider and the second largest particle collider in the world. The second half was completed at CERN (Switzerland), the location of the largest proton collider in the world (Large Hadron Collider (LHC)).

This thesis characterizes a Hollow Electron Beam (HEB) for possible usage at the LHC to enhance its collimation through Hollow Electron Beam Lenses (HEBLs). Collimation is a long established principle in high energy particle accelerators. Hollow Electron Beam Collimation (HEBC) aims to enhance current collimation systems by controlling diffusion of primary halo particles into the limiting aperture. It works on the principle of a transverse radial electric field that kicks the primary halo particles outwards upon each pass in a multi-pass system. The transverse field is produced by a HEB that is coaxially aligned with the accelerator beam, producing a negligible electric field in the center and a strong transverse electric field at amplitudes higher than the inner radius of the electron beam. Ideally, halo particles are affected without perturbation of the beam core. One of the main advantages of this system is to decrease the dependence on instantaneous loss spikes and beam jitter. A solid experimental basis of HEBC was accumulated at the Tevatron. The application of this technique at the LHC is now under investigation.

The aim of this thesis is to present a preliminary report to support a future optimal conceptual design report. It characterizes the available hardware in order to facilitate the design of a Hollow Electron Gun (HEG) for the LHC, characterizes the effect on beam diffusion by determining the transverse electric fields of the electron beam and initiates 3D simulations in order to determine the effect of the HEBL on the beam core. Experiments were conducted in the Tevatron Electron Lens Test Stand (TELTS) at Fermi National Accelerator Lab (Fermilab) (USA) in the lower Linear Accelerator (LINAC) gallery.

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Acronyms

Description	Page List
Accelerator Command Language	29-31
Fermilab Accelerator Control System	20, 29, 30
Brookhaven National Laboratory	6
European Organization for Nuclear Research Computer Graphics Metafile Center Of Mass	I, 1, 5, 56, 57 49 4
Drift Tube LINAC	3
Emittance Growth Rate	44, 47
Fermi National Accelerator Lab	I, 1, 4, 9, 12, 23, 48
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Hollow Electron Beam	I, 1, 6, 7, 14, 15, 20, 33, 40,
Hollow Electron Beam Collimation Hollow Electron Beam Lens	41, 47, 55–57 I, 1, 3, 7–10, 20, 57 I, II, 1, 2, 7, 8, 11, 18, 20, 29, 41, 43, 48, 50, 56, 57
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Linear Accelerator	I, 4
Particle-In-Cell	48, 49, 57
Relativistic Heavy Ion Collider	6
Root Mean Square	5, 43, 44
Space Charge Limited Emission Region	15, 16, 18, 19, 23, 24, 26
	DescriptionAccelerator Command LanguageFermilab Accelerator Control SystemBrookhaven National LaboratoryEuropean Organization for Nuclear Research Computer Graphics Metafile Center Of MassDrift Tube LINACEmittance Growth RateFermi National Accelerator LabGNU's Not UnixHollow Electron BeamHollow Electron Beam Collimation Hollow Electron Gun 0.6 inch Hollow Electron Gun 1 inch Hollow Electron Gun 1 inch Hollow Electron Gun 1 inch Hollow Electron Gun Particle-In-CellRelativistic Heavy Ion Collider Root Mean SquareSpace Charge Limited Emission Region

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Chapter 1

Introduction

This work was completed in the framework of the U.S. Department of Energy collaboration with the Large Hadron Collider (LHC), U.S. LHC Accelerator Research Program (US-LARP). The experiments were completed between February and May 2013 at Fermi National Accelerator Lab (Fermilab) in Illinois, USA. Simulations were done at European Organization for Nuclear Research (CERN) in Switzerland in June and July 2013. The work incorporates aspects of the Tevatron, a proton and antiproton collider, and the LHC, a proton-proton collider.

This thesis addresses the characterization of a new 1 inch Hollow Electron Gun (HG1b) tested at Fermilab for LHC collimation purposes in the framework of Hollow Electron Beam Collimation (HEBC). Considering successful soft collimation experiments using a 0.6 inch Hollow Electron Gun (HG06) in the Tevatron Electron Lens 2 (TEL2) at the Tevatron, this thesis outlines important design parameters for a Hollow Electron Gun (HEG) and an electron lens for the LHC. It measures the transmission, yield and transverse profiles of the new HG1b, which fits the general LHC requirements. Furthermore it determines the transverse electric fields and the beam evolution of the Hollow Electron Beam (HEB). Additionally upper limits for the emittance growth of the proton core due to the HEB are determined.

The first Tevatron electron lens was installed in 2001. It was used on a daily basis for abort gap cleaning for a period of 10 years until the recent shutdown of the Tevatron in 2011. A second Tevatron electron lens was installed in 2006 as a backup for the first lens. It was mainly used for research on beambeam force compensation and space charge compensation [1]. More recently, the lens has been tested as a soft collimator by installing a HEG. Considering the 10 years of prolonged stable operation at the Tevatron, electron lenses have established themselves as a mature technology for beam manipulation.

HEBC is designed to be used together with current collimation systems in order to control the diffusion of halo particles into the limiting aperture through transverse electric fields. The proton core at the same time remains unaffected due to the negligible electric field that is found in a cylindrical hollow electron beam.

HEBC studies have been part of US-LARP since 2009. There is a strong interest to implement two electron lenses in the LHC. The aim is to provide a solid basis for the design of those lenses, in case these are needed when the LHC reaches nominal energy. The TEL2 is now available for use at LHC, but constraints such as cryogenics and a long installation time make this a difficult task. Therefore the plan is to design two new electron lenses, optimized specifically for operation at 7 TeV with improved integration possibilities and instrumentation [2, S.27-28].

The *CERN* strategy for the *Hollow Electron Beam Lens (HEBL)* was presented at the 20th *US-LARP* collaboration meeting in Napa Valley, CA, USA, in April 2013, it was agreed that *Fermilab* would work on an optimum conceptual design for an implementation in the LHC. This study would entail a discussion of the hardware parameters and optimal design parameters for the LHC, the effect of the *HEBL* on beam dynamics and beam core (luminosity) and the enhancement of collimation performance. The effect on beam dynamics was addressed in [3]. This recent paper shows that with an AC beam mode operation, a scraping of 75% of the halo particles was possible in 20 s[3].

This thesis will address the hardware and optimal design parameters as well as the effects of HEBC

on beam dynamics and the beam core. This is done by discussing current Tevatron hardware, measuring the yield and transverse profiles of the electron beam and determining the transverse electric fields of the electron beam. Additionally this thesis initiates a 3D simulation using WARP which will be used in order to obtain a 3D non-linear kick model for implementation in Lifetrack or SixTrack. This will allow the integration of proton trajectories in the LHC passing through the HEBL and thus a analysis of the effect of the electron lens on beam halo diffusion and on the beam core. An overarching aim of this thesis is to present a first study that will support a future optimal conceptual design report.

Chapter 2

Particle Accelerators and Collimation

This chapter introduces particle collimation as a necessity for current high energy particle accelerators. It discusses the limitations of classical collimation systems and elaborates on *Hollow Electron Beam Collimation (HEBC)*.

2.1 Accelerators and Particle Collimation

The origin of particle accelerators can be traced back to Rutherford's 1910 scattering experiment of α particles on gold foil at Cavendish at Cambridge University in England. In 1927, Ernest Rutherford addressed the Royal Society of London, expressing his desire for the development of a device that could continuously produce α and β particles at energies much larger than their natural decay energies. His goal was to disintegrate nuclei with binding energies that exceed that of nitrogen [4]. In 1930, the first particle accelerator was conceived by John D. Cockroft and E.T.S Walton at the laboratory [5]. Since it was yet impossible to accelerate particles such as α particles to energies beyond their radioactive equivalents, Cockroft and Walton decided to work with lighter particles, accelerating protons through a linear discharge tube with a potential of 200 kV. Lacking results with such "low energy" particles, they conceived a voltage multiplier, called the Cockroft-Walton generator (See fig. 2.1), allowing them to accelerate protons to 800 keV, which subsequently allowed them to force the decay of Lithium into two α particles, using protons at 500 keV.



Figure 2.1: Cockroft Walton Generator¹

In 1931, Van der Graaf and Wideröe followed suit. Especially Rolf Wideröe's design, the *Drift Tube LINAC (DTL)*, was a cornerstone for the development of today's particle accelerators. It accelerated particles through axially aligned drift tubes, separated by strong electromagnetic fields. By applying an AC current, he could tune the frequency of the field, such that the particles would be boosted upon every exit of a drift tube (see fig. 2.2). This design removed the need to store huge electric charges, which limited Cockroft and Van der Graaf's designs. While Wideröe's design was limited to around 50 keV of energy, it was his design that inspired Ernest Lawrence to develop the cyclotron, a device that is widely known as the forerunner of particle accelerators (see fig. 2.3). In effect, the cyclotron is equivalent to winding Wideröe's design around and around itself to create a magnetically confined beam orbit with an alternating electric field.

In recent years, the spectrum of different applications of particle accelerators has widened. Leadingedge accelerators are driven by particle physics requirements (beam energy, high luminosity and intensity), but the majority of operating accelerators serve other purposes such as chemical or material research and medical treatments. For a history on particle accelerators, see [6].

¹http://www.visualphotos.com/photo/1x6033020/cockroft-walton_generator_fermilab_a090057.jpg

²http://www.fnal.gov/pub/science/experiments/energy/images/95-1039-small.jpg

Today's state of the art accelerators reach energies that are approximately 10^5 times higher. The highest to date recorded acceleration energy was observed at the *Large Hadron Collider (LHC)* with a value of 4 TeV, giving 8 TeV in the *Center Of Mass (COM)* frame. It is designed to hold a stored energy of ≈ 360 MJ for each beam [7]. Such energies require the use of superconducting magnets to steer the beam, resulting in the need for strong collimation. Even under good stable operational conditions, particles will leave the optimal design path of the accelerator due to beam-gas interactions, intra-beam scattering, beam-beam interactions, RF noises, ground noises and resonances due to imperfections results in a slow diffusion of protons from the axial path and a subsequent growth of the emittance. These particles together form the halo, a set of off-axis particles whose betatron amplitudes increase gradually until they are caught by collimators or other limiting apertures.

The beta function is the envelope around all particle trajectories in the accelerator. The betatron amplitude is the transverse amplitude with respect to the nominal beam orbit of individual particles in the accelerator. The collision of particles with the limiting aperture causes hadronic and electromagnetic showers, which can negatively affect accelerator components and detectors with different severities.

Collimation refers to the spatial alignment and the reduction of the spatial cross-section of a beam halo. It is the removal of the beam halo, in order to protect components against excessive irradiation, prevent quench[es] in the superconductors, minimize background noise in the detectors and reduce the radiation on personnel and the environment [9, 10, p. 3, p. 297]. Collimators can act as diagnostic tools for accelerator admittances, beam vibrations and diffusion rates [11].



Figure 2.2: Drift Tubes in the LINAC facility at $Fermilab^2$



Figure 2.3: Ernest Lawrence with his cyclotron design 3

2.2 Beam Collimation Systems

Collimation of beam halo particles is achieved by limiting the physical aperture of the accelerator beam using solid blocks of absorbing material. Through scattering, a small fraction of halo particles can escape these blocks. We call such particles out-scattered particles. The cleaning efficiency, the leakage of halo protons to sensitive equipment, of the collimation system is improved by using several collimation stages. Such a system consists of primary collimators, whose purpose is to intercept the primary beam halo through multiple Coulomb scattering. Secondary collimators are used to intercept the secondary halo particles which are out-scattered from the primary collimators at higher amplitudes [12]. The LHC uses a complex four stage collimation system. A list of the current LHC collimators is given in tab. 2.2.

2.2.1 Tevatron Collimation System

Tevatron had two Collider Runs. Inbetween the two runs, the collimation system of the Tevatron was upgraded [9]. During Collider Run I, the Tevatron used a single stage collimation system that consisted of 1 m long solid absorbers which allowed *Fermi National Accelerator Lab (Fermilab)* to raise the Tevatron efficiency of the fast resonant extraction system by a factor 5 [13]. Using this system at Tevatron, one obtained a low cleaning efficiency close to 0.5 [13]. This system was insufficient and caused operational limitations, such that in Collider Run II, a two-stage automated collimation system was implemented [9]. At Tevatron collimation is a discrete process in which the beam halo is regularly removed. The new design required that the entire halo removal could be conducted in approximately 5 min. The system incorporated four primary collimators (targets) and eight new 1.5 m long secondary

³http://newscenter.lbl.gov/wp-content/uploads/lawrence-cyclotron.jpg

Interaction	Purpose	Collimator
Regions	Purpose	Types
IR1	ATLAS experiment	TCT, TCL
IR2	ALICE experiment	TCLI, TCT, TDI
	& Injection Beam 1	
IR3	Momentum cleaning	TCP, TCSG, TCLA
IR4	RF cavities	-
IR5	CMS experiment	TCT, TCL
IR6	Beam Dump	TCSG, TCDQA
IR7	Betatron Cleaning	TCP, TCSG, TCLA
IR8	LHC-b experiment	TCLI, TCT, TDI
	& Injection Beam 2	

Functional type	Name	Plane	Num.	Material
Primary IR3	TCP	Н	2	CFC
Secondary IR3	TCSG	Н	8	CFC
Absorbers IR3	TCLA	H,V	8	W
Primary IR7	TCP	H,V,S	6	CFC
Secondary IR7	TCSG	H,V,S	22	CFC
Absorbers IR7	TCLA	H,V	10	W
Tertiary IR1/2/5/8	TCT	H,V	16	W/Cu
Physics debris absor.	TCL	Н	4	Cu
Dump protection	TCSG	Н	2	CFC
	TCDQ	Н	2	С
Inj. prot. (lines)	TCDI	H,V	13	CFC
Inj. prot. (ring)	TDI	V	2	С
	TCLI	V	4	CFC
	TCDD	V	1	CFC

Table 2.1: The 8 interaction regions, their purpose and collimator types.

Table 2.2: Current LHC Collimators. For a complete list of Acronyms, visit the Acronym website of the LHC^4 .

collimators. The primary collimators are 5 mm tungsten plates, placed at 5 σ . σ is the standard deviation of the Gaussian distribution of the proton beam in the LHC. They create a *Root Mean Square (RMS)* transverse kick of 17 μ rad at 980 GeV [11]. The secondary collimators are 1.5 mm steel blocks placed at 6 σ [9]. Additionally a new control system was implemented that would allow the fast processing of beam loss monitor data and a fast beam intensity feedback control, in which the full travel of the collimator took 15 sec [9]. Next to this major upgrade, the system was upgraded on a regular basis, including the installation of a tertiary collimator in 2003.

Measurements have shown that the cleaning inefficiency was 0.001. At the Tevatron collimation is primarily used for background reduction in physics, which was of the order of 6.7×10^{-3} , defined as the ratio of background noise with and without collimation [9, p. 7]. The estimated evolution of beam losses due to $\bar{p}p$ collisions, losses in the RF-bucket and beam-gas scattering are reviewed in [13, p. 14].

2.2.2 LHC Collimation System

Particle accelerators such as the *LHC*, with a nominal stored energy of 360 MJ require more collimation systems. At the *LHC*, the beam is continually cleaned. Even small losses in the *LHC* of the order of 4×10^7 protons, could cause quenches in the superconducting magnets through a deposition of energy in the order of 30 mJ cm⁻³ [14].

The beam cleaning system at the LHC requires a cleaning efficiency of 99.998%. The *LHC* has two cleaning regions in *Interaction Region* (*IR*)3 and *IR*7 and several additional collimators in the other sections, with a total of 100 collimators, absorbers and similar devices [7]. There are 8 *IRs* in the LHC with different purposes. The *IRs* and their purposes are listed in table 2.1. The cleaning insertions have primary (TCP), secondary (TCS) and absorber (TCLA) collimators. Tertiary collimators (horizontal TCTH and vertical TCTV) are placed upstream of the experiments in *IR*1, *IR*2, *IR*5, and IR8. The machine is protected in *IR*6 against beam dump failures through dump protection devices (TCS6 and TCDQ) [7].

2.3 Hollow Electron Beam Collimation (HEBC)

Considering projects such as the High Luminosity LHC (HL-LHC) at European Organization for Nuclear Research (CERN), which aims at increasing the luminosity of the LHC by a factor 10, or the Very Large Hadron Collider (VLHC), designed to operate at 20×20 TeV in Stage 1 and 88×88 TeV in Stage 2 [15], new innovative collimation techniques are required.

Some of the limitations of classical collimators is that they can not operate arbitrarily close to the beam orbit, due to impedance effects and instantaneous loss rates [11]. Additionally, beam loss rates

⁴http://lhccwg.web.cern.ch/lhccwg/Bibliography/UsefulAcronyms.htm



Figure 2.4: Technical Drawing of TEL2 Setup. Source: [8]

increase dramatically when moving collimator beam jaws inwards. Furthermore, classical collimators are affected by beam jitter, oscillations of the accelerator due to mechanical noise or ground motion [8]. Even with active stabilization, periodic bursts at the beam aperture are caused.

Electron lenses employ transverse fields of electron beams to interact with high energy bunches. An electron gun is used to produce an electron beam that is steered into the accelerator beam tube. It is aligned with the proton beam orbit, overlapping over a given length L_e , after which it is extracted from the accelerator beam tube into a collector. The electron beam is magnetically confined through solenoids. The electron gun and therefore the beam can be pulsed. All current electron lenses have the electron gun and the collector positioned outside the accelerator beam tube. A sketch of the *Tevatron Electron Lens 2 (TEL2)* is given in fig. 2.4.

An electron lens can produce different transverse profiles. For example, a Gaussian profile was used for non-linear beam-beam force compensation in the Tevatron and a flat-top profile was used for bunch by bunch tune correction. Further profiles are explained in [16]. An important advantage of the electron lens is that the electron beam current can be adjusted between individual bunches at the Tevatron, allowing single bunch manipulations. A normal figure of merit for electron lenses with Gaussian profiles is the shift of the accelerator beam tune due to the lens. A perfectly steered round electron beam, shifts the betatron tune by [17]

$$dQ_{x,y} = \pm \frac{\beta_{x,y} L_e r_p}{2\gamma ec} j_e \left(\frac{1 \mp \beta_e}{\beta_e}\right)$$
(2.1)

where $\beta_e = v/c$ is the electron beam velocity, γ the Lorentz factor and $r_p = e^2/mc^2 = 1.53 \times 10^{-16}$ is the classical proton radius. The plus sign represents focusing for protons and defocussing for anti-protons.

There are two electron lenses installed in the Tevatron, *Tevatron Electron Lens 1 (TEL1)* and *TEL2*. Both operate around 10 keV and at a couple of amperes. They can produce a tune shift of $dQ_{x,y}^{\max} \approx 0.008$ [16]. They were installed for research on beam-beam compensation [17, 16]. *TEL1* was installed in 2001. It has since been used in normal operation procedure for abort gap cleaning [18]. *TEL2* was installed in 2006 as a backup for *TEL1*. Beam-beam compensation was observed, but was not needed after the installation of electron cooling for antiprotons in the recycler ring. Due to the reliability of *TEL1*, *TEL2* was rarely needed for normal operations and thus used for further experiments on beam-beam compensation with different electron guns [19, 11]. Electron lenses for beam-beam compensation are currently being comissioned at *Relativistic Heavy Ion Collider (RHIC)* at *Brookhaven National Laboratory (BNL)* [20].

With an appropriate gun design, *Hollow Electron Beams (HEBs)* can be produced. Such a beam can be used for enhancing collimation. The *HEB* is aligned coaxially with the proton beam, such that the proton beam core travels through the negligible electric field in the hollow space of the *HEB*. Halo particles of the proton beam at transverse amplitudes bigger than the inner radius R_i experience near-



Figure 2.5: Radial electric field for profiles corresponding to V=500V and 8kV, I=0.073A and 3.88A, $R_i=3.4mm$ and $R_o=6.4$ mm.

linear transverse radial kicks up to the outer radius R_o . Particles above R_o experience a non-linear decay of the transverse kicks with respect to amplitude due to the Biot-Savart decay of the transverse electric field of the *HEB* above its outer radius. A typical electric field profile is shown in fig. 2.5. Its derivation is discussed in chap. 6.

Assuming a cylindrical symmetry of the current density profile, transverse kicks are very controllable and given by [11]:

$$\Theta = \frac{1}{4\pi\varepsilon_0} \frac{2I_r L(1\pm\beta_e\beta_p)}{r\beta_e\beta_p c^2(B\rho)_p}$$
(2.2)

where $\beta_e c$ is the electron velocity, $\beta_p c$ is the particle velocity, I_r is the enclosed electron current, L is the interaction length of the *Hollow Electron Beam Lens (HEBL)* and r is the distance from the axis of the beam. The + applies when the magnetic and electric forces have the same direction. For typical Tevatron parameters, the transverse kick given to 980 GeV particles in Tevatron is 0.2 μ rad.

The magnetic fields in the gun, collector and main solenoid can be varied individually. This allows the operator to compress the *HEB* in the main solenoid. The compression follows the scaling law $\sqrt{B_{main}/B_{gun}}$ and the decompression law $\sqrt{B_{coll}/B_{main}}$. The magnetic field in each solenoid has an upper limit of 6 T, which is easily attainable by superconducting magnets. It has a lower limit of ≈ 0.1 T due to scalloping of the beam. Scalloping refers to transverse oscillations in the beam due to transverse electric fields. Scalloping is reduced by magnetically confining the particles stronger. When increasing the *HEB* size, it becomes necessary to compress the beams. This is shown in sec. 5.2.

The basic concept of the HEBC is that R_i is chosen to be smaller than the half gap of the primary collimators (TCP). It then causes the diffusion of particles into the limiting transverse aperture between R_i and the transverse aperture. A sketch is shown in fig. 2.6. The proton beam is represented through the blue curve with long tails representing the halo. The diffusion coefficient is represented by the green curve. HEBC allows the cleaning of the long tails of the transverse proton beam distribution. This makes the system less dependent on loss spikes such as those caused by beam jitter [8]. It is important to note that HEBC requires that the current collimation system is kept in place.

Various affects can reduce the effectiveness of HEBC. Important requirements are: a sufficient beam current to produce significant transverse kicks, a negligible electric field in the center to reduce emittance growth and a symmetric beam distribution. Beam current should be above 1 A in as was seen from experiments at the Tevatron [11]. The electric field is strongly affected by radial symmetry of the HEB. Furthermore bends in the electron lenses can adversely affect the electron beam distribution and the proton core as it pierces through the HEB wall in the bend (see fig. 2.4.) The effectiveness of the HEBC depends furthermore on the time structures of the pulsed electron beam. Valentina Previtali recently



Figure 2.6: Schematic representation of the effect of the HEBC on the diffusion model of collimation. Source: [8].

completed a study of 3 different time structures; DC current, AC current and a "diffusive" current. These are possible, because the cathode has a rise-time short enough to allow turn by turn modulation in the *LHC*. Her studies show that the AC beam mode is the most efficient with a scraping of 75% of halo particles in less than 20 s. For more details refer to [21].

The HEBLs which were used for HEBC and its characterization are discussed in sec. 3.1. Furthermore a conceptual design of a HEBL where the Hollow Electron Gun (HEG) and the collector are coaxially aligned with the accelerator beam is discussed. HEGs are an important part of the HEBLs and are discussed in sec. 3.2.

2.3.1 Tevatron experience for HEBC at LHC

Tevatron experiments from October 2010 to September 2011 gave an experimental foundation for the HEBC concept in the context of beam studies for the LHC collimation purposes. They were done using a 0.6 inch Hollow Electron Gun (HG06) in the TEL2. The experiments showed that HEBLs are compatible with collider operations and that the alignment of the electron and proton beam is reliable and reproducible. It could be shown that the effects on the beam core were negligible, while a smooth halo removal was observed. Furthermore the loss-rate fluctuations due to beam jitter were reduced. Overall a transverse diffusion of the beam halo was observed. These achievements of the TEL2 at the Tevatron are explained in more detail in [11, 22, 8]. After 10 years of stable Tevatron electron lens operation in the Tevatron (TEL1 and TEL2), it is clear that the technology is reliable and thus ready for operation in the LHC.

Chapter 3

Tevatron Electron Lenses and Guns

The Tevatron Electron Lens 2 (TEL2), installed in 2006, was used as a backup for Tevatron Electron Lens 1 (TEL1) for abort gap cleaning and for studies on beam-beam compensation and Hollow Electron Beam Collimation (HEBC). The HEBC studies at the Tevatron were done using a 0.6 inch Hollow Electron Gun (HG06). Since 2012 a new 1 inch Hollow Electron Gun (HG1b) has been used in the Tevatron Electron Lens Test Stand (TELTS), which is used for yield and transverse profile measurements. It was designed specifically for the Large Hadron Collider (LHC). This chapter discusses the electron lenses and electron guns that have been used for the development of HEBC at Fermi National Accelerator Lab (Fermilab) in the context of collimation studies for the LHC. It furthermore outlines design parameters for the HEBC studies at Tevatron and proposes design parameters for the use of the HG1b at the LHC at CERN.

3.1 Tevatron Electron Lenses

3.1.1 Tevatron Electron Lens 2 Design Considerations

The Tevatron Electron Lens 2 (TEL2) is one of two electron lenses installed in the Tevatron. It was used for beam-beam force compensation and for the experimentation with HEBC. A sketch of the TEL2 is shown in fig. 2.4. Unlike the LHC, the Tevatron circulates a proton and an antiproton beam in opposite directions in the same vacuum pipe. An electron gun, situated left next to the accelerator beam tube, produces an electron beam that is sent on an arc through 3 short solenoids into the accelerator beam tube. After traveling a distance of 2 m, during which the elec-

$I_{max} \pm (\Delta I/I)_{max}$	$1.2 \text{ A} \pm 2\%$
V_max	5 kV
R_1/R_2	4.5 mm/7.5 mm = 0.6
R_{1min}	0.58 mm
Δt_{min}	250 ns

Table 3.1: Design parameters of TEL2 in Tevatron with HG06: maximum electron beam current I_{max} and current jitter $(\Delta I/I)_{max}$, maximum extraction voltage V_{max} , ratio between inner and outer cathode radius R_1/R_2 , minimum achievable inner radius for the electron beam R_{1min} and cathode modulator rising time Δt_{min} [21]

tron beam interacts with the (anti-)proton beam, the electrons are again extracted from the accelerator beam tube using 3 more solenoids [16]. The lens operates with a few amperes and at an energy up to 10 keV [16]. The gun, the interaction region and the collector have 3 separate solenoids, used for the magnetic confinement of the electron beam. Their magnetic field strengths are designated B_{gun} , B_{main} and B_{coll} respectively. Inside the main solenoid, 6 superconducting dipole correctors are found. They are used to align the beam with the (anti-)proton beam. This requires five degrees of freedom, three for the axial position of the electron beam upstream of the main solenoid and two for its angle.

Since the shutdown of the Tevatron at Fermilab, the TEL2s hardware has become available for use. Its efficiency as a scraper was proven through extended tests at the Tevatron collider, driving a 980 GeV anti-proton beam onto the collimators[23]. The TEL2 (using the HG06) was used as a HEBC at currents up to 1.2 A. TEL2 parameters used in the Tevatron are given in tab. 3.1. Significant scraping



Figure 3.1: Sketch of TELTS. [24]



Figure 3.2: Photo of TELTS. Source: Valentina Previtali

could be observed at this current [11]. Similar currents are thus desired for new design considerations. For further information on the scraping studies refer to [23, 8, 11].

3.1.2 Tevatron Electron Lens Test Stand Design Considerations

While HEBC tests with a proton beam in an accelerator are crucial, they are not intended for the characterization of the electron beam parameters. For the study of the transverse profile, the *Tevatron Electron Lens Test Stand (TELTS)* is used. A sketch is shown in fig. 3.1 and a panoramic photo in fig. 3.2. The *TELTS* is a linear test bench, where the pulsed electron gun (1) sends an electron beam downstream through a straight beam tube (5) with pickup electrodes to a water cooled collector (2). The whole beam tube is in a vacuum and is surrounded by 3 solenoids. The gun is placed in a gun solenoid (4) which is followed by a short drift space and the main solenoid (3). The collector (2) at the end of the drift tube is placed in the collector solenoid (4). Magnetic correctors (6) are used to control the beam inside the beam tube. The device is supported using active stabilization (7) on a test bench (8) [24].

The distance between the gun cathode and the collector downstream is 2.86 m. Thereof 35 cm are in the gun solenoid, 23 cm in the first drift space, 192 cm in the main solenoid, 15 cm in the second drift space and the remaining 21 cm in the collector solenoid. The magnetic fields in the 3 solenoids can be set independently. The gun and collector solenoid are 50 cm long. The inner diameter of the two short solenoids is 28 cm and the main solenoid is 20 cm [24].

For detailed information on the power supply, the water cooling system and the vacuum system, please refer to the Electron Lens Test Stand Wiki page¹. The *TELTS* was used for all measurements in chap. 5 and implemented in the simulations in chap. 7.

¹https://cdcvs.fnal.gov/redmine/projects/elens/wiki/Test_Stand

3.1.3 Straight Hollow Electron Beam Lens Design

The straight hollow electron beam lens design eliminates the bends required to steer the electron beam into the accelerator beam tube in the usual electron lenses. It is thus no longer necessary for the proton beam to pass through the hollow electron beam edge when entering the *Hollow Electron Beam Lens* (*HEBL*). A sketch of such a device is shown in fig. 3.3. No such device has been built to date. In the straight hollow electron lens design, the electron gun itself is hollow, not just the cathode, and is coaxially aligned with the hadron beam of the accelerator. The electron beam is coaxial to the proton beam, focusing shortly before the main solenoid and expanding again just after. It is extracted by being radially deflected at the electron collector.



Figure 3.3: Sketch of straight HEBL setup. Source: Gennady Kuznetsov (Fermilab)

3.2 Design of the Electron Guns

An electron gun emits free electrons with a specific energy spectrum. The most common application is the cathode ray tube used in old television sets. Among many different electron sources, the simplest is the planar diode, consisting of two oppositely charged electrodes. The cathode emits the electrons, which are accelerated to the anode via a potential difference. The anode contains a hole or a mesh type structure to allow the electrons to pass and continue downstream of the electron gun [25, p. 7].

The most vital part of the HEBL is the electron gun. Two guns were produced for the TEL2 and TELTS. First a HG06, which was used in the TEL2 setup. Second a HG1b was acquired for testing in the TELTS and subsequent installation in the LHC.

Both cathodes are made out of an 80% density porous tungsten matrix, impregnated with 3BaO-CaO-AL₂O₃. At a cathode temperature of about 2500 K ($k_BT \approx 0.2 \text{ eV}$), the work function of pure tungsten is 4.5 eV and it produces a current density of about 0.5 A cm⁻² [25, p. 9]. The work function is the minimum energy required to remove an electron from a solid to a point immediately outside the solid surface in vacuum. Using tungsten cathodes which are impregnated with barium or strontium oxides, considerably higher current densities of 10 to 20 A cm⁻² can be achieved. Such cathodes are typically operated at 1400 K ($k_BT \approx 0.12 \text{ eV}$) and have a work function of 1.6 eV [25, p. 9]. Pure barium cathodes are not used due to their low melting point of 725°C.



Figure 3.4: Technical Drawing of HG06. The hole drilled through the cathode is not included.Source: Gennady Kuznetsov (Fermilab)

3.2.1 0.6 Inch Hollow Electron Gun

HG06 was manufactured by Hi-Tech Manufacturing, LLC (Shiller Park, IL, U.S.A). The model number is 101507. The cathode is commercially available from Heat Wave Labs, Inc. (Watsonville, CA, U.S.A) [26]. It has an outer diameter of 15.24 mm (0.6 in) and a radius of curvature of 10 mm [26]. A 9 mm hole was drilled into the center of the cathode [26]. The gun is equipped with two control electrodes, situated 0.28 mm and 9.38 mm from the cathode. The anode is 9.84 mm downstream of the cathode and has an aperture of 9 mm at the front. Downstream of the anode is a drift pipe with an aperture of 36.5 mm and a length of 93.1 mm. Fig. 3.4 gives a technical drawing of the gun.

This gun was used for the Tevatron scraping experiments [23]. Through magnetic compression it could be scaled to a minimal inner radius of the electron beam of 0.58 mm. The gun has a perveance of $3.20(7) \times 10^{-6}$ perv (see fig. 3.5) and an optimal operating temperature above 1100° C as seen in fig. 3.6. The unit perv is equivalent to the SI units A V^{$-\frac{3}{2}$}. Perveance is defined as the scaling factor between the beam current, I, and the cathode potential to the power of 1.5, $V_a^{1.5}$. The calculated current density distribution is negligible between 0 < r < 4.5 mm, rising sharply at the inner edge, and falling gradually back to 0 at the outer edge [26].

3.2.2 1 Inch Hollow Electron Gun

Given a maximum current of the HG06 in the Tevatron of 1.2 A, a new gun was designed specifically for LHC parameters. The diameter of the cathode was increased to 1 inch in order to allow higher currents. HG1b, custom designed by *Fermilab* and built by Heat Wave Labs Inc., is made out of the same material as HG06. It has an outer diameter of 25.4 mm (1 inch), an inner diameter of 13.5 mm and a radius of curvature of 10.0 mm between the outer and inner radius. The gun is equipped with two control electrodes, situated 0.4 mm and 7.8 mm from the cathode. The

T	N 9 00 A
Imax	\geq 5.00 A
V_{sug}	3-4 kV
R_1/R_2	$6.75 mm/_{12.7 mm} = 0.53$
f_{comp}	3.4
B_{sug}	0.43-5-0.43 T

Table 3.2: Design parameters for the e-lens installed in the Tevatron: maximum electron beam current I_{max} , suggested voltage V_{sug} , ratio between inner and outer cathode radius R_1/R_2 , compression required to place inner beam radius at 4 σ f_{comp} and a suggested magnetic field at the LHC.

anode is 9.8 mm downstream of the cathode and has an aperture of 14.3 mm upstream. Downstream of the anode is a drift pipe with an aperture of 36.5 mm and a length of 93.0 mm. Fig. 3.7 gives a technical drawing of the gun.



Measurements of perveance of HG06 on 06.07.2010 16:43

Figure 3.5: Perveance of HG06 on 06.07.2010. The perveance is $(3.20 \pm 0.07) \times 10^{-6}$. Time elapsed is measured since heating of the cathode.



Figure 3.6: Performance of the HG06 gun with respect to cathode voltage and filament power. Source: [26]



Figure 3.7: Technical Drawing of HG1b. The center of the drawing shows the cathode with the two electrodes around it and the anode and the drift pipe to its left. Source: Alexander Didenko (Fermilab contractor)

The new gun can use higher magnetic fields up to 6 T in the beam solenoid, to improve transport of the electrons through the lens, by reducing space charge expansion and taking advantage of the E \times **B** drift on the electrons. One should note that design plans place the inner radius of the Hollow *Electron Beam (HEB)* at $\approx 4 \sigma$ of the *LHC* proton beam. The proton beam σ was estimated in sec. 6.3 to be ≈ 0.5 mm at the possible installation site of Interaction Region (IR) 4. Positioning the beam at 4 σ requires a magnetic compression of the electron beam inner radius from 6.75 mm to 2.0 mm by a factor of 3.4 in the main solenoid. Given that the magnetic compression scales as B_{main}/B_{qun} a ratio between the magnetic field in the gun solenoid and main solenoid of ≈ 11.5 is recommended. Given an upper limit of the magnetic field in the LHC of 6 T, a possible magnetic configuration of the electron lens at the LHC is 0.43-5-0.43 T. A high main solenoid magnetic field of 5 T is chosen for its magnetic rigidity. Given this compression, the outer beam radius would be at ≈ 3.7 mm or 7.4 σ . Since primary collimators are placed at 6 σ [7], a large portion of the beam is shielded by the primary collimators. It might thus be possible to retract primary collimators up to 7.4 σ , reducing the impedance on the proton beam. In order to obtain a similar scraping effect as was seen at the Tevatron, a cathode voltage of 3-4 kV is recommended, which provides a beam current of 1-1.5 A as can be seen in appendix tab. B.1. Furthermore, tab. 3.2 provides a list of suggested design parameters for the usage of the HG1b at the LHC.

This gun is the focus of this thesis and was used for all subsequent measurements. HG06 was mentioned for comparison purposes and due to its significance in the Tevatron scraping experiments.

Chapter 4

Thermionic Emission

This chapter introduces the concept of thermionic emission in *Temperature Limited Emission Region* (*TLER*) and *Space Charge Limited Emission Region* (*SCLER*). Thermionic emission governs the emission of particles from the cathode in the 1 inch Hollow Electron Gun (HG1b) and thus is directly responsible for the production of the Hollow Electron Beam (HEB). There are three main emission regimes, *TLER* where the beam current strongly depends on the cathode temperature, *SCLER* where the beam current is independent of the cathode temperature and depends on the cathode potential through Child-Langmuir law and Field Emission Regime. The latter does not apply to our guns, since it only arises at very high electric fields of $\approx 10^{10}$ V m⁻¹. Another important parameter for cathode emission is secondary emission, which is induced through the passing of a primary incident particle with sufficient energy through a material. For linear beam tubes only thermionic emission is considered [27, p.39].

4.1 Temperature Limited Emission

With increasing temperature the probability of electrons escaping the cathode surface increases. Thermionic emission refers to the process of emitting electrons from a surface by heating that surface. The minimum energy required for an electron to escape an emitter is the sum of the Fermi energy and the work function, $E_0 + e\varphi$ [27, p. 42]. The Fermi energy is the energy difference between the highest and lowest occupied single-particle states in a quantum system of non-interacting fermions at absolute zero temperature. The work function is energy difference between the top of the conduction band in the cathode and the vacuum level adjacent to the cathode [27]. The minimum energy required to escape an emitter is schematically shown in fig. 4.1.

Electrons obey the Fermi-Dirac distribution, meaning that each energy state may only be occupied by



Figure 4.1: Scheme of the energy levels in a cathode. At higher energy, the disjunct energy bands merge to form the conduction band just below the Fermi Energy E_0 . Particles require at least $E_0 + e\varphi$ to escape the cathode. [27, p. 42]

one single electron. Electrons emitted from the thermionic cathode are part of the Maxwellian tail of the Fermi-Dirac distribution [25, p. 8]. Their current density is described by the **Richardson-Dushman** equation:

$$J_{therm} = A_0 T^2 e^{-\frac{\Phi}{k_b T}} \tag{4.1}$$

where T is the cathode temperature, Φ is the cathode work function and k_b is the Boltzmann constant. $A_0 = \frac{4\pi e m_e k_b^2}{h^3} = 1.2 \times 10^6 \text{ Am}^{-2} K^{-2}$ is a characteristic constant, where e is the elementary charge, m_e is the mass of the electron and h is Planck's constant. The regime in which the emission has the above temperature dependence is called **Temperature Limited Emission Regime (TLER)**. The beam current density is exponentially dependent on the temperature of the cathode.

Given eq. 4.1, when designing a cathode, one tries to minimize the work function in order to obtain a higher yield and possibly be able to operate at lower temperatures. The work function of our cathode and its operating temperature were discussed in sec. 3.2.

4.2 Schottky-Effect

Richardson-Dushman provides a simple relationship between the temperature, work function and current density, but it does not consider electric fields at the surface of the cathode. A potential difference between the anode and cathode causes charges to diffuse inside the cathode in order to return the electric field inside the conductor to equilibrium. This has the effect of lowering the effective work function by:

$$\Delta \Phi = \left(\frac{eE_a}{4\pi\varepsilon_0}\right). \tag{4.2}$$

 E_a is the applied electric field, e is the electron charge and ε_0 is the electric permittivity of free space. We refer to the effective work function, since the work function is usually an intrinsic quality of the surface and material. This effect is called the Schottky Effect [27, p. 46]

Incorporating the Schottky Effect into the Richardson-Dushman equation provides the improved emission equation:

$$J = A_0 T^2 e^{-\frac{e}{k_b T} \left[\varphi - \left(\frac{eE_a}{4\pi\varepsilon_0}\right)^{1/2}\right]} = J_0 e^{\frac{e}{k_b T} \left(\frac{eE_a}{4\pi\varepsilon_0}\right)^{1/2}}$$
(4.3)

Here J_0 is the Richardson-Dushman current density. It is used to compare the performance of different cathodes [27, p. 47].

4.3 Space Charge Limited Emission

Given an emitting cathode, a substantial electron cloud can build-up near the cathode. The potential outside the cathode is thus reduced by the negative charge presence of the electrons. This effect is shown in fig. 4.2. The amount of space charge at the cathode is increased when increasing the temperature in the TLER and decreased when increasing the cathode potential in the SCLER.

The electron emission rate is limited by the point where the space charge is sufficiently large to depress the potential at the cathode below zero, as is the case for (b) in fig. 4.2b. Consequently emitted electrons need to overcome a potential well, allowing only energetic electrons to be emitted from the cathode. Low energy electrons are returned to the cathode by the space charge. This causes the *SCLER* to be lower than the saturated emission rate, which is found exactly when the potential, and therefore the electric field, at the cathode is 0. The saturated emission rate is equal to the Richardson-Dushman emission rate J_0 , mentioned in sec. 4.2. If the potential difference between the diode increases, more electrons are emitted and in turn the potential at the cathode is decreased. If the potential difference decreases, less electrons are emitted, reducing the space charge effect and thus increasing the potential at the cathode. This effect is called the *Space Charge Limited Emission Region (SCLER)*. During



(a) If no space charge is present, the potential between a planar diode changes linearly. Because of the space charge, the gradient of the potential changes shape to that of the dashed line. Source: [27, p. 49].

(b)As the electron emission rate is increased, the reduction of the potential continues until it reaches case (b)where the potential is negative just outside the cathode, pushing electrons back into the cathode. Source: [27, p. 50]

Figure 4.2: Schemes showing the effect of space charge build-up on the potential outside a cathode.

this emission regime, cathode temperature and surface effects no longer play a role in the emission rate [27, p. 50]. This emission regime can be reached by increasing the cathode temperature in the *TLER*, thus increasing the space charge build-up at the cathode. This transition is depicted in figure 4.3b. The space charge limited current density is given by [25, p. 10]:

$$J = 1.67 \times 10^{-3} \left(\frac{q}{mc^2}\right)^{1/2} \frac{V_a^{3/2}}{d^2} \quad [A/m^2]$$
(4.4)

where V_a is the potential difference between anode and cathode, d is the spacing between the cathode and the anode, q and m are the particle charge and mass and c is the speed of light. All constants are given in MKS units [27, p. 51]. From eq. 4.4 the beam current of a cylindrical hollow thermionic dispenser cathode can be determined [25, p. 10].

$$I_{beam} = 1.67 \times 10^{-3} \pi \left(\frac{q}{mc^2}\right)^{1/2} \frac{V_a^{3/2}}{d^2} \left(r_{ext}^2 - r_{int}^2\right) \quad [A]$$
(4.5)

where r_{ext} is the outer radius and r_{int} the inner radius of the cathode. This law is called the **Child-Langmuir law**, from which a quantity, dependent on the electron gun geometry, called the **perveance** is defined as the ratio (distance d is difficult to determine for complex diodes):

$$P = \frac{I}{V_a^{3/2}} = 1.67 \times 10^{-3} \pi \left(\frac{q}{mc^2}\right)^{1/2} \frac{r_{ext}^2 - r_{int}^2}{d^2} \qquad \left[\frac{A}{V^{3/2}}\right]$$
(4.6)

For the purpose of this thesis the unit perv is defined in SI units as $AV^{-\frac{3}{2}}$.

If V_a exceeds a certain level, the space charge condition can not be upheld and the beam enters the temperature limited region, diverging from the Child-Langmuir law. This is depicted in fig. 4.3a.

Considering residual gases in the drift tubes and relativistic effects, the concept of a generalized perveance is introduced [25, p. 196].

$$K = \frac{I}{I_0} \frac{2}{(\beta \gamma)^3} (1 - \gamma^2 f_e)$$
(4.7)

where $I_0 = \frac{4\pi\varepsilon_0 mc^3}{q} = 17kA$ is the characteristic current, I is the beam current, f_e is the neutralization factor and β and γ are the relativistic factors. Unlike the perveance, the generalized perveance is a



(a)In general the current density is given by the child langmuir law (see eq. 4.6). At high potential differences, the space charge condition is broken and the beam enters the Temperature Limited (TL) region again. Source: [27, p. 53]

(b) With increasing temperature, the emission becomes temperature independent. The higher the cathode potential, the higher the temperature at which this transition occurs. Source: [27, p. 53]

dimensionless quantity. Given that $\beta = \frac{v}{c}$ and $v = \left(\frac{2qV}{m}\right)^{\frac{1}{2}}$, we can relate to P by

$$K = P \times \left[\frac{(1 - \gamma^2 f_e)}{4\pi\varepsilon_0 \gamma^3 (2q/m)^{1/2}} \right]$$
(4.8)

Considering a non-relativistic situation equation 4.8 can be simplified to:

$$\frac{K}{(1-f_e)} = P \times \left[\frac{m^{1/2}}{4\pi\varepsilon_0 (2q)^{1/2}}\right] = P \times 1.515 \times 10^4$$
(4.9)

Given a neutralization free beam, the correlation is $K = 1.515 \times 10^4 P$ [25].

4.4 Field Emission

For the sake of completion, we also introduce field emission. At electric fields much higher than that of the SCLER, $\approx 10^9 - 10^{10} \text{ Vm}^{-1}$, electron emission increases rapidly, due to a low work function. Quantum mechanics thus dictates that even particles with insufficient kinetic energy to overcome the barrier by classical theory, have a non-zero probability of passing through the barrier due to the tunneling effect. Such an electron emission is called field emission. This process is schematically shown in fig. 4.4. This type of emission does not play a role in our *Hollow Electron Beam Lenses (HEBLs)*, since we deal with a potential difference of $10^3 - 10^4$ over 1 cm and thus electric fields of $\approx 10^5 - 10^6 \text{ Vm}^{-1}$.

4.5 Summary of Thermionic Emission

In summary, at very low temperatures, the cathode operates in the TLER. When the temperature is high enough, a substantial space charge builds up and the emission regime transitions into SCLER. This is depicted in figure 4.3. The optimal point of operation is given by the green bar and is only considered for voltages up to 8 kV. Figure 4.5 gives a summary of the various emission regions and the respective potential fields needed to achieve them. The emission from the thermionic cathode first enters the SCLER, rising with voltage, until the cathode potential is strong enough in order to transport away the space charge. At this point it enters the *TLER*. Ultimately when the voltage is high enough, field emission takes place, since particles have a non-vanishing probability of tunneling through the work function. The optimal area of operation is again indicated by the green bar.



Figure 4.4: An electron with a wave function whose amplitude is lower than the potential barrier at the surface of the cathode, can tunnel through the barrier given a very strong electric field [27, p. 48].



Figure 4.5: Summary of the various thermionic emission regions with increasing anode voltage. The optimal area of operation is indicated by the green bar. Source: [27, p. 55]

Chapter 5

1 Inch Hollow Electron Gun Characterization

Having discussed the necessary theory, explained the terms *Hollow Electron Beam Collimation (HEBC)* and *Hollow Electron Beam Lens (HEBL)* and described the relevant electron lenses and guns, the measurements at the *Tevatron Electron Lens Test Stand (TELTS)* are presented. First transport measurements are discussed, these measurements describe the amount of beam that is transmitted through the main solenoid and arrives at the collectors. Then the yield measurements are discussed; a discussion of the amount of current produced by the gun. Lastly the method of transverse current distribution profile measurement is discussed and the evolution of the profiles through the beam tube with respect to the magnetic field and the cathode potential is presented.

Transport and yield measurements are done using a Tektronix TDS3034B oscilloscope¹. The cathode and collector current are read using 3 induction coils, 2 dedicated to the collector current which shows a systematic error of $\approx 1\%$. Since all induction coils are the same, this systematic error can be transferred to the cathode current readout. Additionally a statistical error on the oscillator readout of a fraction of a percent is observed $\approx 0.5\%$. Profile measurements are made through the average current in a pinhole in the collector. The transverse profile of the beam is scanned by moving the beam over the pinhole using 6 horizontal and vertical magnetic correctors in the beam tube. Profile measurements are explained in detail in sec. 5.4. The transverse fields are calculated using a simulation code called WARP. This is discussed in detail in sec. 6.

5.1 Operational Procedure of the TELTS

All measurements throughout this thesis were made in the *TELTS*. Turning on the test stand can take 1-2 days, due to the long time it takes to heat the cathode and re-establish the vacuum. Before any parameters are changed on the *TELTS*, the vacuum in the beam tube is measured. The vacuum must remain below 10^{-7} mbar. Next the cathode is heated through a filament. The filament current is gradually increased with long pauses inbetween, so that the vacuum can recuperate. It is desirable to leave the cathode at a lower temperature than the operational temperature of 400 °C in order to evaporate all liquid constituents from the surface. The filament current is then increased to an operational temperature of 9.25 A. Later on we will explain why this should be lowered to 8.5 A. This process can take several days. It is important to keep an eye on the vacuum through *Fermilab Accelerator Control System (ACNET)*.

For future reference, we also mention the cathode temperature dependence on the filament current used to heat it. The scale of temperature to filament resistance is given in [28] as a linear local approximation:

$$T = \beta R + T_0 \tag{5.1}$$

¹http://www.tequipment.net/TektronixTDS3034B.asp



Figure 5.1: Cathode temperature in terms of filament resistance and filament current. [28]

where $\beta = 1765.7 \text{ K} \Omega^{-1}$ and $T_0 = -462.29 \text{ °C}$ and R is the filament resistance. The relation of filament resistance vs. filament current is given in figure 5.1a and the relation of cathode temperature to filament resistance in 5.1b.

Next the Tektronix TDS3034B oscilloscope is turned on. The gun, main and collector solenoids are set to the necessary magnetic fields. The main solenoid uses 2 power supplies that are linked to each other. It is thus sufficient to set the current on one of the power supplies to half that which is used to drive the main solenoid. The other power supply is set automatically. Having set the magnetic fields, the cathode voltage is set and a DEI PVX-4110 10-kV modulator is turned on in order to pulse the electron beam. The pulse width and repetition rate are set. The pulse width is usually 8 μ s and the repetition rate is set in order to obtain an appreciable signal.

During measurements, the pulse from the modulator is measured using 2 induction toroids at the collector, 1 induction toroid at the cathode and the pinhole in the collector of the *TELTS*. The data is recorded through a *Programmable Logic Controller (PLC)* interface for profile measurements. Beam current measurements are done using the oscilloscope.

5.2 Transport Measurements

Transport refers to the migration of particles or energy within a host medium, considering effects such as absorption, emission and scattering. Here the medium is a combination of beam plasma, neutralized particles and a confining magnetic field. We define transmission T as the ratio of beam current at the collector to the cathode:

$$T = \frac{I_{collector}}{I_{cathode}} \tag{5.2}$$

This measure should ideally be 100%. The new 1 inch Hollow Electron Gun (HG1b) is the largest gun that was ever used in the test stand. After a month shut-down of the e-gun in April, profiles taken on 02.05.2013 showed significant asymmetries (see fig. 5.2). Bad conditioning of the cathode surface was suspected. Following 11 additional days of conditioning at 9.25 A and 10.0 A, the profiles only improved slightly.



Figure 5.2: The figures show the normalized current density in Acm^{-2} of the measured profiles in the transverse plane. Current densities below $6.25 \times 10^{-4} Acm^{-2}$ are negligible and ignored. Profile (a) shows the electron beam distribution on 02.05.2013. It is strongly asymmetric. After 11 more days of conditioning, the asymmetries have not improved as seen in (b). The production of these plots is explained in detail in section 5.4

During the follow-up a discrepancy between collector and cathode current was noted. The cathode and collector current were 68.6 mA and 50.6 mA respectively, giving a transmission of 73.8%. The cathode potential was 500 V and the magnetic field was 0.3 T for this measurement. Causes for the low transmission can either be particle-particle interactions, such as neutralization, or scraping at the beam pipe. To lower the effect of scraping, the beam can be compressed in the main solenoid, and to reduce neutralization, a better vacuum could be provided.

	B_{gun} [kG]	B_{main} [kG]	$B_{collector}$ [kG]
	1	1	1
1.5 Pre 2		1.5	1.5
		2	2
	2.5	2.5	2.5
	3	3	3
	0.4	1.6	0.4
	0.6	2.4	0.6
Post	0.8	3.2	0.8
	1.0	4.0	1.0

Table 5.1: Pre and post transmission improvementsolenoidal magnetic field settings

Trying several combinations of B_{gun} , B_{main} , B_{coll} , it was found that transport is best with $B_{gun} = B_{collector} = 0.25B_{main}$. This provides a transport of $\approx 100\%$. Table 5.1 shows the improved solenoid settings that were used. Using the older 0.6 inch Hollow Electron Gun (HG06), such transmission issues were not observed and the gun could be used under equal magnetic fields ($B_{gun} = B_{main} = B_{coll}$). Because the improvement in transmission for the HG1b was achieved by compressing the beam, the most likely cause of the losses of current using equal magnetic fields is beam scraping.

5.3 Yield Measurements

Yield measurements are made by increasing the potential difference between the cathode and anode in regular steps and using the Tektronix oscilloscope to measure the yield at the collector. The oscilloscope averages results taken during a few seconds, reducing the statistical error significantly to a fraction of a percent. The systematic error is given by about 1%.

The yield at the collector is defined by the acceleration potential V_a , the perveance P and the transmission in the beam tube T as:

$$Y = P * V_a^{3/2} * T (5.3)$$



Measurement Date	Perveance $[\times 10^{-6} \text{ A/V}^{3/2}]$
23.05.2013 15:02	5.36 ± 0.03
22.05.2013	5.27 ± 0.04
21.05.2013	5.35 ± 0.03
Expectation	5.33 ± 0.14
07.05.2013	4.20 ± 0.01
20.02.2013	4.23 ± 0.01
14.12.2012	4.22 ± 0.01
Expectation	4.22 ± 0.03

Figure 5.3: Effect of transport improvement on the yield of the test-bench. The yield follows the Child-Langmuir law and measurements are thus taken in SCLER. The statistical error is estimated at 0.5%. Systematic errors are not considered, because they affect each measuring point equally.

Table 5.2: Measured global perveances and their expectation values. The new measurements are $\approx 26\%$ higher than the old perveance measurements, prior to May 2013. Error estimates are given through the statistical uncertainty and calculated by the fit.

We define the global perveance to be the fit between all data points of the peak collector current and the cathode potential according to the model in eq. 4.6; $I \propto V^{\frac{3}{2}}$. The peak collector current is the highest current obtained from the pulsed beam shape. The pulse shape is shown in [28, p. 13]. This is an overall average which is independent of the cathode potential. In addition to the global perveance we define a quantity called the local perveance which describes the fraction of $I_{peak}/V^{\frac{3}{2}}$ for each measurement point.

5.3.1 Global Perveance

Fig. 5.3 compares the global perveance of measurements made after the transport improvements², $5.33 \pm 0.14 \ \mu$ perv, and those made before the improvement, $4.22 \pm 0.03 \ \mu$ perv. Measurements prior to the transport improvement confirm measurements done earlier by Siqi Li [28]. Li obtained a perveance of $4.13 \pm 0.01 \ \mu$ perv in February 2012 and $4.15 \pm 0.02 \ \mu$ perv during the measurements of June 2012. The gun is relatively stable over long time periods (see fig. 5.3). All global perveance measurements after the asymmetric profiles, experienced beginning of May 2013, seem to show a slight deviation from the Child-Langmuir law. The significance of the transmission improvement is apparent through a yield increase of $\approx 26\%$.

5.3.2 Global Perveance from SAM Simulations

SAM is a code developed at Budker Institute of Nuclear Physics for simulations of stationary axialsymmetric electron optical systems and electron guns. The code distributes a surface charge across electrodes and dielectrics, which are simulated using line and arc segments. A curved mesh is used to describe the space charge density which is constant within one cell. Further information on this code can be found in [29].

²Measurements continued by: G. Stancari, A. Valishev and M. Chung during the days after my departure from Fermialb.



Magnetic Field [T]	Perveance $[\times 10^{-6} \text{ A/V}^{3/2}]$
0.05	5.796 ± 0.079
0.1	6.322 ± 0.057
0.2	6.381 ± 0.051
0.32	6.371 ± 0.051
0.5	6.337 ± 0.060
Expectation	6.321 ± 0.110

Figure 5.4: Perveance plots for the SAM simulations at 5 different solenoidal field strengths. [30]

Table 5.3: SAM simulated perveance and expectation value of $6.3(1) \times 10^{-6} \text{ AV}^{-\frac{3}{2}}$.

In January 2012, using SAM, Leonid Vorobiev at *Fermi National Accelerator Lab (Fermilab)* simulated the electron gun current for a given acceleration potential at a magnetic field of 0.05, 0.1, 0.2, 0.32 and 0.5 T in the gun solenoid. From this, we calculated the global perveance of the electron gun, depicted in fig. 5.4 and table 5.3. Our measurements were within 15% of the SAM simulations and satisfactory.

The expectation value of $6.3(1) \times 10^{-6}$ perv includes all magnetic fields except that of 0.05 T, which is insufficient to force the electrons to go on a straight path. The transverse electric fields between the cathode and anode cause the particles to undergo strong transverse oscillations, called scalloping [27, p. 138]. This results in the partial scraping of the electron beam and reduces the yield. The scalloping can be observed in fig. 5.5, where (a) is the electron beam at a solenoidal field of 0.05 T and (b) is at 0.32 T. In the stronger field, the magnetic confining force is too strong to allow scalloping.

5.3.3 Local Perveance

Following the improvement of the transport mentioned in sec. 5.2, yield measurements were taken on 14.05.2013. Fig. 5.6a shows the electron gun peak current [A] as a function of the cathode potential [V]. Fitting a line to all data points (red and blue) provides us with a global perveance of $5.12 \pm 0.03 \mu$ perv. Without a discrepancy from the Child-Langmuir law, this graph should portray a straight line. This is the case up to 4000 V. At higher voltages a gradual deviation from the Child-Langmuir law can be observed. A fit of the data up to 4000 V provides us with a perveance of $5.70 \pm 0.02 \mu$ perv, $\approx 90\%$ of the SAM simulations perveance. The discrepancy between the Child-Langmuir law and the measurement is portrayed through a local perveance plot with respect to the cathode potential in fig. 5.6b.

The yield is measured at the end of the beam tube. Therefore it can be negatively affected by neutralization of the beam by residual gases in the beam tube, scraping of the beam at the beam tube edge and a departure from the *SCLER* to the *Temperature Limited Emission Region (TLER)*.

When the potential is high enough, space charge effects occur further downstream, decreasing the space charge close to the gun. Consequently the SCLER cannot be sustained. Under these conditions, the emission returns to the TLER. This effect is shown in fig. 4.5 and 4.3a. Sec. 5.3.4 tells us that the voltage at which the SCLER transitions into the TLER, rises with increasing filament current. Seeing that this break down has not occurred for a filament current of 8.25 A in fig. 5.8c, the drop in yield



Figure 5.5: SAM simulations of HG1b from January 2012 [30]. The blue lines represent the electron trajectories while the green lines represent the electric field at the cathode.



Figure 5.6: Global (a) and local (b) perveance measurements of the HG1b. A discrepancy was seen in fig. (a) where the perveance departed from the Child Langmuir law at 4 kV. This is most likely due to beam scraping as explained in the text. Fig. (b) shows that the change of local perveance with voltage due to this discrepancy is linear.

seen in fig. 5.6a is probably not due to the transition into the TLER.

Neutralization refers to the absorption of electrons by background gases. There are many reasons for background gases [27, p. 173]. In the *TELTS*, the predominant background gas is Hydrogen at 1×10^{-8} mbar. Fig. 5.7 shows the number of ions of various gases formed per $cm \times mbar$ for electron-gas interactions. As can be seen, the cross section drops to values under unity when the electron energy exceeds 10^3 eV. Assuming a pressure of 1×10^{-8} mbar, at 5000 V only 5×10^{-8} ions are formed per cm. We can thus safely assume that neutralization only plays a role for low energy electrons, below 4000 V. This was also confirmed in Siqi's paper [28].

The sudden change of slope between 4 kV and 5 kV in fig. 5.6a is thus likely explained by the scraping of the beam at the beam tube with a radius of 3 cm. The effect thereof seems to be a linear change of perveance with voltage after 4 kV



Figure 5.7: Number of ions formed per cm per mbar for various gases. Source: [31, p. 104]

with a slope of $-1.27(4) \times 10^{-10}$ perv V⁻¹, as seen in fig. 5.6b. The linear change could be explained by the beam reaching a transverse beam size at $V_a=4$ kV such that the beam starts scraping on the beam tube edge. As the acceleration potential V_a is increased, more current is produced and the scraping increases, further lowering the perveance.

Assuming that the decrease in yield is mostly due to scraping, the fit for the whole data set (green) in fig. 5.6a is the best assumption for the yield at the collector. This gives a perveance of $(5.12\pm0.03) \mu$ perv. The fit for the set of lower data best describes the perveance of the gun itself, which excludes losses downstream of the gun.

5.3.4 Dependence of Perveance on Filament Current

This section discusses the filament current dependence of the perveance. A scaling law between filament current and temperature of the cathode is given in eq. 5.1.

Additional measurements were made at a magnet setting of 1-4-1 kG for linearly changing filament currents from 6.25 A to 9.75 A in steps of 0.5 A. Using this data, we expect to be able to validate our current measuring point of 9.25 A and possibly deduce a lower filament operating current, without a significant loss in yield due to the temperature-limited regime. The cathode yield was acquired at a repetition rate of 4 Hz. Fig. 5.8 shows the measurement results.

Fig. 5.8a shows the transition from SCLER to TLER and the voltage at which this transition occurs. In the TLER, the emission current is independent from the cathode potential, when the emission has entered the TLER completely. The transition does not take place below 10 kV for a filament current between 8.25 A and 9.7 A. Interestingly, the emission current seems to drop again at higher cathode potentials as one surpasses 8.75 A of filament current. This effect is analyzed more closely in fig. 5.9a.

Fig. 5.8b shows the electron gun current as a function of the filament current. Considering measurements on the full potential voltage scope (0.1-10 kV), the maximum "knee current" is at 8.25 A. The "knee current" is given by the filament current at which the transition into the *SCLER* takes place. If we consider a safety filament current increment of 0.5 A, an operating filament current of at least 8.75 A is recommended. If one wishes to only measure profiles with a cathode potential of up to 7 kV, this recommended filament operating current can be lowered to 8.25 A and for up to 4 kV to 7.75 A. The filament current increment is used, because the yield tends to fall with time as described in [27, p. 73]. By taking a safety filament current increment, we guarantee that the current density produced by the cathode does not fall and measured profiles are consistent.



(a)Peak current vs. cathode potential for various fila- (b)Peak current vs. filament current for various cathode potentials.



(c)Perveance vs. cathode potential for various filament (d)Perveance vs. filament current for various cathode currents. potentials.

Figure 5.8: Electron gun peak current (top graphs) and perveance (bottom graphs) as a function of acceleration potential and filament current. All measurements were taken at 1-4-1 kG at a repetition rate of 4 Hz.


(a) Close-up of peak current vs. cathode potential for the highest filament currents.

(b)Close-up of perveance vs. cathode potential for various filament currents.

Figure 5.9: Close up of the cathode-potential dependence of the electron gun peak current and the perveance. The left plot shows that the yield is actually lower for the higher temperature cathode with a filament current of 9.75 A. The right plot, allows one to deduce what temperature to use at given cathode voltages. Below 6 kV, the yield is greatest for the highest filament current of 9.25 A. Above 6 kV, the yield is highest for a filament current of 8 to 9 A.

Fig. 5.8c, similar to fig. 5.8a, shows the transition into the TLER and the drop in perveance. This plot confirms that the local perveance drops slightly at high cathode potentials when exceeding 8.75 A filament current and that the cathode should be operated above 8 A. Fig. 5.8d shows the transition of the perveance from TLER to SCLER.

Fig. 5.9 shows that above a given filament current, the yield is lower than for smaller filament currents. The perveance at 9.75 A falls below that of 8.75 A and 8.25 A at 6 kV and the perveance at 9.25 A falls below that of 8.75 A and 8.25 A at 3 kV as seen in fig. 5.9b. Running the cathode at a too high temperature will deplete the BaO in the porous tungsten faster and decrease the lifetime of the cathode [27, p. 63].

The above analysis proposes that one operates the cathode not at 9.25 A, as done up to now, but at a lower filament current of ≈ 8.5 A.

5.3.5 Generalized Perveance

The generalized perveance provides a dimensionless unit used to compare different cathodes (sec. 4.3). In order to calculate this quantity, we first need to assert that we can operate in non-relativistic assumptions. Given the simple relation $\gamma = \frac{E}{E_0}$ one can derive a relation for the velocity of a particle having passed the anode.

$$\sqrt{\frac{1}{1-\beta^2}} = \gamma = \frac{E_0 + qV_a}{E_0} \Rightarrow \beta = \sqrt{1 - \left(\frac{E_0}{E_0 + qV_a}\right)^2} \tag{5.4}$$

 E_0 is the rest energy of the electron, γ and β are relativistic parameters, q is the electron charge, V_a is the acceleration potential and E is the total energy of the electron, which is given by the acceleration potential. The electron velocity is given by βc , where c is the speed of light. Using conservative values for V_a of 8000 V, which is the maximum we have measured, we obtain a β of 0.17 and can thus proceed by calculating the general perveance in a non-relativistic manner.

In order to determine the general perveance (see eq. 4.9), we would plot the right hand side of that equation vs. the voltage and see that the right hand side equals the general perveance when the neutralization is 0. Sigi Li's paper [28, p. 15] states that the neutralization approaches 0 at very high acceleration potentials. An upper limit for the generalized perveance is determined by utilizing the local perveances at 10 kV from fig. 5.9b, where neutralization is clearly 0. Averaging these values gives a local perveance of $5.14(5) \times 10^{-6}$ perv. Using eq. 4.9 the generalized perveance is $7.93(7) \times 10^{-2}$. This quantity will be important when comparing the measurements to the numerical simulations in chap. 7.

5.4 Profile Measurements

Transverse profiles are measured in order to analyze the transverse beam distribution and any azimuthal asymmetries in the transverse profile. They are subsequently used to determine the transverse electric fields of the measured profiles and an upper limit for the emittance growth due to the *HEBL*.

5.4.1 Transverse Profile Measurement

Transverse profile measurements are acquired using 6 horizontal and 6 vertical beam correctors that deflect the beam in the transverse plane. A 0.2 mm pinhole at the center of the collector measures the current density as the beam is moved along a grid, defined in the acquisition script (ACL). The readout is then saved as a text file and processed by a R Script. This chapter discusses the profile measurement method and the profile evolution as it travels through the beam tube. Additionally a scaling law is deduced and discussed.

ACL Script

The full Accelerator Command Language (ACL) script is provided in appendix C.1.1.

An ACL script [32] controls the TELTSthrough the ACNET for profile scans. After initializing the run, the script defines the list of horizontal and vertical correctors that are used and the variables that are to be read (Code Line: 18) For future reference the most important variables are listed in table 5.4. It then defines the extension of the grid over which the scan is to be performed. This grid size is given in units of ampere. It then defines a quantity called *step* which defines the step size made in each direction. The

Variable Name	Variable Descriptor
Z:BC1AD3	probe current
Z:BC1AD5	tube current
Z:BC2AD5	diode pump vacuum
Z:HCORxx	horizontal corrector positions
Z:VCORxx	vertical corrector positions

Table 5.4: List of variables recorded by ACNET and saved during the transverse profile measurements.

total size in horizontal or vertical direction of the grid has to be an integer multiple of the step size. The variable dt defines how much time is spent per measurement point. For post transmission upgrade measurements a 31×31 pixels with 2 s interval per pixel were used. Pre-transmission upgrade measurements used a 61×61 pixel grid with a 1 s interval per pixel. The change was made purely for the purpose of reducing measurement time. Table 5.5a shows the optimal grid sizes in A for the various magnetic field configurations. Additionally tab. 5.5b gives the correction parameters for the grid not being centered around (0,0) A.

The script first runs a few safety calculations (Code Line: 55) to see that all entries are correct and physical. After this it proceeds by calculating the expected runtime. For our measurements this

	-	
Solenoidal	Horizontal	Vertical
configuration	settings [A]	settings [A]
[T]		
1.0-1.0-1.0	[-1.5, 1.5]	[-1.5, 1.5]
1.5 - 1.5 - 1.5	[-2.25, 2.25]	[-2.25, 2.25]
2.0-2.0-2.0	[-3,3]	[-3,3]
2.5 - 2.5 - 2.5	[-4,3.5]	[-3.75, 3.75]
3.0-3.0-3.0	[-4.8,4.2]	[-4.5, 4.5]
0.4-1.6-0.4	[-1.8,0.6]	[-1.4,1.0]
0.6-2.4-0.6	[-2.8,0.9]	[-2.1, 1.5]
0.8-3.2-0.8	[-3.6, 1.2]	[-2.8, 2.0]
1.0-4.0-1.0	[-4.5, 1.5]	[-3.5, 2.5]

Solenoid con-	Horiz. correc-	Vert. correc-
figuration [T]	tors [A]	tors [A]
0.4-1.6-0.4	-0.645	-0.2
0.6-2.4-0.6	-0.968	-0.3
0.8-3.2-0.8	-1.290	-0.4
1.0-4.0-1.0	-1.613	-0.5

 $(a) Corrector\ settings\ for\ grid\ scans.$

(b) Corrector center of mass settings.

Table 5.5: The corrector settings in the beam tube and center of mass of the measured profiles for the different magnetic field configurations $(B_{gun} - B_{main} - B_{coll})$.

was usually of the order of one hour. Lastly the script defines a quantity called **stepsign** (Code Lines: 70-73) which gives the direction of the scan. The script scans one row after the other in the vertical direction, changing scan direction for each row. The directionality of the scan is important and can cause artifacts in the transverse scan profile.

The script then terminates by resetting all the correctors to 0 (Code Lines: 130-132). The measured data is saved in a text file. Code 5.1 is an example of such a text file, for a measurement that was taken on 14.05.13 at a filament current of 9.25 A, magnetic field of 0.8-3.2-0.8 kG, a cathode potential of 500 V and a peak current of 73 mA.

```
#=== Electron lens test stand profile scan ===
  # 14-MAY-2013 17:47:54
  # Scan parameters:
  #
      mix = -3.6
                     A
  #
      max = 1.2
                     A
  #
      miy
          = -2.8
                     A
           = 2
  #
      may
                     A
 8 #
      step = .16
                     A
  #
      dt
           = 2
9
                     s
10
11
12 # xset[A] yset[A] Z:BC1AD3[UA] Z:BC1AD5[UA] Z:BC2AD5[MA] Z:HCOR02[A] ... Z:VCORB5[A] ctime[s]
13
    -3.6
            -2.8
                   .0012207031 .0001953125 6.51550293 -3.603515625 ... -2.80090332 1368553731
14
                   .0014648437 .0001953125 6.572570801 1.192626953 ... 1.995239258 1368555846
     1.2
             2
                              Code 5.1: "Sample of profile measurement text file"
```

The text file first contains a preamble that lists the settings with which the ACL script was run. In this case we used a 31 x 31 grid with a 2 s measurement per grid point. Following the preamble, the measurements are listed. First the x and y position are given. Followed by the probe current, the tube current and the diode pump vacuum. It then lists the positions of all 6 horizontal and 6 vertical correctors. Lastly it provides the time stamp of the measurement.

R Script

R is an open-source GNU's Not Unix (GNU) language for statistical computing and graphical output. It is based on the similar S language and environment which was developed by Bell Laboratories³.

³www.r-project.org

The R script starts of with the usual initialization (Code Lines: 8-12), such as the loading of library files, defining very long vectors of all input files, their beam parameters and their transverse offsets. Furthermore some processing and plotting options are set. The script uses two types of color coding in contour plots. One is the classical MATLAB color series with blue being the lowest value and red the highest. It was found that this color scheme provided too little contrast for the areas of interest. A second scheme named COLOR.CHART was created to provide more color contrast in the mid-range (Code Line: 175).

The script calculates a few important global quantities such as the electron frequency, flight time, cyclotron and plasma frequency and evolution number (Code Lines: 321-339). These are printed to the standard output. The R script can also handle measurements made with previous versions of the above mentioned *ACL* script. The new *ACL* script provides vacuum data. The R script thus extracts the data if necessary. The measured data is then stored into 3 arrays: **Zprobe**, **Ztube** and **Zvac** (Code Lines: 421-423). The quantity that is ultimately plotted is **Zprobe**. The background of the measurement is calculated by averaging the fourth lowest quantile of the measurements. It is then subtracted from the measurements are normalized and the center of mass of the profiles are calculated. At this point, the transverse grid is still given in units of amperes, which are provided to the correctors to deviate the beam. To convert them to mm, either the calibration coefficients saved in calx.v and caly.v are used, or they can be recalculated by turning on the option CALIBRATE.CORRECTORS (Code Line: 564). This will ask the user to define the outer and inner beam by defining 8 points on each edge. Given the current deviation and the normal beam size, the calibrations coefficients are calculated and printed to the standard output.

The major calculations have now been completed and the script creates several plots of the measurements (Code Line: 681). First a 1D cross-section of the beam, showing the variation of current density along the horizontal direction with the vertical position set to 0. It then produces 2D contour plots (Code Line: 720). These come in two variations. One with annotations and one without. Lastly it produces a 3D plot of the profile (Code Line 800).

Having produced all the plots, the script, if the respective options are turned on, outputs; the calibrated profile and a radial plot of current density vs radius from the beam center, which is the average over all azimuthal angles, a particle plot together with a text file containing all the coordinates and if selected a polar decomposition plot (Code Line: 931). The full R Script is provided in app. C.1.2.

Measured Transverse Profiles

Profiles were measured throughout the course of this thesis in the *TELTS*. For a list of all measured profiles, see appendix B.1. Profiles can be sorted by the magnetic fields in the three big solenoids and the cathode voltage. The cathode current is dependent on the cathode voltage through the Child-Langmuir law. In general it was observed that profiles were reproducible when measured with the same parameters. Considering space charge evolution and the angular rotation of the profiles, a chart of all profiles with respect to the magnetic field in the main solenoid and the cathode potential was made. This allows us to uniquely observe profile evolution through these two parameters. Fig. 5.10 depicts such a chart with profiles made after the transmission upgrade mentioned in sec. 5.2. The red lines indicate the expected scaling laws for a constant ratio of pinhole current to total current, which is discussed in section 5.4.3. We can compare this to the profiles that were measured before the transmission upgrade had a magnetic configuration of $B_{gun} = B_{coll} = 0.25B_{main}$ while measurements before the transmission upgrade had equal magnetic fields in all 3 solenoids. The lack of magnetic compression in the main solenoid has prevented the rapid space charge evolution seen in fig. 5.10 in the measurements made before the transmission upgrade shown in fig. 5.11.

The profiles that are of interest to us are those in the top right corner. Profiles below 0.5 kV tend to become unstable, due to insufficient removal of space charge from the surface of the cathode. At the LHC, the maximum magnetic confinement will be of the order of 6 T, which is the easy attainable







Figure 5.12: Beam profile at $B=1-4-1 \ kG$, $V=500 \ V$ and $I_{peak}=73 \ mA$. The beam shows a good symmetry, being squeezed a little in the horizontal direction. fig. (b) shows slight slippage through two radially opposing centers. The radial profile in fig. (c) shows a smooth distribution, indicating that the beam is radially similarly distributed for all azimuthal angles.

magnetic field with today's superconducting magnets. Given this and the assumption that the red lines in fig. 5.10 hold, we see that the profiles in the top right corner are those that represent the profiles that will ultimately scale to LHC parameters. The cathode potential at the LHC will most likely be operated around 4-5 kV. Fig. 5.12 shows the 2D, 3D and radial profile of an optimal measurement point at B=1-4-1 kG and V=500 V.

5.4.2 Beam Evolution

This section derives an equation for the angular rotation velocity of the transverse profiles. This is done as a sanity check for the evolutions observed in figures 5.10, 5.13 and 5.14.

The e-beam is essentially a non-neutral plasma that exhibits collective behavior. In order for plasmas to be in equilibrium, they always require a restoring (space charge) and a disturbing self-field force. The radial motion is given by (see: [25, p. 175])

$$\ddot{r} = \frac{\omega_p^2}{2}r \qquad \omega_p^2 = \frac{q^2n}{\varepsilon_0\gamma^3m} \left(1 - \gamma^2 f_e\right) \tag{5.5}$$

Here r is the transverse position of the profile, ω_p is the plasma frequency, q is the elementary charge, n is the particle density, ε_0 is the permittivity of vacuum, γ is the relativistic factor and f_e is the neutralization factor. As discussed earlier, we can assume non-relativity and thus set $\gamma = 1$.

In the case of our electron beam, the restoring force is the $E \times B$ twist, where the magnetic field B is applied externally through the solenoids and the electric field E is radially outward, given by the self fields of the space charge in a 2D approximation. Equilibrium between these two forces is given when the beam is matched (beam envelope on straight line). The beam is matched for a oscillation frequency ω due to the applied focusing force and the space charge force of (see: [25, p. 191]):

$$\omega^2 = \omega_L^2 - \frac{\omega_p^2}{2} = \left(\frac{eB}{2m_e}\right)^2 - \frac{ne^2}{2\varepsilon_0 m_e}$$
(5.6)

 ω_L is the larmor frequency, $\frac{qB}{2m_e}$. Due to the radial oscillations, particles have a non-zero transverse velocity and thus exhibit larmor oscillations in the axial magnetic field, causing the beam to rotate. The angular rotation velocity ω_r is given by the dioctron frequency ω_D [33, p. 298]:

$$\omega_r = \omega_D = \frac{\omega_p^2}{2\omega_c} = \frac{ne}{2\varepsilon_0 B} \tag{5.7}$$

where ω_c is the cyclotron frequency which is 2 times the larmor frequency. For a full cylindrical electron beam the angular rotation velocity ω_r is independent of the radial position of the particle because the radial electric field rises linearly with the electric field. Considering a *Hollow Electron Beam (HEB)* configuration and including neutralization f and a contribution due to a rotation of the beam induced by the axial electron current and the azimuthal self-magnetic field β_z^2 , we obtain the radially dependent azimuthal rotation velocity [33, p. 308]:

$$\omega_r = \omega_D (1 - f - \beta_z^2) \left[1 - \left(\frac{R_i}{r}\right)^2 \right], \quad \forall R_i \le r \le R_o$$
(5.8)

where R_i is the inner beam radius, R_o is the outer beam radius and r is the particle radius. We assume neutralization to be negligible [28, p. 15], f = 0, and neglect self-magnetic fields, $\beta_z^2 = 0$. It is thus possible to deduce the angle through which the radial profile layers evolve by multiplying with the time needed to transverse the electron beam tube. The distance between the cathode and the collector L is 2.86 m in the test stand. Assuming that the particle velocity v_z is constant throughout the beam tube, we can give the angle or rotation by:

$$\varphi_r = \frac{\omega_D L}{v_z} \left[1 - \left(\frac{R_i}{r}\right)^2 \right] = \frac{n_e e L}{2\varepsilon_0 B v_z} \left[1 - \left(\frac{R_i}{r}\right)^2 \right], \quad \forall R_i \le r \le R_o$$
(5.9)

Given electrons, the number density of electrons n_e and the transverse charge density ρ of electrons are related by $n_e=6.24\times10^{18} \rho \,\mathrm{L}^{-1}$, where $\rho = \frac{I}{v_z \pi (R_o^2 - R_i^2)}$. Furthermore using equation 5.4, one obtains:

$$\varphi_r = \frac{IL}{2\pi\varepsilon_0 B(R_0^2 - R_i^2) v_z^2} \left[1 - \left(\frac{R_i}{r}\right)^2 \right]$$
(5.10)

$$=\frac{IL}{2\pi\varepsilon_0 B(R_0^2-R_i^2)c^2eV_a\left(\frac{(2E_0+eV_a)}{(E_0+eV_a)^2}\right)}\left[1-\left(\frac{R_i}{r}\right)^2\right],\quad\forall R_i\leq r\leq R_o\tag{5.11}$$

where c is the speed of light, e the electron charge and V_a the acceleration potential between the anode and the cathode. Given that $\frac{eV_a}{E_0} \ll 1$, we can approximate $\frac{2E_0+eV_a}{(E_0+eV_a)^2} \approx \frac{2}{E_0}$ and thus write:

$$\varphi = \frac{IE_0L}{2\pi\varepsilon_0 B(R_0^2 - R_i^2)2c^2 eV_a} \left[1 - \left(\frac{R_i}{r}\right)^2 \right], \quad \forall R_i \le r \le R_o$$
(5.12)

Having a dependency upon B, V_a , I and r at this point, one can eliminate the dependency on I through the Child-Langmuir law. As perveance, the measured perveance of the test stand in fig. 5.2 of 5.33×10^{-6} perv was used. Furthermore $E_0=8.19\times10^{-14}$ J, $\varepsilon_0=8.85\times10^{-12}$ A² s⁴ m⁻³ kg⁻¹, $c=3.00\times10^8$ m s⁻¹ and $e=1.60\times10^{-19}$ C. The inner and outer radius must be taken as an average of the beam radii before, during and after the magnetic compression. Given a magnetic compression of a factor 2, the average beam radii are 0.63 times the cathode beam radii. This gives an average outer beam radius of 8.0 mm and an inner radius of 4.3 mm. Using these values we can thus write:

$$\varphi = \frac{PE_0L}{4\pi\varepsilon_0 c^2 (R_0^2 - R_i^2)e} \times \frac{\sqrt{V_a}}{B} \left[1 - \left(\frac{R_i}{r}\right)^2 \right]$$
(5.13)

$$= 2.7 \times 10^{-2} \frac{\sqrt{V_a}}{B} \left[1 - \left(\frac{R_i}{r}\right)^2 \right], \quad \forall R_i \le r \le R_o$$
(5.14)

If we consider particles halfway between the average inner and outer cathode radius, r=6.15 mm, the angular rotation velocity is given by:

$$6.15\varphi = 1.4 \times 10^{-2} \frac{\sqrt{V_a}}{B}, \quad \forall R_i \le r \le R_o \tag{5.15}$$

Angle of rotation [°]		Acceleration Potential [kV]								
Magnetic Field [T]	1	2	3	4	5	6	7	8		
0.40	72	101	124	143	160	175	189	202		
0.32	89	127	155	179	200	219	237	253		
0.24	120	169	207	239	267	292	316	337		
0.16	179	253	310	358	400	438	473	506		

Table 5.6: Transverse rotation angles for particles at outer cathode amplitude as a dependence of acceleration potential V_a and magnetic field strength in the main solenoid B_{main}

Table 5.6 gives a list of all angles of rotation in radians due to the dioctron profile evolution for profiles with an acceleration potential of $V_a=1$ kV to 8 kV and $B_{main}=0.16$ T to 0.4 T. We can use this equation as a sanity check for the beam evolution and beam profiles we have measured. It is important that B in eq. 5.15 is the average magnetic field throughout the beam tube, which is 0.87 times the magnetic field in the main solenoid.

Evolution as a function of V

The beam evolution is portrayed in fig. 5.13 for a constant magnetic field of B = 0.32 T. With increasing cathode potential, the beam diverges from a uniform distribution (V = 500 V) to a two-island shape (V = 3 kV) and eventually into one major peak (V = 8 kV). The radial profiles portray an average over all azimuthal angles of the beam profile. At 125 V and 250 V a clear beam edge can be seen at a radius of about 3 mm. As the cathode potential increases, this blurs out. At low voltages, the beam seems to be very stable and axis-symmetrical. As expected, the area underneath the radial profile increases, signalizing a higher beam current. The beam profiles rotate in fig. 5.13 with a radially dependent frequency, indicated by the change in beam profile with changing V. Given that $\omega_0 \propto \sqrt{V_a}$, doubling the cathode potential should make the profiles evolve 1.4 times as fast. According to table 5.6, going from 2 kV to 3 kV causes a rotation of 28°. This can be qualitatively confirmed.

Evolution as a function B

Fig. 5.14 shows the scaling of the transverse profiles with constant cathode potential. Given that $\omega_0 \propto B^{-1}$, a change of the magnetic field from 3.2 kG to 4 kG means that the evolution should be slower by 20%. According to table 5.6, transitioning from $B_{main}=3.2$ kG to 2.4 kG should make the profile turn by 42°. This seems to be the case.

5.4.3 Beam Scaling

Considering tab. 5.6 and fig. 5.10, we can assume that profiles will be frozen at magnetic fields of a few Tesla. Still we attempt a first quantitative analysis of dynamical similarity between the profiles, by finding an approximate scaling law for equal profiles with respect to B and V.

In 1914, Mr. E. Buckingham published a paper on dimensional analysis that is still cited regularly today (see [34]). He discusses several techniques that allow one to draw fundamental, but necessary conclusions just by considering the dimensional units of equations.

Assuming the necessity to describe a system solely based on n physically different quantities, a physically complete relation can be formulated as

$$f(Q_1, Q_2, \dots, Q_n, r^1, r^2, \dots) = 0$$
(5.16)

where r^1, r^2, \ldots signify a number of ratios between dimensionally equal quantities. Assuming that these remain constant, one can reduce the above to

$$f(Q_1, Q_2, \dots, Q_n) = 0 \tag{5.17}$$





Figure 5.14: Transverse profile series at V = 2 kV.

which describes the complete system and is called a physical equation. Each such equation can be expressed in the more specific form

$$\sum MQ_1^{b_1}Q_2^{b_2}\dots Q_n^{b_n} = 0 \tag{5.18}$$

where terms such as logarithms and sin functions have been replaced by their respective expansions. If we now divide through by any one term, this becomes

$$\sum MQ_1^{a_1}Q_2^{a_2}\dots Q_n^{a_n} + 1 = 0.$$
(5.19)

Each term in this equation must have the same dimension and therefore, one can write the dimensional equations

$$[Q_1^{a_1}Q_2^{a_2}\dots Q_n^{a_n}] = [1].$$
(5.20)

By defining the dimensionless products $\Pi = Q_1^{a_1}, Q_2^{a_2}, \ldots, Q_n^{a_n}$ we can rewrite eq. 5.19 as

$$\sum \mathbb{N}\Pi + 1 = 0. \tag{5.21}$$

Considering that Π is dimensionless, we can rewrite eq. 5.21 as

$$\sum \mathbb{N}\Pi_1^{x_1} \dots \Pi_i^{x_i} + 1 = 0 \tag{5.22}$$

and thus for every complete physical equation of the form 5.16 we can write a relation in terms of the n - k dimensionless quantities Π , combined from the n physical quantities Q_1, Q_2, \ldots, Q_n and the k fundamental units entailed in theses physical quantities, as

$$\Psi(\Pi_1, \Pi_2, \dots, \Pi_{n-k}) = 0 \tag{5.23}$$

$$[\Pi_1] = \dots = [\Pi_{n-k}] = [1] \tag{5.24}$$

In order to obtain a perception for the profiles at higher magnetic fields, as used at the LHC, a scaling law for a constant profile

ing law for a constant profile evolution was derived using a reduced dimensional analysis, where the permittivity of free space ε_0 was omitted. Constant profile evolution requires that the ratio of $I_{\text{pinhole}}I_{\text{total}}$ remains constant. Given the assumption that we operate in the space charge limited regime, we assume Child-Langmuir's law, eq. 4.6, and thus consider the following 5 independent variables

boundary conditions due to the conductors were neglected.

Given 5 independent physical quantities and 4 SI units, one can define 1 dimensionless parameter that describes the system through the equation $(1 + 1)^{1/2}$

$$\Psi(\Pi) = R,\tag{5.25}$$

where $R = I_{\text{pinhole}}/I_{\text{total}}$. We chose the 4 fundamental units not to be the SI units, but the magnetic field B, the cathode potential V, the system length L and the electron mass e_m allowing us to write the following according to equation 5.20:

$$\begin{bmatrix} B^{\alpha} \times V^{\beta} \times L^{\gamma} \times e_m^{\zeta} \times e^{-} \end{bmatrix} = [1]$$
(5.26)

$$\left[\left(\frac{kg}{A\times s^2}\right)^{\alpha_1} \left(\frac{kg\times m^2}{A\times s^3}\right)^{\beta_1} (m)^{\gamma_1} (kg)^{\zeta_1} (A\times s)\right] = [1]$$
(5.27)

$$\Psi\left(\frac{B^2 \times L^2 \times e^-}{V \times e_m}\right) = R \tag{5.28}$$

$$\frac{B^2 \times L^2 \times e^-}{V \times e_m} = \Psi^{-1}(R)$$
(5.29)

$$B = \sqrt{V} \times \frac{\sqrt{e_m}}{L \times \sqrt{e^-}} \Psi^{-1}(R)$$
(5.30)

Considering that e_m and e^- are constant and we do not change the system length, the scaling law thus states that:

$$B \propto \sqrt{V}$$
 (5.31)

This scaling law coincides with the formula for the angle of rotation in eq. 5.13, obtained from the dioctron frequency ω_D . If we wish to determine the scaling for equal profiles, the angle of rotation of the different profiles must be equal. Keeping φ constant in eq. 5.13, we obtain the same scaling law of $B \propto \sqrt{V}$.

Having discussed a possible scaling law, fig. 5.15 sorts measured profiles by their magnetic field B in the main solenoid and the scaling factor. The scaling can not be replicated perfectly, but is observable to a given degree. Especially at the lowest scaling factor of $4.47 \times 10^{-3} \text{ TV}^{\frac{1}{2}}$. Here the only profile diverging from the scaling is that at B = 0.4 - 1.6 - 0.4 kG and V=500 V. Qualitatively the scaling laws holds, but profiles break down at high cathode potentials, making it difficult to confirm the scaling law.

Quantity	Symbol	Units	SI Units
Electron charge	e^-	[C]	$[A \times s]$
Electron mass	e_m	[amu]	[kg]
Magnetic field in solenoids	В	[T]	$\left[\frac{kg}{A \times s^2}\right]$
Cathode potential	V	[V]	$\left[\frac{kg \times m^2}{A \times s^3}\right]$
System length	L	[m]	[m]

Table 5.7: Physical quantities used for the dimensional analysis of the constant

beam profile. Neutralization, which would be in the form of a cross-section, and



Chapter 6

Transverse Fields of Measured Profiles

This chapter discusses the transverse electric fields of the electron beam and an upper limit for the emittance growth due to the electric field of the *HEB*.

Using WARP¹ the charge density, electric potential and electric field of the measured profiles are calculated. It is important to note that WARP is only being used as a tool for the calculations. It is used in a 2D setup and no 3D simulations are done. Therefore the detailed discussion of WARP is left to chap. 7, where full 3D simulations are conducted. The transverse fields will be used in tracking studies to examine the effect on the proton beam core and halo. As a preliminary check, the upper limits of the emittance growth due to the transverse electric fields is calculated by assuming the E-field fluctuations in the hole act like random kicks.

WARP was developed in the pursuit of heavy-ion driven inertial confinement fusion. It has been greatly enhanced since and is suited for the simulation of a *HEB*. It contains three coordinate systems: a 3D model (WARP3d) we use in chap. 7; an axis-symmetrical model (WARPrz) and a transverse slice model (WARPxy), which is used here. For further information refer to [35, p. 9]

6.1 Transverse Field of Hollow Electron Beams

The field of a *HEB*, can be described using the first Maxwell equation, Gauss' Law in integral and microscopic form.

$$\oint_{\partial\Omega} E \cdot \mathrm{d}S = \frac{Q_{\mathrm{encl},\Omega}}{\varepsilon_0} \tag{6.1}$$

The left hand side is an integral over the electric field E of the surface of a cylinder $\partial\Omega$. The electric field is assumed to be constant everywhere on the surface $\partial\Omega$. $Q_{encl,\Omega}$ is the charge enclosed by $\partial\Omega$ and ε_0 is the vacuum permittivity. In a 2D approximation, where the cylinder is assumed to be infinitely long, the front and back parts of the cylinder can be neglected. The right hand side, assuming a uniform line charge distribution, is written as an integral over the charge density ρ .

$$E \cdot 2\pi rL = \frac{\int_0^R \rho(r') 2\pi r' L \mathrm{d}r'}{\varepsilon_0}$$
(6.2)

 $r = r(\partial \Omega)$ is defined as the radius at the transverse surface of the cylinder and L as the length of the beam. The electric field E is singled out and some constants are eliminated.

$$E = \frac{\int_0^R \rho(r') r' \mathrm{d}r'}{\varepsilon_0 r} \tag{6.3}$$

 $^{^1\}mathrm{A}$ particle in cell code developed at Lawrence Berkley National Laboratory.

The inner beam radius is R_i and the outer R_o . Assuming a uniform charge distribution $\rho(r') = \rho$ in the hollow beam, 3 cases are considered.

$$r \le R_o \qquad \qquad E = 0 \tag{6.4}$$

$$R_{i} \ge r \le R_{o} \qquad E(r) = \frac{\rho(r^{2} - R_{i}^{2})}{2\varepsilon_{0}r} = \frac{IL(r^{2} - R_{i}^{2})}{2\pi v_{z}\varepsilon_{0}r(R_{o}^{2} - R_{i}^{2})}$$
(6.5)

$$R_o \ge r \qquad \qquad E(r) = \frac{\rho(R_o^2 - R_i^2)}{2\pi\varepsilon_0 r} = \frac{IL}{2\pi v_z \varepsilon_0 r} \tag{6.6}$$

These 3 equations fully qualify the electric field of a hollow uniform cylindrical charge distribution. Fig. 2.5 shows the transverse electric field distributions for the profiles corresponding to V=500 V and V=8 kV at B=1-4-1 kG ($B_{gun} - B_{main} - B_{coll}$).

6.2 Emittance Growth

From the radial electric field of the beam, the energy of the protons and the length during which the protons are exposed to the electric field, an upper boundary for the additional emittance growth of the proton beam can be estimated. The purpose hereof is to obtain a rough estimate of the emittance growth of the proton beam in the Large Hadron Collider (LHC) due to the HEBL.

Assuming the transverse electric field to be the only source of a transverse force on the proton particles, the following equation holds:

$$\frac{\mathrm{d}}{\mathrm{d}t}p_r = F_y = q\vec{E}_r \tag{6.7}$$

Here the subscript r indicates the radial direction in the transverse plane, p is the momentum q the elementary charge and \vec{E} the radial electric field. Assuming that the particles are on a perfect design orbit $(p_r = 0)$, and that the electric field is constant in time, the above equation can be integrated.

$$p_r = q\vec{E}_r t \tag{6.8}$$

We define θ to be the inscribed angle between the total momentum and the axial momentum along z. Expressing p_r in terms of the total momentum and the momentum along z, $p_r = p \sin(\theta) = p_z \tan(\theta)$ gives:

$$p_z \tan(\theta) = q\vec{E}_r \tag{6.9}$$

$$\rightarrow \tan(\theta) = \frac{q\vec{E}_r t}{p_z} \tag{6.10}$$

$$\rightarrow \tan(\theta) = \frac{q E_r L}{p_z v_z} \tag{6.11}$$

$$\rightarrow \tan(\theta) = \frac{\gamma E_0 q \vec{E}_r L}{p_z^2} \tag{6.12}$$

$$\rightarrow \tan(\theta) = \frac{E_{tot}q\vec{E_r}L}{(E_{tot}^2 - E_0^2)} \tag{6.13}$$

Given that the kick in the transverse direction is x' = dx/dz, it follows that $x' = \tan \theta \stackrel{!}{=} \vartheta$. Here γ is the relativistic factor, E_0 is the rest energy of the proton, v_{0z} is the velocity along z, q is the elementary charge, L is the length of the *Hollow Electron Gun (HEG)* and E_{tot} is the total energy of the proton beam. The transverse kick on a particle due to a transverse electric field is thus:

$$\vartheta = \frac{E_{tot}qE_rL}{E_{tot}^2 - E_0^2} \tag{6.14}$$

A normal transfer equation for a particle in an accelerator is given by [36]:

$$\begin{pmatrix} x_{n+1} \\ x'_{n+1} \end{pmatrix} = M \begin{pmatrix} x_n \\ x'_n \end{pmatrix}$$
(6.15)

where M is the transfer matrix. If the particles are not disturbed by the electron lens, the Courant Snyder invariant (see: [36, p. 105])

$$\varepsilon = \gamma x_{n+1}^2 + 2\alpha x_{n+1} x_{n+1}' + \beta x_{n+1}'^2 = \gamma x_n^2 + 2\alpha x_n x_n' + \beta x_n'^2$$
(6.16)

is conserved. Considering the kick that is introduced by the electron lens, eq. 6.15 changes to:

$$\begin{pmatrix} \tilde{x}_{n+1} \\ \tilde{x}'_{n+1} \end{pmatrix} = M \begin{pmatrix} x_n \\ x'_n \end{pmatrix} + \begin{pmatrix} 0 \\ \vartheta \end{pmatrix}$$
(6.17)

Assuming that the electron lens is turned on after n turns of the beam, the change of emittance per turn can be found. The emittance after turn n + 1 is:

$$\varepsilon = \gamma \tilde{x}_{n+1}^2 + 2\alpha \tilde{x}_{n+1} \tilde{x}_{n+1}' + \beta \tilde{x}_{n+1}'^2 = \gamma x_{n+1}^2 + 2\alpha x_{n+1} (x_{n+1}' + \vartheta) + \beta (x_{n+1}' + \vartheta)^2$$
(6.18)

The difference is given by

$$\Delta \varepsilon = 2\vartheta (\alpha x_{n+1} + \beta x'_{n+1}) + \beta \vartheta^2 \tag{6.19}$$

Given Floquet Transformations (see [37, p.37]), the emittance is retained, but projected onto a circle in a transformed phase space with the axis x and $\alpha x + \beta x'$. In this system, the contribution made to the emittance growth by the first term in eq. 6.19 is always orthogonal to the growth and therefore the first term can be neglected. The emittance growth can thus be given as:

$$\Delta \varepsilon = \beta \vartheta^2 = \beta \left(\frac{E_{tot} q E_r L}{(E_{tot}^2 - E_0^2)} \right)^2 \tag{6.20}$$

6.3 Method and Script

WARP is a Python package that includes Fortran² in order to facilitate the simulation of particle beams ³. Scripts are written in Python, using WARP specific syntax. The script used to calculate the transverse fields and the emittance growths (Fieldsolver_single.py) is included in appendix C.2.

After loading the appropriate Python packages and setting options, the code defines a list of all the text files containing the particle positions, produced by **profile_analysis**. R (see sec. 5.4.1) as well as lists containing the parameters of those profiles, such as the magnetic fields, the cathode potentials and the beam currents. The script has to be executed individually for each profile, setting the profile to be executed through the variable n (Code Line: 104). It is not recommended to execute several simulations in one run. Upon starting a new run, the charge distribution is not correctly deleted by WARP. One should thus quit Python and restart it between simulations or run Python in a non-interactive mode.

Having defined author information and titles for WARP, it defines the species that is to be loaded, electrons (Code Line: 123). The code then defines the lattice which constitutes of a drift pipe with a transverse number of mesh cells defined by nx and ny and a longitudinal mesh size of 1 cell (Code Line: 138-146). The drift pipe is installed as a conductor to enforce the boundary conditions of $\vec{E_r} = 0$. Consecutively, beam parameters such as the current and the beam energy are set. With all preliminary settings completed, the script now calls the 2D WARP package and generates. The conductor is installed, the particles are loaded and the charge is deposited onto them by loadrho() (Code Line: 202). Particle coordinates are loaded into transverse coordinate vectors xinit and yinit. The particle positions are

²A programming language specifically suited for numerical computing.

³WARP is available at warp.lbl.gov. For a Willson-Cluster specific WARP, I have compiled a zip-file with installation instructions that can be found here .*TELTS* wiki page

scaled to the correct positions, in order to accommodate magnetic (de)compression. This scaling is done using $\sqrt{B_{coll}/B_{main}}$.

At this point seven types of plots are made: particle plot, charge density plot, electric potential plot, 3 transverse electric field plots and an electric field line plot along the x and y-axis. Furthermore the script writes the electric fields to a file which has the name _fields.txt appended (Code Line: 262). The transverse fields have thus been completely characterized and the script starts with the determination of the upper limit for the emittance growth.

In order to discuss the emittance growth under expected operating conditions of the *HEBL* at the LHC, the emittance growth has to be calculated with the inner radius of the electron beam set at 4 σ and 6 σ of the LHC beam. σ is the standard deviation of the Gaussian distribution of the proton beam in the LHC.

To calculate the σ of the LHC, we use the two relations (see [10, p. 66])

$$\varepsilon_{RMS} = \frac{\sigma_x^2}{\beta} \qquad \varepsilon_N = \varepsilon_{RMS} \times \gamma \beta_c \tag{6.21}$$

where ε_N is the normalized emittance, ε_{RMS} is the *Root Mean Square (RMS)* emittance as defined by [38], γ and β_c are the relativistic parameters, σ_x is the standard deviation from the beam in the x direction and β is the beta-function of the beam at the point of insertion of the HEBL. We set β to 250 m [39]. The normalized emittance is 3.75 μm as specified in [7, p. 243]. σ is thus given by

$$\sigma_x = \sqrt{\varepsilon_{RMS}\beta} = \sqrt{\frac{\varepsilon_N\beta}{\gamma\beta_c}} = 4.69 \times 10^{-4} \ m \tag{6.22}$$

at $\beta = 250$ m, $\varepsilon_N = 3.75 \ \mu m$ and $\gamma = 4264.39$.

Given σ and the cathode radius, the script applies a Gaussian weighting with a standard deviation of σ to the measured electric field. This factors in the probability of a particle from the LHC proton beam interacting with a given point of the electric field produced by the *HEBL*. Considering that the transverse profile has been compressed by the magnetic field, $\sqrt{B_{gun}/B_{main}}$, the script now selects all points in the 80%ile of the cathode radius. It then takes the *RMS* value of all those points and uses them as the transverse field in eq. 6.20. This can be done, because we assume that the electric field in the center acts like random kicks on the core protons and thus only the *RMS* value of the electric field is significant.

At this stage the transverse field is given at the magnetically compressed cathode radius, $6.75 \times 10^{-3} \sqrt{B_{coll}/B_{main}}$ and stored in the variable Er_center (Code Line 326). The scaling for the radius goes as R⁻¹ as shown in equation 6.3. The emittance is thus scaled further down to 4σ and 6σ and stored in the variables Er_center_4 and Er_center_6 (Code Line: 327-328).

Using eq. 6.20, we can thus calculate the emittance growth per turn. In order to obtain the emittance growth in units of s⁻¹, we divide $\Delta \varepsilon$ by the normalized emittance and the revolution time. This quantity is denoted *Emittance Growth Rate (EGR)* (Code Line: 337. The transverse electric field, emittance growth per turn and *EGR* are then saved in a table for the cathode radius, and a scaled radius of 4 σ and 6 σ . It is important to note that this estimate of the emittance growth is only a rough upper limit, because in reality multipole components will be present in the electric field that do not cause an emittance growth.

6.4 Results

Here the general shape of the electric field and the upper limit for the emittance growth are discussed.

6.4.1 Field Measurements

Fig. 6.1 shows the plots obtained from WARP for the transverse electric fields of the profile shown in fig. 5.12 at V=500 V and B=1-4-1 kG. Since the charge density in fig. 6.1a is symmetric with a slight

Emittance [µm]							EGR	[1/s]	Diffusion [μ m/s]		
Plane	ϵ_{ws}	$\delta \epsilon_{ws}$	ϵ_{sl}	$\delta \epsilon_{\scriptscriptstyle WS}$	ϵ	$\delta \epsilon$	γ	$\delta\gamma$	D'	$\delta D'$	
B1 H	1.92	1.3×10^{-2}	2.15	$9.6 imes 10^{-3}$	2.03	0.07	5.23×10^{-5}	$1.50 imes 10^{-5}$	2.49×10^{-8}	7.4×10^{-9}	
B1 V	1.36	$2.0 imes 10^{-3}$	1.29	$4.7 imes 10^{-3}$	1.33	0.02	$1.70 imes 10^{-5}$	$1.00 imes 10^{-5}$	$5.27 imes 10^{-9}$	3.2×10^{-9}	
B2 H	1.69	$6.6 imes 10^{-3}$	2.43	1.3×10^{-2}	2.06	0.21	5.37×10^{-5}	$1.90 imes 10^{-5}$	$2.60 imes 10^{-8}$	$9.6 imes 10^{-9}$	
B2 V	1.73	$4.6 imes 10^{-2}$	1.98	$1.2 imes 10^{-2}$	1.85	0.09	$1.10 imes 10^{-5}$	$1.70 imes 10^{-5}$	$4.79 imes 10^{-9}$	$7.2 imes 10^{-9}$	

TABLE VI. Initial emittances (wire scan, synch-light, average), emittance growth rates (EGR) γ , and diffusion slopes D'.

Table 6.1: Current EGR at the LHC. Source: [39, p. 8]

$EGR [\times]$	$10^{-5}\frac{1}{s}$]				Cathod	e Voltag	ge [kV]			
Radius [mm]	B_{max} [kG]	500	1000	2000	3000	4000	5000	6000	7000	8000
2 205	4	0.028	0.076	0.344	30.2 (3125 V)	26.6	25.6	24.5 (6250 V)	-	691
	3.2	0.008	0.029	9.05	15.1	22.9	276	617	978	1050
5.525	2.4	0.003	1.42	5.85	70.3	192	318	-	-	-
	2.4	0.414	1.40	44.7	-	-	-	-	-	-
	4	0.039	0.109	0.495	43.4 (3125 V)	38.3	36.8	35.2 (6250 V)	-	994
6 7	3.2	0.011	0.041	13.0	21.7	33.0	397	887	1410	1510
00	2.4	0.005	2.04	8.41	101	276	-	-	-	-
	2.4	0.595	2.02	64.4	-	-	-	-	-	-
4 σ	4	0.089	0.246	1.11	97.8 (3125 V)	86.2	82.8	79.3 (6250 V)	-	2240
	3.2	0.025	0.093	29.3	48.7	74.3	893	2000	3170	3390
	2.4	0.010	4.59	18.9	227	621	-	-	-	-
	2.4	1.34	4.54	144	-	-	-	-	-	-

Table 6.2: The EGR due to the transverse fields caused by the HEB at the given inner beam radius, magnetic field strength and cathode potential. Green cells give acceptable EGR.

elongation in the y-direction, the resulting electric potential and electric fields are symmetric around the z-axis. The electric field in the center is $\approx 1 \text{ kV m}^{-1}$, as shown in fig. 6.1b. Comparing this plot to fig. 2.5a, shows that the measured and theoretical electric fields coincide very well. This furthermore indicates that our assumptions in the theoretical deduction in section 6.1 are correct and thus the charge density is near uniform.

In order to contrast these results at a low cathode potential with higher cathode potentials, the plots for V=8 kV and B=1-4-1 kG are shown in fig. 6.2. The charge distribution in fig. 6.2a is clearly asymmetric. The charge is concentrated in the lower right corner, causing the electric field in the center to be of the order of $\approx 300 \text{ kV m}^{-1}$. A small area of significantly lower electric field, $\leq 100 \text{ kV m}^{-1}$, is apparent in the lower right corner in fig. 6.2f. One would thus expect the proton beam to be much more perturbed at a cathode potential of -8 kV instead of -500 V.

6.4.2 Upper Limit for Emittance growth

As a reference for the EGR values published on 19 February 2013 in Physical Review Special Topics Accelerators and Beams are used [39, p. 8]. They are presented in fig. 6.1. One should note that the nominal emittance of the LHC of 3.75 μ m has not yet been reached.

Table 6.2 presents the EGR in units of 10^{-5} s⁻¹. The quantity of 1×10^{-5} s⁻¹ was chosen as an upper limit below which the EGR is acceptable, on the basis that EGR values in table 6.1 are just above that threshold.

When one compares table 6.2 with fig. 5.10, we can see that the EGR tends to rise above 1×10^{-5} s⁻¹ only when the charge distribution deviates strongly from an azimuthally uniform distribution. Values then tend to rise significantly.

Given that the scaling law $B \propto \sqrt{V}$ in section 5.4.3 holds to a certain degree, we are interested in profiles with low cathode potential and high magnetic field. Because the *EGR* values are only upper boundaries, due to the presence of multipole components, no error calculation was done.







(c)Electric potential Electric field strength in x-y plane





(f)Electric field in center





(b)Electric field Line Plot Electrostatic potential in x-y plane



(d)Electric potential in center Electric field strength in x-y plane





(a)Charge density Electrostatic potential in x-y plane



(c)Electric potential Electric field strength in x-y plane







Electrostatic potential in x-y plane



(d)Electric field Line Plot Electric field strength in x-y plane



(e)Electric field in y direction

(f) Transverse electric field



Chapter 7

Numerical Simulations of Electron Beam Dynamics

Having discussed all experimental measurements, the first attempt of 3D numerical simulations of the Hollow Electron Beam Lens (HEBL) through a Particle-In-Cell (PIC) code called WARP¹ will be discussed. This serves as a basis for future 3D beam simulations in order to implement a non-linear map for core and halo tracking studies in SixTrack or Lifetrack. Additionally such simulations are promising for calculating the 3D magnetic and electric fields for the electron lens and for examining the effect of the bends in Tevatron Electron Lens 2 (TEL2) on the beam core. During this thesis, WARP was installed on the Accelerator Simulations Cluster (TEV)² and the first full simulations were completed.

WARP was developed at Lawrence Berkeley National Laboratory in California, USA, by Dave Grote³, Jean-Luc Vay⁴ and Alex Friedman⁵. WARP is a *PIC* code that was designed in order to simulate high current particle beams. It runs in the highly interactive Python interpreter, using precompiled packages. The interpreter allows one to control the simulations interactively.

7.1 Introduction to WARP

WARP was installed on the TEV at Fermi National Accelerator Lab (Fermilab). Version-compatibility issues with Python installed on the cluster prevented the installation of WARP. Therefore Python was installed in the home directory with the necessary packages. This new installation was used to compile WARP. Installation instructions and files can be found on the Electron Lens Wiki⁶. Future users of WARP on the TEV may choose to link to my local installation of WARP instead of reinstalling. For this the **\$PATHs** must be adapted as explained in the instructions on the Wiki.

During this thesis, WARP codes were developed on local machines using a low number of iterations. Normal laptops were mostly sufficient for this. Full simulations were run on the TEV at Fermilab. The TEV cluster has 32 AMD nodes and 12 Intel nodes. Each AMD node has 32 processors and 64 GB of RAM. The WARP simulations are highly memory intensive. In order to run a full simulation, 12 nodes were used with 2 processes per node. This allowed us to use 24 parallel processors and 768 GB of memory. At lower amounts of memory, the code crashed due to heavy disk swapping. This is an issue that will have to be investigated but could not be done in the scope of this thesis.

As mentioned in chapter 6, WARP contains several basic models, including a transverse slice model (WARPxy) and an axis symmetric model (WARPrz). Models exist in 2D and 3D. Additionally, three coordinate frames are available. The lab frame in which all transverse positions are given. In this frame,

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⁶https://cdcvs.fnal.gov/redmine/documents/619

the user sets a 0-point and the axial-coordinate increases continually along the axial position. The lab frame follows bends when they are used. In closed accelerators, the lab frame axial coordinate will increase lap by lap. The beam frame and grid frame move along with the beam. The location of the beam frame is given by zbeam.

PIC codes are well known from plasma physics and are used to simulate neutral and non-neutral plasmas. In a *PIC* code, macromolecules are assigned to cells on a discrete mesh and advanced using the Lorentz equation of motion. Often such codes are also called "Nearest-Grid-Point Particle Mesh models". From the distribution of the macromolecules on the mesh, the charge density is calculated, using a linear interpolation of the particles on the mesh. By solving Poisson's equation, the electrostatic potential and ultimately the electric field are calculated. These calculations are done for each time step, after which the velocities and positions are updated. As explained in [40, p. 28], while positions are usually updated during each full time step, the velocities of the particles are updated every half time step. This can cause errors in diagnostics. Whenever a diagnostic is made in WARP, the isochronous leap frog system is used, for which velocities are only updated at half time steps. For a detailed explanation of the time-step layout, refer to [35, Chap. 3].

WARP combines the power of *PIC* codes with the easy implementation of accelerator lattices. An extensive set of lattice elements are precompiled and included in WARP. These include dipole, quadrupole and sextupole magnets, accelerator cavities and solenoids. Additional elements can easily be implemented through surfaces of revolution. Furthermore WARP includes several diagnostics packages and allows one to either use a set of precompiled plotting methods or develop individual plotting methods. These can then be executed at predefined time points during the simulation and interactively after the simulation.

7.2 General Structure of the WARP Script

The code used for the WARP simulations is listed in appendix C.3. This section discusses the general layout of a WARP code. We have used two injection methods throughout the simulations. First a direct injection from a replica of the 1 inch Hollow Electron Gun (HG1b) and second an injection of the measured profiles from the test bench at the z-coordinate of the cathode. The injection mechanisms will be discussed more closely in sec. 7.3 and 7.4.

The WARP code initially calls the WARP package. The code is separated into a few sections: file loading, options, headers, variables and the actual scripting, which is divided into further sections. The file loading sets the profile that is to be injected or simulated. In options (Code Line: 115) the injection type is defined in the variable machine_injtype and the cathode emittance method is set in machine_injtype. For gun injection, the cathode emits according to the Child-Langmuir law. For profile injection, it emits at a constant current. WARP requires the user to define a variable called runid, which sets the name of the current run. Additionally one can set some variables that are used as descriptors in the plots and the variable runmaker which defines the author of the run. In the section variables (Code Line: 144), the parameters of all lattice elements and some beam parameters are defined, in order to prevent the user from having to deal with the more complicated WARP codes in the section Script (Code Line: 267).

Having defined the **runid** and the **runmaker**, the **setup** (Code Line: 271) command must be called in order to initialize the graphic output file. The output is saved as *Computer Graphics Metafile* (*CGM*). A software called gist⁷ is used to view and manipulate the plots from the *CGM* files. If profile injection is chosen, the selected measured profile from the test bench⁸ is loaded and saved in an array with the first column being the x-coordinates and the second column the y-coordinates. The number of particles that were loaded is printed to the standard output.

⁷Available at warp.lbl.gov

⁸Profile measurements in the test bench are explained in section 5.4

At this stage the code defines the particle species, in our case electrons are designated elec (Code Line: 303). It also defines the number of particles per macro-particles as sw. It then defines beam parameters such as the beam current, the kinetic energy and the time step used in the simulation. The time step is significant, because it needs to be small enough in order to consider all physical effects, such as the plasma frequency ω_p^{9} and the cyclotron frequency ω_c^{10} . For the *HEBL* simulations, the cyclotron frequency exceeds the plasma frequency by a few orders of magnitude. In the case of a 500 V beam with a magnetic configuration of 1-4-1 kG, $\omega_c=7.0 \times 10^{11} \text{ s}^{-1}$ and $\omega_p=4.6 \times 10^8 \text{ s}^{-1}$. The time step is thus determined by the cyclotron frequency. In order to be certain that we use a small enough time step, it was chosen that

$$\frac{\omega_c \Delta t}{2\pi} = 0.25 \quad \Rightarrow \quad \Delta t = \frac{\pi}{2\omega_c} \approx 1 \times 10^{-12} \tag{7.1}$$

The electron velocity is set to 0 by the code and deduced from the electron energy through the command derivqty() (Code Line: 315). Next the code sets whether calculations should be done in a relativistic manner. The thermal velocity of the particles is set to 0. This should be changed in future versions of this code. Having set all the beam parameters, the number of steps necessary to complete the simulation are calculated via:

$$n_{steps} = 1.1 \times \frac{L}{v_e \Delta t} \tag{7.2}$$

Here L is the distance the beam travels from gun to collector, v_e is the electron velocity and Δt is the time step in eq. 7.1. The factor of 1.1 is used in order to increase the number of steps by 10% to accommodate for the fact that particles move slower than v_e while they are in the gun. Next the code defines the size of the electron beam and where it is to be injected (323-357). Amongst others it defines what type of injection is to be used.

Having set up the particle injection, the code now defines the solenoids. It is important to note that these solenoids do not manifest themselves as conductors in the code, but only provide the necessary magnetic fields. It is not necessary to include the solenoids as conductors, since they are positioned outside the drift tube and thus do not affect the electric field inside of the beam tube and are not involved in scraping. Next the script defines the lab frame mesh. When running the code in parallel mode, it is important that the number of grid cells in each direction exceeds the number of processors used to run the script.

In WARP it is possible to set up diagnostic windows, limiting the number of particles that will be involved in the diagnostics. For example is it possible to set up a window so that only the particles in the first 2 time steps after the cathode are plotted. For further information on diagnostic windows please refer to [35, Chap. 7]. Once diagnostic windows are set, it is important to define how often simulation data should be saved (nhist) and plots are made (itplps, itplfreq and itplalways) (Code Lines: 417-433). One can then choose a couple of precompiled plots. These plots are generally less useful and it is recommended to create ones own plots. For more information on plotting refer to [35, Chap. 7].

Having set up the diagnostics and the mesh, the field solver needs to be chosen. First of all the conductor boundary walls are set to absorb particles. Several field solvers are available with WARP. For more information on the individual field solvers, refer to [35, p. 61]. The field solver used is set by the variable top.fstype (Code Line: 442) which is set to 7, 3D multi-grid solver. The multi-grid solver is an extension of the Successive Over-Relaxation Field Solver (SOR). It uses the finite difference form of Poisson's equation in combination with a relaxation parameter that is given by:

$$\varphi_{ijk} = \omega \Delta^2 \left(\frac{\rho_{ijk}}{\varepsilon_0} + \frac{\varphi_{i+1jk} + \varphi_{i-1jk}}{\Delta x^2} + \frac{\varphi_{ij+1k} + \varphi_{ij-1k}}{\Delta y^2} + \frac{\varphi_{ijk+1} + \varphi_{ijk-1}}{\Delta z^2} \right) + (1 - \omega)\varphi_{ijk}$$
(7.3)

 ${}^{9}_{10}\omega_p = \frac{nq^2}{\varepsilon_0 m}$

$$\omega_c = \frac{q_B}{\gamma m}$$



Figure 7.1: Technical drawing and plot of the gun that was implemented into WARP.

here φ is the potential of the respective cell designated by ijk, ω is the relaxation parameter, ρ is the charge density and $\frac{1}{\Delta^2} = \frac{1}{\Delta x^2} + \frac{1}{\Delta y^2} \frac{1}{\Delta z^2}$ where $\Delta(x, y, zy)$ are the respective grid cell sizes [35, p. 69]. At this point the code is ready to initialize. A package is initialized using the command generate and a time step is made using the command step. The 3D model is initialized (allocating storage, load particles, etc...) using package("w3d");generate() (Code Line: 464)). It is time to implement all the conductors. Both injection methods require the implementation of the drift tube which has a radius of 3 cm. On top of that, the gun injection requires the implementation of all necessary conductors in the gun. The drift tube can be implemented using a precompiled version of a solenoid ZCylinderOut, which defines the outside of a cylinder as a conductor. All conductors that are defined must be installed using installconductors. Lastly self made plots are defined for the diagnostics. Plots that are to be repeated at the time steps given by itplalways and itplseldom must be defined as a function and installel using installplalways and installplseldom (Code Line: 567).

Having completed all the preparatory work, the simulation can now be executed. With the command step one can define how many time steps are to be made. By writing steps(nsteps) (Code Line: 570) the full simulation is done. In order to test new settings, it might be useful to just run a few iteration steps. Lastly a dump of the simulation can be made. This has not yet been implemented, since when running the code on 32 processors in parallel mode, 32 dump files are created, which is counterproductive.

7.3 Gun Injection

Gun injection refers to the injection of particles directly of the cathode of the gun. It is thus necessary to implement the gun in the code according to the specifications given in the technical drawings of the gun, fig. 3.7. Fig. 7.1a shows the parts that were implemented and fig. 7.1b shows a plot of the conductors in the WARP code. The most important dimensions of these conductors are listed in appendix D. All 4 conductors were installed using the function ZSrfrv which creates surfaces of revolution. The function takes 3 vectors containing the radial coordinates, the axial coordinates and the radii of curvature for the sections between the specified points. The surface spanned by the points in the vectors is then spun around the lab frame beam axis to produce the conductors in 3D space. The function immediately sets the voltage of the conductors. The conductors must be passed on as a list to the command installconductor in order to be implemented.

On top of installing the gun, the injection of the particles must be specified. This is done earlier in the code in the section labelled injection. First the cathode emission is set to Child-Langmuir law. This is set in the variable top.inject. Additionally the surface from which the particles are injected, their

>0 00

-0.02

0.00

0.05

(c)Electric field in z of gun.

7

0.10

0.15

>0 00

-0.02

0.00

0.05

(a)Electric potential of qun.

7

0.10

0.15



drawing that are implemented.

(b)Gun as it is implements in WARP.

(c)Gun as it is implements in WARP



0.05

(b)Electric field of gun.

Figure 7.3: Plot of the electrostatic potential and electric field of the gun implemented in WARP.

7

0.10

0.15

>0 00

-0.02

0.00

Figure 7.2: Technical drawing and plot of the gun that was implemented into WARP.

angles and their velocities are defined. Lastly the voltage of the injector source is set with vinject. An additional variable called elec.npinject defines how many macro particles are injected during each time step. In the case of Child-Langmuir law injection, this value may be set to an approximate value, since the law will change the number of particles injected as necessary.

Fig. 7.2 shows the settings used for a normal run on the TEV. The generalized perveance in fig. 7.2b serves as a reference to our measured gun perveance. In section 5.3.5 we measured a generalized perveance of $7.93(7) \times 10^{-3}$. The WARP simulation gives us a generalized perveance of 9.89×10^{-2} . Our measured generalized perveance is $\approx 80\%$ of our simulated generalized perveance.

The electrostatic potential and electric field of the gun that was implemented in WARP is shown in fig. 7.3. It is important to notice that one needs to increase w3d.nz to a very high number in order to present the field correctly (2048). It is thus recommended to reduce the system length machin_syslen and only simulate the gun for these plots. The plots in fig. 7.3 were made with w3d.nx=128, w3d.ny=128, w3d.nz=512.

Lastly fig. 7.4 shows the full trajectory of the beam. Three main features can be recognized. First, we see the clear magnetic compression of the beam as it enters the main solenoid that has a 4-fold magnetic field, meaning a 2-fold compression. Additionally we see that the beam is nicely hollow, portrayed through the two yellow bars. And lastly we see some suspected micro-bunching as the beam reaches the end of the beam tube.





Figure 7.4: Full electron beam as implemented in WARP due to gun injection. Three important features are visible: the magnetic compression of the beam before and after the main solenoid, the hollow center of the beam and micro bunching towards the end of the beam.

7.4 Profile Injection

Profile injection refers to the injection of particle distributions measured in the test stand. The assumption is that the measured profile at high magnetic field and low cathode potential has not yet undergone significant space charge evolution (see sec. 5.4.2) and is thus very similar to profiles that are emitted directly from the gun. The measured profile that is used for profile injection is that at a magnetic field of B=1-4-1 kG and a cathode potential of -500 V. This measured profile distribution is shown in fig. 7.5.

The injection of this profile is a user defined injection method in which a vector **fns** passes the path of the profile to the code. The code then imports the particle positions through the command **fromfile** into a vector **posi**. The x and y coordinates are read into one long vector in sequence, requiring one to split the single vector into an array with a column of x coordinates and a column of y coordinates. This is done using **reshape**. During profile injection, a test is conducted to verify that the right particles are injected, which is given through the variable **w3d.inj_js**, and the num-



Figure 7.5: Particle distribution of measured profile at $B=1-4-1 \ kG$ and $V=500 \ V$

ber of particles to be injected each time step is defined. Next the coordinates are passed on to w3d.xt, w3d.yt and w3d.zt which define the particle positions in the 3D lab frame. Additionally the initial



Figure 7.6: Full electron beam as implemented in WARP due to profile injection. The electron density is too high using this injection.

velocities are set to 0 for the transverse velocities and to the beam velocity for the axial velocity. To finalize injection, the source has to be installed using installuserparticlesinjection.

The injection method is not finalized at this stage, but initial runs have been conducted. The main issue with this injection type is that it injects too many particles. This is seen when comparing figure 7.6 with fig. 7.4, which has a ≈ 60 times lower particle count. The high particle count makes it nearly impossible to process these profiles. It is assumed that this is a consequence of the whole loaded profile being injected at each time step. Further development on this injection was not done, since gun injection was considered to be more important.

7.5 Directions for further studies on 3D WARP simulations

WARP has successfully been installed on TEV and the first few runs have been completed. Nonetheless, the simulations are far from complete and many things need to be improved. This section outlines some of the steps that need to be implemented in the current code:

- First of all the code needs to be optimized in order to use less computational resources.
- Additional diagnostics and plotting tools should be implemented. For example it would be interesting to implement a few diagnostic windows that allow us to compare the transverse profiles from the simulations with those from the measurements in *Tevatron Electron Lens Test Stand* (*TELTS*).
- As discussed previously, the profile injection needs to be optimized in order to reduce the number of particles that are injected. Additionally to test whether the perveance and the beam current resemble measured values needs to be performed.
- The dumping of the simulations needs to be implemented, so that runs, which sometimes take a whole day, do not need to be repeated.

• Extraction of simulation data is possible through the package ENV and should be implemented.

Furthermore, the *TEL2* lattice or an optimal electron lens lattice for the *Large Hadron Collider* (LHC) should be implemented. Using this deductions can be made on the effect of the electron lens bends on the *Hollow Electron Beam* (*HEB*). Furthermore it is possible to implement multiple species and sources in WARP, allowing the implementation of the accelerator proton beam and thus the direct simulation in WARP of the effect of the *HEB* on the proton beam in one single pass.

Chapter 8

Conclusion

The studies discussed in this thesis were completed in the context of supporting a conceptual design report on a Hollow Electron Beam Lens (HEBL) for proton beams at the Large Hadron Collider (LHC) at European Organization for Nuclear Research (CERN). Experiments conducted at the Tevatron over recent years have provided a solid basis for this beam scraping and halo control technique. Nonetheless, several further issues need to be addressed in the extrapolation of the HEBL from the Tevatron to the LHC. In this thesis, we have discussed measurements and considerations that pertain to the design and feasibility of Hollow Electron Gun (HEG) hardware parameters, and to the effects of the Hollow Electron Beam (HEB) on the dynamics of the circulating proton beam.

For halo scraping, the inner radius of the *HEB* is envisioned to be placed between $\approx 3 \sigma$ and 6σ of the *LHC* proton beam size. For the purpose of this thesis, σ was set to 4.7×10^{-4} m. For stability reasons, the field in the guiding solenoids should be as large as possible. Considering previous experience and technical feasibilities, we considered a magnetic configuration of B=0.43-5-0.43 T ($B_{gun} - B_{main} - B_{coll}$). This implies a magnetic compression of a factor 3.4, which sets the cathode inner radius to 4σ and the outer radius to $\approx 7.5 \sigma$.

A 1 inch Hollow Electron Gun (HG1b) based on a tungsten dispenser cathode was developed and tested. The perveance of this gun was determined to be $5.3(1)x10^{-6}$ perv, which is within 15% of numerical SAM simulations. This perveance implies a peak yield of over 5 A at 10 kV. Similar beam currents as were produced by the 0.6 inch Hollow Electron Gun (HG06) of 1.2 A, were measured using the HG1b at cathode voltages of 3-4 kV. Based upon Tevatron experiments and numerical simulations for the LHC, this yield should be sufficient to smoothly scrape 7-TeV protons.

During *HEG* characterization studies at the *Tevatron Electron Lens Test Stand (TELTS)*, it was found that this larger *HG1b* required a compression in order to obtain a transmission of 100% in the beam tube. Previous cathodes did not require such compression. The best achievable configuration in the *TELTS* was found to be $B_{gun} = B_{coll} = 0.25B_{main}$. This configuration results in a 2-fold compression of the *HEB*. Because the actual *HEBL* will have a larger inner aperture, we do not expect transmission degradation to be an issue. By examining the dependence of the cathode yield on filament current, it was found that the optimal filament current is about 8.5 A.

The transverse *HEB* profiles were measured to estimate the symmetry of the resulting electric fields and possible adverse effects on the proton beam core. Extensive profile measurements as a function of beam current and solenoidal magnetic field were completed. These allowed us to study the evolution of the magnetically confined *HEB* due to space charge. The profiles showed a scaling proportional to the accelerating voltage V_a and inversely proportional to the square of the axial magnetic field *B*, which is in qualitative agreement with the theory of the diocotron instability. Consequently, profiles measured at very low cathode voltages and at the highest possible magnetic fields in the conventional test stand solenoids will be similar to the profiles in the actual *HEBL* for the *LHC* using the parameters suggested above.

Effects on the beam dynamics were studied by calculating the transverse electric fields generated by the measured profiles, and by estimating the resulting emittance growth. For typical operational parameters, the effect was much smaller that the natural emittance growth in the LHC. At higher HEB currents or very low magnetic fields, the HEB loses its azimuthal symmetry and electric fields in the center can no longer be considered negligible. This, together with the Tevatron experiments, provides a strong indication that it is possible to operate the HEBL without negatively affecting the LHC proton core. Nonetheless, this aspect will have to be further investigated by implementing the fields generated by the HEB in a numerical tracking code.

In view of a realistic representation of the HEBL in tracking simulations, the first fully 3D calculation of the magnetized electron beam dynamics was completed using the WARP *Particle-In-Cell (PIC)* code. The HEG was implemented in WARP together with the geometry of the *TELTS*. Three main effects could be verified: the magnetic compression of the beam, the evolution of the hollow profile, and some longitudinal micro-bunching towards the end of the drift tube.

This thesis work advanced the understanding of the physics and technology of *Hollow Electron Beam* Collimation (HEBC) and provided the basis for further studies on the implementation of this technique at CERN.

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Appendices

Appendix A

LHC and Tevatron Parameters

Quantity	Symbol	Unit	Tevatron [16, 9]	LHC $[7]$
Collider type			$p-\overline{p}$	p-p
Circumference	C	[m]	6.28 km	$26.66 \mathrm{~km}$
# Beam Pipes			1	2
# of IRs	N_{IR}		2	8
Injection Energy	E_{inj}	[GeV]	120/150	450
Nominal Energy	E_{nom}	[TeV]	0.98	7
Achieved Energy	E	[TeV]	0. 980	4.5
# of bunches	N_b		36 (3 bunch trains)	2808
Protons per bunch	N_p		$2.7 \times 10^{11} / 1.35 \times 10^{11}$	$1.15 imes 10^{11}$
Bunch spacing	T_b	[ns]	396	25
Nominal Beam Intensity	$N_p, N_{\overline{p}}$		$9.72 \times 10^{12}, 4.86 \times 10^{12}$	3.23×10^{14}
Peak Luminosity	L_0	$[\rm cm^{-2} s^{-1}]$	3.15×10^{32}	$2.3 imes 10^{34}$
Nominal Emittance (RMS)	ε	$[\mu m]$	2.8/1.4	3.75
Nominal Energy Density	U	$\left[\mathrm{MJ/mm^{2}}\right]$	2	360
Cryogenic Temperature	T	[K]	4	1.9

Table A.1: Important parameters of Tevatron and LHC

It is important to notice that the proton and anti-proton beams in the Tevatron share the same beam pipe, in which they are transversely displaced. Beam-Beam effects are thus significantly more important at the Tevatron. In the LHC, both beams have their own channels, which intersect at 4 Interaction Regions (IRs), where the experiments ATLAS, CMS, ALICE, LHCb and others can be found.

Appendix B

Measured Profiles

B.1 List of measured profiles

#	Date	Magnetic Field	Cathode	Cathode	Avg.	Collector	Rep.	Vacuum	Pulse
		[T]	Poten-	Current	Current	Current	Rate	[V]	Length
			tial	[mA]	[mA]	[mA]	[Hz]		[mus]
			[kV]						
1	20.02.13 11.35	0.1-0.1-0.1	0.9	N/A	N/A	118	500	N/A	8
2	21.02.13 11.00	0.2-0.2-0.2	3.6	N/A	N/A	928	50	N/A	8
3	21.02.13 11.40	0.1-0.1-0.1	2.0	N/A	N/A	398	200	N/A	8
4	21.02.13 13.05	0.1-0.1-0.1	3.0	N/A	N/A	726	100	N/A	8
5	04.04.13 09.00	0.25 - 0.25 - 0.25	5.6	2322	1.21	1770	20	6.48	8
6	04.04.13	0.25 - 0.25 - 0.25	5.0	1980	1.25	1484	30	6.41	8
7	$04.04.13\ 12.25$	0.25 - 0.25 - 0.25	2.1	548	1.33	412	200	6.47	8
8	$04.04.13\ 13.45$	0.25 - 0.25 - 0.25	3.5	1156	1.35	865	80	6.41	8
9	04.04.13 15.08	0.3-0.3-0.3	5.0	1964	1.24	1460	30	6.41	8
10	$05.04.13\ 08.30$	0.2-0.2-0.2	1.3	282	1.33	212	400	6.57	8
11	05.04.13 10.00	0.2-0.2-0.2	2.2	592	1.335	448	180	6.57	8
12	05.04.13 11.20	0.2-0.2-0.2	3.1	972.6	1.35	744.4	100	6.52	8
13	05.04.13 12.45	0.15-0.15-0.15	0.75	120.8	1.26	90.4	800	6.52	8
14	05.04.13 14.15	0.15-0.15-0.15	2.0	506	1.25	390	200	6.51	8
15	05.04.13 15.47	0.15-0.15-0.15	1.3	254	1.21	192	400	6.52	8
16	02.05.13	0.25-0.25-0.25	0.50	69.6	0.96	51	1000	6.64	8
17	02.05.13	0.2-0.2-0.2	0.50	68.4	0.94	49	1000	6.60	8
18	02.05.13	0.3-0.3-0.3	0.50	68.4	0.94	56	1000	6.60	8
19	07.05.13	0.3-0.3-0.3	0.50	70.8	0.98	50.4	7000	6.58	8
20	13.05.13	0.3-0.3-0.3	0.50	74.4	1.05	50.2	1000	6.64	8
21	13.05.13	0.3-0.3-0.3	0.50	N/A	0.36	N/A	300	6.64	8
22	13.05.13	0.3-0.3-0.3	0.50	70.4	0.99	50.8	1000	6.60	8
23	13.05.13	0.3-0.3-0.3	0.50	N/A	0.34	N/A	300	6.60	8
24	13.05.13	0.1-0.3-0.1	0.50	72.4	1.02	69.8	1000	N/A	8
25	14.05.13 20.21	0.08-0.32-0.08	0.50	74.4	1.02	72.8	1000	N/A	8
26	14.05.13 20.22	0.08-0.32-0.08	1.0	198	0.98	195	400	N/A	8
27	14.05.13 20.24	0.08-0.32-0.08	2.0	528	1.07	526	150	N/A	8
28	14.05.13 20.31	0.08-0.32-0.08	4.0	1460	1.13	1450	40	N/A	8
29	15.05.13 11.49	0.08-0.32-0.08	8.0	3760	1.57	3700	10	N/A	8
30	15.05.13 12.31	0.08-0.32-0.08	7.0	3160	1.33	3100	10	N/A	8
31	15.05.13 13.22	0.08-0.32-0.08	0.25	27.0	1.09	25.8	2500	N/A	8
32	15.05.13	0.08-0.32-0.08	0.125	9.72	0.91	9.32	4000	N/A	8
33	15.05.13	0.08-0.32-0.08	3.0	956	1.32	944	100	N/A	8
34	15.05.13 15.39	0.08-0.32-0.08	5.0	7990	1.01	7990	15	N/A	8
35	15.05.13	0.08-0.32-0.08	6.0	2560	1.14	2530	10	N/A	8
36	15.05.13 17.53	0.1-0.4-0.1	1.0	197	0.97	196	400	N/A	8
37	15.05.13 18.51	0.1-0.4-0.1	3.125	1024	1.4	1022	100	N/A	8
38	15.05.13 19:20	0.1-0.4-0.1	6.250	2780	1.16	2754	10	N/A	8
39	16.05.13 10:45	0.06-0.24-0.06	1.0	798	0.99	796	400	N/A	8
40	16.05.13 11:02	0.04-0.16-0.04	1.0	797	0.99	795	400	N/A	8
41	16.05.13 11:46	0.04-0.16-0.04	0.50	74.4	1.02	72.4	7000	N/A	8
42	20.05.13 11:40	0.1-0.4-0.1	0.50	73.6	1.02	72.8	1000	N/A	8

 $Table \ B.1: \ Table \ showing \ all \ profile \ measurements \ made.$
43	20.05.13 11:42	0.1-0.4-0.1	2.0	530	1.06	534	750	N/A	8
44	20.05.13 12:19	0.1-0.4-0.1	4.0	1463	1.13	1440	40	N/A	8
45	20.05.13 12:59	0.1-0.4-0.1	5.0	1990	1.01	1980	15	N/A	8
46	20.05.13 13:40	0.1-0.4-0.1	8.0	3880	1.56	3880	10	N/A	8
47	20.05.13 14:33	0.04-0.16-0.04	2.0	498	1.03	488	150	N/A	8
48	20.05.13 15:21	0.04-0.16-0.04	0.75	127.2	0.83	125.5	500	N/A	8
49	21.05.13 10:57	0.04-0.16-0.04	1.5	338.5	0.92	332.5	200	N/A	8
50	21.05.13 14:10	0.04-0.16-0.04	2.5	682	1.04	658	100	N/A	8
51	21.05.13 14:50	0.06-0.24-0.06	0.50	74.0	0.98	73.2	7000	N/A	8
52	21.05.13	0.06-0.24-0.06	2.0	528	1.07	526	150	N/A	8
53	21.05.13 15:31	0.06-0.24-0.06	3.0	928	1.32	923.8	100	N/A	8
54	21.05.13 16:18	0.06-0.24-0.06	4.0	1384	1.11	1368	40	N/A	8

Appendix C

Codes

C.1 Profile Measurement

C.1.1 Data Acquisition - ACL script

```
31 # define scan parameters
2 #+description
                                                    32 # mix = minimum value of horizontal correctors[A]
3 #
                                                    33 # max = maximum...
4 # Measurement of current-density electron beam
                                                    34 # miy = minimum value of vertical correctors[A]
                                                    35 # may = maximum...
   profile
                                                    36 # step = step size [A]
5 # in electron lens test stand (linac lab)
6 # Raster scan of magnetic correctors, recording
                                                    37 # dt = time spent at each point [s]
   pinhole current and vacuum
                                                    38 input/user/float/default=-1 inmix "Min. H value (
                                                        mix) [A]"
  #
                                                    39 mix = <inmix>
8 # v1 - Giulio Stancari, 18 Dec 2012
                                                    40 input/user/float/default=1 inmax "Max. H value (
9 #
10 #-description
                                                        max) [A]"
41 max = <inmax>
                                                    42 input/user/float/default=-1 inmiy "Min. V value (
12
13 # initialize
                                                        miy) [A]"
14 fst = newline + newline + "#=== Electron lens
                                                    43 miy = <inmiy>
   test stand profile scan ===" + newline + "# " +
                                                    44 input/user/float/default=1 inmay "Max. V value (
   dateToString()
                                                        may) [A]"
15 print fst
                                                    45 may = <inmay>
                                                    46 input/user/float/default=0.5 instep "Step size [A
16 print/output=file:scan.txt fst
17
                                                        1"
18 # define list of horizontal and vertical
                                                    47 step = <instep>
                                                    48 input/user/float/default=1 indt "Time interval [s
   correctors
19 # (only 6 --7th and 8th H and V correctors are
                                                        ייך י
                                                    49 dt = <indt>
   not wired)
20 deviceList/create hcor devices='Z:HCOR02,Z:HCOR06 50 fst = "# Scan parameters:" + newline + "# mix = "
                                                         + toString(mix,8) + " A" + newline + "# max = "
   ,Z:HCORB2,Z:HCOR10,Z:HCOR14,Z:HCORB6'
                                                         + toString(max,8) + " A" + newline + "# miy = "
21 deviceList/create vcor devices='Z:VCOR01,Z:VCOR05
                                                         + toString(miy,8) + " A" + newline + \
   ,Z:VCORB1,Z:VCOR09,Z:VCOR13,Z:VCORB5'
                                                         "# may = " + toString(may,8) + " A"+newline+"
22
                                                          # step = "+toString(step,8)+" A"+newline+"# dt
23 # define list of variables to read
24 # Z:BC1AD3 = probe current
                                                           = "+toString(dt,8)+" s"
25 # Z:BC1AD5 = tube current
                                                    52 print fst
26 # Z:BC2AD5 = diode pump vacuum
                                                    53 print/output=file:scan.txt fst
27 deviceList/create devin devices='Z:BC1AD3,Z:
                                                    54
   BC1AD5,Z:BC2AD5,Z:HCORO2,Z:HCORO6,Z:HCORB2,Z:
                                                    55 # validating scan parameters
   HCOR10,Z:HCOR14,Z:HCORB6,Z:VCOR01,Z:VCOR05,Z:
                                                    56 if (mix < -5) or (max > 5) or
   VCORB1,Z:VCOR09,Z:VCOR13,Z:VCORB5'
                                                          (miy < -5) or (may > 5) or \setminus
                                                    57
28 deviceList/size/output=variable:ndevin devin
                                                          (mix > max) or (miy > may) or \setminus
                                                    58
29 declare rd float[15]
                                                    59
                                                          (step <= 0) or (step > 5) or \setminus
                                                          (dt <= 0) or (dt > 600)
30
                                                    60
```

```
exit "Invalid scan parameters."
                                                              else
                                                        101
62 endif
                                                        102
                                                               y = may
63
                                                        103
                                                              endif
64 # calculating scan duration
                                                             while (y >= miy) and (y <= may)</pre>
                                                        104
65 \text{ nx} = (\text{max-mix})/\text{step} + 1
                                                               ramp deviceList=vcor y
                                                        105
66 \text{ ny} = (\text{may-miy})/\text{step} + 1
                                                               wait/sec dt
                                                        106
67 \text{ dur} = \text{nx*ny*dt}
                                                        107
                                                               read_list/nodisplay/values=rd device_list=
68 print "Estimated scan duration: " dur " s"
                                                                devin
                                                        108 # format file output
69
70 # initialize direction of raster scan
                                                               fst = " "+toString(x,6)+" "+toString(y,6)
                                                        109
71 # stepsign = +1 -> positive
                                                               loop ndevin i
                                                        110
                                                                fst = fst + " " + toString(rd[i],12)
72 # stepsign = -1 -> negative
                                                        111
73 stepsign = 1
                                                               endloop
                                                        112
                                                               fst = fst + " " + toString(ctimeNow)
74 enable settings
                                                        113
75 print "Ramping correctors to initial settings..." 114
                                                               print/output=file:scan.txt fst
76 x = mix
                                                        115 # format screen output
77 \mathbf{y} = \mathbf{miy}
                                                               print pos(2) width(x,7) \
                                                        116
78 ramp/duration=2 deviceList=hcor x
                                                                    pos(10) width(y,7) 
                                                        117
79 ramp/duration=2 deviceList=vcor y
                                                        118
                                                                     pos(20) width(rd[0],12) \
                                                                   pos(35) width(rd[1],11) \
80 print "Starting raster scan..." newline newline
                                                        119
81
                                                        120
                                                                     pos(50) width(rd[2],10)
                                                        121 # increase or decrease vertical setting
82 # screen header
83 print pos(2) " x [A] " \
                                                               y = y + stepsign*step
                                                        122
       pos(10) " y [A] " \
                                                        123 endwhile
84
        pos(20) "probe [arb.]" \
                                                        124 # revert direction of vertical scan
85
        pos(35) "tube [arb.]" \
                                                        125 stepsign = -1 * stepsign
86
        pos(50) "vacuum [V]"
                                                        126 x = x + step
87
                                                        127 endwhile
88
89 # file header
                                                        128 exit
90 fst = newline + newline + "# xset [A] yset [A]"
                                                        129
                                                        130 # always clean up
91 loop ndevin
    fst=fst+" "+name(Q:DEVIN)+" ["+units(Q:DEVIN)+" 131 always:
     ייך -
                                                             print "Ramping correctors back to zero..."
                                                        132
    deviceList/inc devin
                                                             ramp/duration=2 deviceList=hcor 0
93
                                                        133
                                                             ramp/duration=2 deviceList=vcor 0
94 endloop
                                                        134
95 fst = fst + " ctime [s]"
                                                             fst = newline + "# === End of profile scan ==="
                                                        135
96 print/output=file:scan.txt fst
                                                               + newline + \setminus
97 while x <= max
                                                               "# " + dateToString()
                                                        136
    ramp deviceList=hcor x
                                                             print fst newline newline
                                                        137
98
                                                             print/output=file:scan.txt fst
99
    if stepsign = 1
                                                        138
     y = miy
100
```

C.1.2 Data Processing - R script

```
"../Data/HG/HG_091023_775A_250V_303030kG_16mA.
                                                16
2 # NAME: profile_amalysis.R
                                                     dat", #1
3 # AUTHOR: Vince Moens, Giulio Stancari
4 # LAST MODIFIED: 05.07.2013
                                                                        . . .
5 # PURPOSE: Profile Calibration, Particle Export,
   Transverse Profile Calculation, Radial Profile 152
                                                    "../Data/HG1/HG1b_130521_9p25A_06-24-06kG_4kV_
   Calculation, Polar Decomposition
                                                     1368mA.txt.gz" #136
153)
                                               154
8 # Initialization
                                               155 ### CONTROL SWITCHES ###
9 library(fields)
                                               156 profile.list <- 99:100 #length(fns)
10 require(MASS); require(rgl); require(RSEIS)
                                               157
11 T <- TRUE; F <- FALSE
                                               158 CORRECT.SCALLOPS <- T
12 par(ask=FALSE)
                                               159 FIND.SCALLOP.CORRECTION.PAR <- F
                                               160 SUBTRACT.BACKGROUND <- T
13
                                               161 CORRECT.BOTTOM.LEFT <- T
14 # Profile Loading
15 fns <- c(
                                               162 NORMALIZE.PROFILE <- T
                                               163 CALIBRATE.CORRECTORS <- F
```

```
164 MAKE.PLOTS <- T
                                                         207
165 EXPORT.CALIBRATED.PROFILE <- T
                                                         208
166 EXPORT.RADIAL <- T
                                                         209
167 EXPORT.PARTICLES <- T; Nmax <- 8192
                                                         210
168 POLAR.DECOMPOSITION <- T
                                                         211
169 RUBBER.STAMP <- T
                                                         212
170 BKGR.BLACK <- F
                                                         213
171 BKGR.TRANS <- T
                                                         214
172 COLOR.CHART <- T
                                                         215
173 D3.AXIS <- T
                                                         216
174
175 if(COLOR.CHART) {out.dir <- "../Results/Chart_
                                                         217
     Colors/"
                                                         218
                     polar.dir <- "../Polar/Results/</pre>
176
                                                         219
                      Chart_Colors/"
                                                         220
177 } else {out.dir <- "../Results/Matlab_Colors/"</pre>
                                                         221
           polar.dir <- "../Polar/Results/Matlab_</pre>
                                                         222
178
            Colors/"
                                                         223
179 }
                                                         224
180 radial.dir <- "../Radial/Results/"</pre>
                                                         225
181 fname.id <- "test"</pre>
                                                         226
182
183 # log file
                                                         227
184 logfile <- "log.txt"</pre>
185 unlink(logfile)
                                                         228
186 sink(logfile, split=TRUE)
                                                         229
187 cat("\nAnalysis of electron beam profiles\n")
                                                         230
188 cat("Start: ", date(), "\n")
                                                         231
189
                                                         232
190 # plot options
191 if(BKGR.BLACK) par(bg="black",fg="white",col="
                                                         233
     white", col.axis="white", col.lab="white", col.
                                                         234
     main="white", col.sub="white")
                                                         235
192 if(BKGR.BLACK) pbgr <- "black" else pbgr <- "</pre>
                                                         236
     white"
193 if(BKGR.TRANS) par(bg="transparent")
                                                         237
194 if (BKGR.TRANS) pbgr <- "transparent"
                                                         238
195 pwid <- 3.25; phei <- 3.25
                                                         239
196 reso <- 300; pwpx <- pwid*reso; phpx <- phei*reso 240
197 resoRGL <- 300; pwpxRGL <- pwid*resoRGL; phpxRGL 241
     <- phei*resoRGL
                                                         242
198 pfon <- "Times"; ptsz <- 10
                                                         243
199 quartz.options(width=pwid, height=phei, family=
                                                         244
     pfon,pointsize=ptsz,bg=pbgr,reset=TRUE)
                                                         245
200 pdf.options(width=pwid,height=phei,family=pfon,
                                                         246
    pointsize=ptsz,bg=pbgr,reset=TRUE)
                                                         247
201 setEPS(width=pwid,height=phei,family=pfon,
                                                         248
    pointsize=ptsz,bg=pbgr,reset=TRUE)
                                                         249
202 matlab.colors <- colorRampPalette( c("#00007F","</pre>
                                                         250
     blue","#007FFF","cyan","#7FFF7F","yellow","#
                                                         251
     FF7F00", "red", "#7F0000"))
203
                                                         253
204 # filled contour with no key
205 filled.contour.nokey <- function (x=seq(0,1,
    length.out=nrow(z)),y=seq(0,1,length.out=ncol(z) 256 Xcorrl <- c( -1.210455, -2.381263, -3.072298, -
     ),z,xlim=range(x,finite=TRUE),ylim = range(y,
     levels = pretty(zlim, nlevels), nlevels = 20,
     color.palette = cm.colors,col = color.palette(
                                                        258
     length(levels) -1), plot.title, plot.axes,key.
     title, key.axes, asp = NA, xaxs = "i", yaxs = "i 260 # filament heater current [A]
     ", las = 1,axes = FALSE, frame.plot = axes, ...) 261 IH <- c(rep(7.75,20), rep(5.79,2), rep(7.75,6),
206 {
```

```
if (missing(z)) {
                                                              if (!missing(x)) {
                                                                  if (is.list(x)) {
                                                                     z <- x$z
                                                                     y <- x$y
                                                                     x <- x$x
                                                                  }
                                                                  else {
                                                                     z <- x
                                                                     x <- seq.int(0, 1, length.out =</pre>
                                                                      nrow(z))
                                                                  }
                                                              }
                                                              else stop("no 'z' matrix specified")
                                                          }
                                                          else if (is.list(x)) {
                                                              y <- x$y
                                                              x <- x$x
                                                          }
                                                          if (any(diff(x) \le 0) || any(diff(y) \le 0))
                                                              stop("increasing 'x' and 'y' values
                                                               expected")
                                                          mar.orig <- (par.orig <- par(c("mar", "las", "</pre>
                                                           mfrow")))$mar
                                                          on.exit(par(par.orig))
                                                          par(las = las, xaxs = "i", yaxs = "i")
                                                          plot.new()
                                                          plot.window(xlim, ylim, "", asp = asp)
                                                          if (!is.matrix(z) || nrow(z) \leq 1L || ncol(z)
                                                           <= 1L)
                                                              stop("no proper 'z' matrix specified")
                                                          if (!is.double(z))
                                                              storage.mode(z) <- "double"</pre>
                                                           .Internal(filledcontour(as.double(x), as.
                                                           double(y), z, as.double(levels),
                                                              col = col))
                                                          if (missing(plot.axes)) {
                                                              if (axes) {
                                                                  title(main = "", xlab = "", ylab = "")
                                                                  Axis(x, side = 1)
                                                                  Axis(y, side = 2)
                                                              }
                                                          }
                                                          else plot.axes
                                                          if (frame.plot)
                                                              box()
                                                          if (missing(plot.title))
                                                              title(...)
                                                          else plot.title
                                                          invisible()
                                                   252 }
                                                   254 # Centering corrections
                                                   255 Seq <- c(1.6, 2.4, 3.2, 4)
                                                       3.208143)
finite = TRUE), zlim = range(z, finite = TRUE), 257 Ycorrl <- c( -0.4880404, -0.6774335, -0.926263, -
                                                       1.03759)
                                                  259 # experimental conditions
                                                       rep(4,2), rep(8.75, 17), rep(6.25, 5), rep(8.75,
```

2), rep(9.25, 10), rep(7.25, 2), rep(9.25, 30), 274 caly.v <- rep(0.14, length(fns)) 275 calx.v[which(NC==3)] <- 0.1285 # checked rep(10.0,4),rep(9.25,36)) 262 # cathode voltage [V] 0.1526 # NC=3 but different 276 calx.v[20] <-263 V <- c(250, 500, 1000, 1500, 2000, 3000, 4000, settings 5000, 6000, 7500,9000, 3e3, 3e3, 9e3, 9e3, 9e3, 277 calx.v[which(NC==4)] <- 0.1313 # checked 250, 250, 1e3, 6e3, 500, 7500, 500,2000, 7500, 278 calx.v[which(NC==6)] <- 0.1706 # checked 7500, 50, 2000, 100, 5000,500, 1e3, 2e3, 4e3, 8 279 calx.v[29:30] <- 0.199 # 3.5-4.0-1.5 kG, Gaussian e3, 1e3, 2e3, 500, 4e3, 8e3, 500,500, 1e3, 2e3, qun 4e3, 8e3, 3e3,8e3, 2e3, 500, 500, 500, 8e3, 500, 280 caly.v[which(NC==3)] <- 0.1389 # checked 1e3, 5e3,500, 8e3, 4e3, 2e3, 1e3, 6e3, 3e3, 7e3 281 caly.v[20] <-0.1436 # NC=3 but different , rep(8e3,2), 889, 2e3, 3e3, 3556, 5556, 5000, settings 2083, 3472, 5000, 1333, 2222, 3111, 750, 2000, 282 caly.v[which(NC==4)] <- 0.1450 # checked 1250, rep(500,23) ,1e3 ,2e3, 4e3, 8e3, 7e3, 250, 283 caly.v[which(NC==6)] <- 0.1793 # checked 125, 3e3, 5e3, 6e3, 1e3, 3125, 6250, rep(1e3,5) 284 caly.v[29:30] <- 0.216 # 3.5-4.0-1.5 kG, Gaussian , 500, 500, 2000, 4000, 5000, 8000, 2000, 750, qun1500, 2500, 500, 2000, 3000, 4000) 285264 # peak beam current at collector [A] 286 # offsets [A] 265 I <- c(0.016, 0.044, 0.121, 0.220, 0.325, 0.582, 287 xoff.v <- rep(0, length(fns)) 0.876, 1.180,1.490, 2.020, 2.490, 0.564, 0.600, 288 yoff.v <- rep(0, length(fns)) 2.330, 2.470, 2.550, 0.016,0.015, 0.118, 1.46, 289 xoff.v[which(NC==4)] <- -0.4109 0.023, 0.074, 0.044, 0.330, 2.040, 1.984, 1460e- 290 xoff.v[which(NC==3)] <- c(-0.0199, -0.0895, -6,0.348, 0.00198, 1.670,0.050, 0.134, 0.386, 0.3217, -0.0787, -0.0081, -0.1633, 0) # NC=3 1.030, 2.850, 0.134, 0.394, 0.049, 1.064,2.860, 291 xoff.v[20] <- 0.5942 # NC=3 special case 0.050,0.050, 0.135, 0.376, 1.056, 2.940, 292 xoff.v[which(NC==6)] <- -0.3265 # NC=6 0.692,0.140, 0.098, rep(0.048, 3), 2.860, 0.050, 293 xoff.v[which(NC==8)] <- -0.1772 # NC=8 294 xoff.v[29:30] <- -1.2 # 3.5-4.0-1.5 kG 0.004, 0.013,0.051, 2.94, 1.1, .388, .14, 1.924, .694, 2.394, .440, .027,0.118, .398, 295 xoff.v[c(31:35,47:54,57:66)] <- -0.419 # 1-in .726, .928, 1.770, 1.484, .412, .865, 1.460, hollow gun, 3 kG .213, .448, 0.744, .090, .390, .192, rep(.051,3) 296 xoff.v[c(36:40,67:69)] <- -0.114 # 1-in hollow , rep(.049,5), rep(.051,6), rep(0.050,5), rep qun, 1 kG (0.070,2), rep(0.073,2), 0.195, 0.526, 1.45, 297 xoff.v[41] <- -0.58 # 1-in hollow gun, 4 kG, to 3.7, 3.1, 0.026, 0.009, 0.944, 1.99, 2.53, be verified 0.195, 1.022, 2.754, 0.196, rep(0.195,4), 0.072, 298 xoff.v[c(42:46, 55:56,70)] <- -0.23 # 1-in hollow 0.073, 0.0534, 1.44, 1.98, 3.88, 0.488, 0.125, gun, 2 kG 0.332, 0.658, 0.073, 0.526, 0.924, 1.368) 299 yoff.v[which(NC==4)] <- -0.1157 300 yoff.v[which(NC==3)] <- c(-0.0533, -0.0872, -266 # solenoidal field [T] 267 B <- c(rep(.3,11), 0.1, 0.4, 0.1, 0.2, 0.4, 0.1, 0.0786, -0.0469, -0.0330, -0.0582, 0) # NC=3 0.033, 0.133,0.223, rep(0.3,7), 0.4, rep(0.4,2), 301 yoff.v[20] <- 0.1957 # NC=3 special case rep(0.3, 5), rep(0.1, 5), 0.4, rep(0.2, 5), 0.3, 302 yoff.v[which(NC==6)] <- -0.1000 # NC=6 rep(0.3,7), rep(0.2,2), rep(0.3,10), rep(0.1,3) 303 yoff.v[which(NC==8)] <- -0.0510 # NC=8 , 0.2, rep(0.25,4), 0.3, rep(0.2,3), rep(0.15,3) 304 yoff.v[29:30] <- -0.40 # 3.5-4.0-1.5 kG , rep(0.25,3), rep(0.2,5), rep(0.3,13), rep 305 yoff.v[c(31:35,47:54,57:66)] <- 0.081 # 1-in (0.32,12), rep(0.4,3), 0.24, rep(0.16,5), rep hollow gun, 3 kG (0.4,5), rep(0.16,4), rep(0.24,4)) 306 yoff.v[c(36:40, 67:69)] <- 0.033 # 1-in hollow 268 # number of correctors gun, 1 kG 269 NC <- c(rep(4,11), 3, 6, 3, 3, 6, 3, 3, 3, 3, 3, rep 307 yoff.v[41] <- 0.067 # 1-in hollow gun, 4 kG, to (6,7), 4, rep(4,2),rep(6, 106)) be verified 270 # raster-scan correction Parameters 308 yoff.v[c(42:46, 55:56,70)] <- 0.07 # 1-in hollow 271 besttp <- c(-0.6069, -0.5819, -0.7744, -0.8363, gun, 2 kG 0.5632, -0.5392, -0.3877, -1.017, -0.5627, -309 0.5051, -0.7392, 0.0392, -1.9296, -0.1264, -310 # inner (r1) and outer (r2) qun radii [m]0.3909, -1.631, -0.2347, -0.2685, -0.1528, -311 r1 <- c(rep(0.00450, 28), rep(0.00000, 2), rep (0.00675, 106))0.0197, -1.7174, -0.7895, -1.2949, -1.5334, -312 r2 <- c(rep(0.00762, 28), rep(0.00508, 2), rep 1.6833, -1.2, -1.19, -4.02, -0.37, -0.44, -0.3901180, -0.4279687, -0.4502339, -0.3418769, -(0.01270, 106)) 0.2973467,-0.3418769, -0.3203540, -0.3233227, - 313 xylim.fudge <- 1.2 # fudge factor for xy limits = 0.2928937, -0.3671107, -0.4932798, -0.4583977, r2 * xylim.fudge0.4895689, -0.4784364, -0.2832455, -0.2832455, -3140.3878915, -0.350783, rep(-0.4, 6), rep(-0.62, 315 # constants 2), rep(-0.62, 8),-.18,-.62, -0.3418769, -316 **c** <- 299792458 0.3203540, -0.3, rep(-0.2832455,67)) 317 eps0 <- 8.854187817e-12 272 # calibration constants [mm*T/A/corr] 318 me <- 9.1093826e-31 273 calx.v <- rep(0.13, length(fns))</pre> 319 q <- 1.60217653e-19

```
_{320} L <- rep(2.86, length(fns)) # anode-collector
                                                         374
     distance [m]
321 # electron velocities
                                                         375
322 gamma <- 1 + q*V/(me*c^2)
323 beta <- sqrt(gamma*gamma-1)/gamma
                                                         376
324 v <- beta*c
325 # flight times
                                                         377
326 t <- L/v
327 # average densities
                                                         378
328 lambda <- I/(q*v) # linear
329 n <- lambda/(pi*(r2*r2-r1*r1)) # volume
                                                         379
330 # plasma frequencies
331 wp <- sqrt(n*q*q/(eps0*me))
                                                         380
332 # cyclotron frequencies
                                                         381
333 wc <- q*B/me
334 # diocotron frequencies
                                                         382
335 wD <- wp*wp/(2*wc)
336 # evolution numbers
                                                         383
337 nev <- wD*t/(2*pi)
338 # assemble data frame
                                                         384
339 DSET <- data.frame(fns, IH, B, V, I, NC, besttp,
     calx.v, caly.v,xoff.v, yoff.v, r1, r2, L, beta,
                                                         385
     t, lambda, n, wc,wp, wD, nev)
340 write.table(DSET, file=paste(out.dir, fname.id, "386
     _data_set.txt", sep=""))
341
                                                         387
342 # loop over profiles
                                                         388
343 for(l in profile.list) {
                                                         389
     if(B[1]==0.16){
344
                                                         390
       Xcorr=Xcorr1[1]
345
       Ycorr=Ycorr1[1]
346
                                                         391
     }
347
                                                         392
     else if(B[1]==0.24){
348
       Xcorr=Xcorr1[2]
349
                                                         393
       Ycorr=Ycorr1[2]
350
351
     }
                                                         394
     else if(B[1]==0.32){
352
                                                         395
       Xcorr=Xcorr1[3]
353
       Ycorr=Ycorr1[3]
354
                                                         396
     }
                                                         397
355
     else if(B[1]==0.4){
                                                         398
356
       Xcorr=Xcorr1[4]
357
                                                         399
       Ycorr=Ycorr1[4]
358
                                                         400
     }
359
                                                         401
     else {
360
       Xcorr=0
361
       Ycorr=0
362
                                                         402
     }
363
                                                         403
     fname.id <- substr(fns[1],13,nchar(fns[1])-7) #</pre>
364
       filename prefix
                                                         404
     message("fname = ", fname.id)
365
                                                         405
     graphics.off()
366
                                                         406
     if(rgl.cur()>0) rgl.close()
367
368
                                                         407
      # read scan data
369
                                                         408
     message("\n\n\n==== Reading file", fns[1], "
370
      ...=====")
     cat("Hollow gun inner radius:", format(r1[1]*1 409
371
      e3, di=3), "mm\n")
                                                         410
     cat("Hollow gun outer radius:", format(r2[1]*1
                                                         411
372
      e3, di=3), "mm\n")
                                                         412
     cat("Anode-collector distance:", format(L[1],
373
                                                         413
      di=3), "m\n")
                                                         414
```

```
cat("Filament heater current:", format(IH[1],di
 =3), "A \ n")
cat("Solenoidal field:", format(B[1]*10,di=4),
 kG(n)
cat("Cathode voltage:", format(V[1]/1e3,di=4),
 "kV\n")
cat("Peak collector current:", format(I[1]*1e3,
 di=4), "mAn")
cat("Number of correctors:", format(NC[1]), "\n
 ")
cat("Electron velocity:", format(v[1],di=3), "m
/s (beta =",format(beta[1],di=3), ")\n")
cat("Flight time:", format(t[1],di=3), "s\n")
cat("Linear number density:", format(lambda[1],
 di=3), "/m n")
cat("Volume number density:", format(n[1],di=3)
 , "/m3\n")
cat("Cyclotron frequency:", format(wc[1]/(2*pi)
 ,di=3), "Hz\n")
cat("Plasma frequency:", format(wp[1]/(2*pi),di
 =3), "Hz n")
cat("Diocotron frequency:", format(wD[1]/(2*pi)
 ,di=3), "Hz\n")
cat("Evolution number:", format(nev[1],di=3), "
rev/flight\n")
# Finding the number of data points to plot.
message("Obtaining scan parameters...")
xraw <- round(D$V1, digits=3) # horz. deflector</pre>
 , rounded
Xmin <- min(xraw); Xmax <- max(xraw)</pre>
Nx <- nlevels(as.factor(xraw)); Xstep <- (Xmax-</pre>
 Xmin)/(Nx-1)
yraw <- round(D$V2, digits=3) # vert. deflector</pre>
 , rounded
Ymin <- min(yraw); Ymax <- max(yraw)</pre>
Ny <- nlevels(as.factor(yraw)); Ystep <- (Ymax-
 Ymin)/(Ny-1)
# define grid
X <- seq(Xmin, Xmax, length.out=Nx)
Y <- seq(Ymin, Ymax, length.out=Ny)</pre>
cat("X scan: from", Xmin, "to", Xmax, "(", Nx,
 "points)\n")
cat("Y scan: from", Ymin, "to", Ymax, "(", Ny,
 "points)\n")
# new profiles have vacuum data in 5th column (
 ACL scripts)
VACUUM <- FALSE
if(1>=57) VACUUM <- TRUE
# calibration of 20-l/s Tevatron diode pump
 aauae
Vga <- c(9,8.4,8,7,6.9,6,5.4,5,4,3.85,3) # V
logPtorr <- log10(c(4e-10, 1e-9, 2e-9, 8.5e-9,
 1e-8, 4e-8, 1e-7, 1.8e-7, 8e-7, 1e-6, 5.5e-6)
 ) # [torr], log scale
logPvsVsm <- smooth.spline(Vga, logPtorr)</pre>
PvsV <- function(V=6){</pre>
 Vmin <- 2.8; Vmax <- 9.9
 P <- 1e-4 + 0∗V
 P[V>=Vmax] <- 1e-10
```

P[V<=Vmin] <- 1e-4

```
Appendix C. Codes
```

```
good <- Vmin<V & V<Vmax
415
                                                             471
        P[good] <- predict(logPvsVsm, x=V[good])$y</pre>
416
                                                             472
        P[good] <- 10^P[good] * 1.333 # [mbar]
417
                                                             473
        return(P)
418
                                                             474
      }
419
                                                             475
     message("Defining profile matrices...")
420
                                                             476
      Zprobe <- matrix(0, nrow=Nx, ncol=Ny)</pre>
421
      Ztube <- matrix(0, nrow=Nx, ncol=Ny)</pre>
422
                                                             477
      Zvac <- matrix(NA, nrow=Nx, ncol=Ny)</pre>
423
      # assign value according to position in grid
424
                                                             478
      for(k in 1:length(D$V3)) {
425
                                                             479
        i <- round((xraw[k]-Xmin)/Xstep) + 1</pre>
426
                                                             480
        j <- round((yraw[k]-Ymin)/Ystep) + 1</pre>
427
        Zprobe[i,j] <- D$V3[k]</pre>
                                                             481
428
        Ztube[i,j] <- D$V4[k]</pre>
429
                                                             482
        # new profiles have vacuum data
430
        if(VACUUM) Zvac[i,j] <- PvsV( D$V5[k] )</pre>
431
                                                             483
      }
                                                             484
432
433
                                                             485
      # choose which variable to analyze and plot
                                                             486
434
      Z <- Zprobe
                                                             487
435
436
                                                             488
437
      # correct for scallops due to raster scan?
      if(CORRECT.SCALLOPS){
438
                                                             489
       message("Correcting for scallops due to raster
439
                                                             490
         scan...")
                                                             491
       # find best correction parameter?
440
                                                             492
       if(FIND.SCALLOP.CORRECTION.PAR){
441
        message("Looking for best correction parameter493
442
         ...")
                                                             494
        # tpval <- seq(-2, 0, by=0.05)
443
                                                             495
       tpval <- seq(besttp[1]-0.2,besttp[1]+0.2,by</pre>
444
                                                             496
        =0.05)
       hfc <- 0*tpval
445
                                                             497
       for(t in 1:length(tpval)) {
                                                             498
446
447
        tp <- tpval[t]</pre>
                                                             499
        Ztmp <- Z
448
                                                             500
        for(i in seq(1,Nx)) {
449
                                                             501
         if(floor(i/2) == (i/2)) \{ \# odd rows \}
450
                                                             502
          for(j in 2:(Ny)) {
451
                                                             503
           Ztmp[i,j] <- Z[i,j-1] + (Z[i,j] -Z[i,j-1])
452
            / (1-tp) }}
                                                             504
         else{ # even rows
453
                                                             505
          for(j in 1:(Ny-1)) {
454
                                                             506
           Ztmp[i,j] <- Z[i,j+1] + (Z[i,j] -Z[i,j+1])
                                                             507
455
            / (1-tp) }}}
                                                             508
          # evaluate scallop frequency component
456
                                                             509
          # by looking at FFT of aligned rows
457
                                                             510
          if(F){
458
          if(floor(Nx/2) != (Nx/2)) {
459
                                                             511
            Zv <- Ztmp[1:(Nx-1),1]
460
                                                             512
            for(j in 2:Ny){Zv <- c(Zv,Ztmp[1:(Nx-1),j 513</pre>
461
             ])}}
          else{
462
                                                             514
            Zv <- Ztmp[1:Nx,1]
463
            for(j in 2:Ny){Zv <- c(Zv,Ztmp[1:Nx,j])} } 516</pre>
464
          lFZv <- length(Zv)</pre>
465
                                                             517
          win <- get.slepians(lFZv, 1, 3)[,1]</pre>
466
                                                             518
          FZv <- abs(fft(Zv*win))^2</pre>
467
          centr <- 1FZv/2
                                                             519
468
          hwdth <-3
469
          hfc[t] <- sum(FZv[floor(centr-hwdth):ceiling520</pre>
470
           (centr+hwdth)])
                                                             521
```

```
}
    # by looking at spatial FFT
    if(T){
    icmtmp <- sum(t(1:Nx) * Ztmp) / sum(Ztmp)</pre>
    jcmtmp <- sum(Ztmp * 1:Ny) / sum(Ztmp)</pre>
    scallop.filter <- abs((1:Nx-icmtmp) %o% (1:</pre>
    Ny-jcmtmp))
    scallop.filter <- scallop.filter / max(</pre>
    scallop.filter)
    scallop.filter <- 1 -scallop.filter</pre>
    scallop.filter <- (scallop.filter>0.8)
    win <- get.slepians(Nx, 1, 3)[,1] %0% get.
    slepians(Ny, 1, 3)[,1]
    FZ <- Mod(fft(Ztmp*win))</pre>
    hfc[t] <- sum( FZ[scallop.filter] ) / sum(</pre>
    FZ[!scallop.filter] )
    }
    # "maximum entropy"
    if(F){
    Zprb <- Ztmp / sum(Ztmp)
    hfc[t] <- sum(Zprb)*log(sum(Zprb))</pre>
    for(i in 1:Nx){ for(j in 1:Ny){hfc[t] <- hfc</pre>
     [t] -Zprb[i, j]*log(Zprb[i, j]) }}
   hfc[t] <- -hfc[t]</pre>
   }
 }
 cat("Please select minimum of figure of merit
   (1 \text{ click}) \setminus n")
 par(mar=c(5,4,4,2)+0.1)
 plot(tpval, hfc, type="p", pch=16)
  spf <- splinefun(tpval, hfc)</pre>
  curve(spf(x), min(tpval), max(tpval), add=TRUE
  )
 tp <- locator(1)$x</pre>
 besttp[1] <- tp</pre>
}
else {
 tp <- besttp[1]</pre>
}
cat("Raster-scan correction parameter:", tp, "\
n")
# correct profile
Zc < - Z
for(i in seq(1,Nx)) {
if(floor(i/2) == (i/2)) \{ \# odd rows
 for(j in 2:(Ny)) {
  Zc[i,j] <- Z[i,j-1] + (Z[i,j] -Z[i,j-1]) / (1)
    -tp) }}
 else{ # even rows
 for(j in 1:(Ny-1)) {
  Zc[i,j] <- Z[i,j+1] + (Z[i,j] -Z[i,j+1]) / (1
    -tp) }}}
Z <- Zc
}
# fix 250-V profile (power supply trip)?
if(T & fns[1]=="../Data/HG/HG_091023_775A_250V_
 303030kG 16mA.dat") {
 message("Correcting profile with HVPS trip..."
 itrip <- 63; ic <- 54; ioff <- 1; imag <- 1
  jc <- 52.8; joff <- 3; jmag <- 1
```

```
Zc <- Z
522
                                                            575
        for(riga in itrip:min(Nx,(Nx-ioff))) {
523
          for(colo in 1:min(Ny,(Ny-joff))) {
524
                                                            577
            ixcor <- ic+(riga-ic)*imag+ioff</pre>
                                                            578
525
            jycor <- jc+(colo-jc)*jmag+joff</pre>
526
                                                            579
            icor <- round(ixcor)</pre>
527
                                                            580
            if(icor < itrip) icor <- 1</pre>
528
            if(icor > Nx) icor <- Nx</pre>
529
                                                            581
            jcor <- round(jycor)</pre>
530
                                                            582
            if(jcor < 1) jcor <- 1</pre>
531
                                                            583
            if(jcor > Ny) jcor <- Ny</pre>
532
                                                            584
            Zc[riga,colo] <- Z[icor,jcor]</pre>
                                                            585
         }
                                                            586
       }
535
                                                            587
       Z <- Zc
536
                                                            588
      3
                                                            589
                                                            590
      # subtract background?
                                                            591
539
      if (COLOR.CHART & SUBTRACT.BACKGROUND) {message( 592
540
       "Subtracting background..."); Z <- Z -mean(Z[Z 593
       <=as.numeric(quantile(Z,probs=0.25))]);Z[Z<0]< 594
       -0}
      else if(!COLOR.CHART & SUBTRACT.BACKGROUND){
541
                                                            595
       message("Subtracting background..."); Z <- Z - 596</pre>
       mean(Z[Z<=as.numeric(quantile(Z,probs=0.25))])</pre>
                                                           597
       } #Should we average over the few minimum
                                                            598
       places?
542
                                                            599
      # eliminate fake signals at bottom LEFT
543
      if(CORRECT.BOTTOM.LEFT){Z[1:2,1:2] <- min(Z)}</pre>
544
545
                                                            601
      # normalize?
546
      Zu <- Z # save unnormalized profile
547
                                                            603
      if(NORMALIZE.PROFILE){message("Normalizing...")
548
                                                           604
       ; Z <- Z/(sum(Z))
549
                                                            606
      # find "center of mass" in terms of coordinates 607
550
        and array indices for plotting slices
                                                            608
     message("Finding center of mass...")
                                                            609
551
      xcm <- sum(t(X) * Z) / sum(Z)
                                                            610
552
      icm <- sum(t(seq(1,Nx)) * Z) / sum(Z)
553
                                                            611
      ycm <- sum(Z * Y) / sum(Z)
554
                                                            612
      jcm <- sum(Z * seq(1,Ny)) / sum(Z)
555
                                                            613
      cat("xcm =", xcm, "(row", icm, ")\n")
556
                                                            614
      cat("ycm =", ycm, "(column", jcm, ")\n")
557
558
                                                            615
      # horizontal and vertical scales in mm
559
      # corrector calibration: mm*T/A/ncorr
560
                                                            617
     message("Axis calibration...")
561
                                                            618
      calx <- calx.v[l]; caly <- caly.v[l]</pre>
562
                                                            619
      Xoff <- xoff.v[l]; Yoff <- yoff.v[l]</pre>
563
                                                            620
      if(CALIBRATE.CORRECTORS){
564
       message("Calibration by fitting ellipse...")
565
        rp <- c(r1[1], r2[1], 0)*1e3 # mm
566
                                                            623
        rx <- rep(0,3); ry <- rep(0,3)
567
                                                            624
        xcent <- rep(0,2); ycent <- rep(0,2)</pre>
568
        message <- c("inner", "outer")</pre>
569
                                                            625
        for(k in 1:2){
570
                                                            626
          image(X, Y, Z, col=matlab.colors(64))
571
                                                            627
         np <- 8
                                                            628
          cat("Please choose", np, "points on",
                                                            629
573
          message[k], "edge\n")
                                                            630
          po <- locator(np)</pre>
574
                                                            631
```

```
x <- po$x; y <- po$y
e \leftarrow rep(0.01, length(x)) # errors
rguess <- mean(sqrt((x-xcm)^2+(y-ycm)^2))</pre>
ipar <- c(rguess, rguess, 0, xcm, ycm)</pre>
# parameter scales
pscl <- c(rguess, rguess, pi, 3*Xstep, 3*</pre>
Ystep)
# initial step sizes
dpar <- rep(0.01, 5)
# define function to be optimized
fopt <- function(p) {</pre>
 a2 <- p[1]*p[1]; b2 <- p[2]*p[2]
 st <- sin(p[3]); ct <- cos(p[3])</pre>
 x0 <- p[4]; y0 <- p[5]
 A <- ct*ct/a2 + st*st/b2
 B <- 2*st*ct*(1/a2-1/b2)
 C \leq st*st/a2 + ct*ct/b2
 D <- -2*x0*A -B*v0
 E <- -B*x0 -2*C*y0
 F <- -1+A*x0*x0+B*x0*y0+C*y0*y0
 fopt <- sum((((A*x*x + B*x*y + C*y*y + D*x</pre>
    + E*y)/F + 1)/e)^2)
 return(fopt)
7
# minimize function
fit <- optim(ipar, fopt, hessian=TRUE,</pre>
method="BFGS".
control=list(trace=6, parscale=pscl, ndeps=
dpar,reltol=1e-5, REPORT=1))
# fitted parameters
fpar <- fit$par</pre>
# covariance matrix
emat <- solve(fit$hessian)</pre>
# parameter uncertainties
epar <- sqrt(diag(emat))</pre>
# matrix of correlation coefficients
cmat <- emat
for (i in 1:length(fpar)) {
 for (j in 1:length(fpar)) {
   cmat[i,j] <- emat[i,j]/epar[i]/epar[j]</pre>
 1
}
cat("\n", "Parameters:\n")
pars <- data.frame(fitted=fpar, errors=epar,</pre>
  initial=ipar)
print(round(pars,6))
cat("\n", "Covariance matrix:\n")
write.matrix(emat)
cat("\n", "Correlation matrix:\n")
write.matrix(round(cmat,3))
# Results
a <- fpar[1]; b <- fpar[2]
th <- fpar[3]
xc <- fpar[4]; yc <- fpar[5]</pre>
if(a>=b) {c <- sqrt(a*a-b*b)} else {c <-
sqrt(b*b-a*a)}
ecc <- c/a
cat("\n\nRotation =", th*180/pi, "deg\n")
cat("Axes =", a, "and", b, "n")
cat("Center-foci distance =", c, "\n")
cat("Eccentricity =", ecc, "\n")
cat("Center = (", xc, ",", yc, ")\n")
# draw fitted ellipse
```

```
plot(x, y, pch=16)
632
                                                            679
          phi <- seq(0, 2*pi, length.out=300)
633
          rho <- sqrt(a*a*b*b/(a*a*sin(phi)*sin(phi) +680
634
            b*b*cos(phi)*cos(phi)))
                                                            681
          xx <- rho*cos(phi+th) + xc</pre>
635
                                                            682
          yy <- rho*sin(phi+th) + yc</pre>
636
                                                            683
          points(xx, yy, type="1")
                                                            684
637
          points(xc, yc, pch="+")
638
          # save results for linear interpolation
639
                                                            685
          rx[k] \leftarrow abs(a); ry[k] \leftarrow abs(b)
640
          xcent[k] <- xc; ycent[k] <- yc</pre>
641
        3
642
                                                            686
        CALx <- lm(rp ~ rx); CALy <- lm(rp ~ ry)
643
        plot(range(c(rx,ry)), range(rp), type="n")
644
                                                            687
        points(rx, rp); points(ry, rp)
645
        curve(CALx$coef[1] + CALx$coef[2]*x, min(rx),
646
         max(rx), add=TRUE)
                                                            688
        curve(CALy$coef[1] + CALy$coef[2]*x, min(ry),
                                                            689
647
         max(ry), add=TRUE)
                                                            690
        calx <- CALx$coef[2] * B[1] / NC[1]</pre>
                                                            691
648
        caly <- CALy$coef[2] * B[1] / NC[1]</pre>
                                                            692
649
650
        Xoff <- mean(xcent); Yoff <- mean(ycent)</pre>
651
      }
                                                            694
      cat("Axis calibration coefficients:\n","\tcalx
652
       = ", calx, "mm*T/A/corr\n","\tcaly = ", caly,
                                                            695
       "mm*T/A/corr\n")
      calx.v[l] <- calx
653
                                                            696
      calv.v[1] <- calv</pre>
                                                            697
654
      cat("Offsets:\n","\tXoff = ", Xoff, "A\n","\
655
       tYoff = ", Yoff, "A n"
                                                            698
      xoff.v[l] <- Xoff</pre>
656
                                                            699
      yoff.v[1] <- Yoff</pre>
657
                                                            700
      # calibrated coordinates and step sizes [mm]
658
                                                            701
      Xc <- (X-Xoff) * calx / B[1] * NC[1]</pre>
659
                                                            702
      Yc <- (Y-Yoff) * caly / B[1] * NC[1]
                                                            703
660
      Xcstep <- median(abs(diff(Xc)))</pre>
661
                                                            704
      Ycstep <- median(abs(diff(Yc)))</pre>
662
                                                            705
      # calibrated centers of mass [mm]
663
                                                            706
      Xcm <- (xcm-Xoff) * calx / B[1] * NC[1]</pre>
664
      Ycm <- (ycm-Yoff) * caly / B[1] * NC[1]</pre>
665
      cat("Calibrated center of mass (should be zero) 707
666
       :\n","\tXcm = ", Xcm, "mm\n","\tYcm = ", Ycm,
       "mm\n")
667
                                                            708
      # physical units [A/cm<sup>2</sup>]
668
                                                            709
      Zphys <- Z*I[1] / (Xcstep/10) / (Ycstep/10) #Z/
669
       (Xcstep*Ycstep)
                                                            710
670
      ### profile plots ###
671
                                                            711
      # plot settings
672
      if(COLOR.CHART) {;Zlim <- range(Zphys);Nc <-</pre>
673
                                                            712
       100 # of colors; palette <- matlab.colors(Nc);
       palette <- c(rgb(0,0,0,0),rgb(0,0,0,0.5),rgb
                                                            713
       (0,0,0,1),palette[c(seq(1,29,by=3),seq(30,40,
                                                            714
       by=3), seq(44,56, by=4), seq(60,80, by=2), seq
       (80,100,by=2))]);Nc <-length(palette)}
                                                            715
      else {Zlim <- range(Zphys)</pre>
674
                                                            716
        Nc <- 64 # of colors
675
                                                            717
        palette <- matlab.colors(64)}</pre>
676
                                                            718
      icol <- round((Zphys-Zlim[1])/(Zlim[2]-Zlim[1])</pre>
677
       *(Nc-1)) + 1
                                                            719
      if(COLOR.CHART) icol[icol<=0]=1</pre>
678
```

```
mcol <- matrix(palette[icol], nrow=Nx, ncol=Ny)</pre>
       # color matrix
     # plot 1D slices
     if(MAKE.PLOTS){
       dev.new()
       par(oma=rep(0, 4), mar=c(3.5, 3.5, 0.5, 0.5),
        mgp=c(2, 0.5, 0), tcl=-0.2)
       if(BKGR.BLACK) par(fg="white", col="white",
        col.axis="white", col.lab="white", col.main="
        white",col.sub="white")
       xx <- Xc; yy <- Zphys[,jcm] # horz. along c.m.</pre>
         slice
       plot(xx, yy, type="n", axes=FALSE,xlab="
        Horizontal position [mm]", ylab=expression(
        paste("Current density [", A/cm<sup>2</sup>, "]")))
       axis(1); axis(2, las=1)
       grid(lty="solid")
       box()
       points(xx, yy, pch=16)
       f <- splinefun(xx, yy)</pre>
       curve(f(x), min(xx), max(xx), n=301, add=TRUE)
       file.prefix <- paste(out.dir, fname.id, "1Dx_"
        , sep="")
       fout <- paste(file.prefix,formatC(1, width=2,</pre>
        flag="0"), ".pdf", sep="")
       dev.print(pdf, file=fout)
       fout <- paste(file.prefix,formatC(1, width=2,</pre>
        flag="0"), ".eps", sep="")
       dev.print(postscript, file=fout)
     }
     # --- 2D PLOTS ---#
     if(MAKE.PLOTS){
       # raw 2D measurement
       dev.new()
       par(oma=rep(0,4), mar=c(4.2, 4.2, 0.5, 0.5))
       if(BKGR.BLACK) par(fg="white", col="white",
        col.axis="white", col.lab="white", col.main="
        white", col.sub="white")
       image(X, Y, Zprobe, col=matlab.colors(64), asp
        =1,xlab="H. corrector setting [A]",ylab="V.
        corrector setting [A]")
       # save to file
       file.prefix <- paste(out.dir, fname.id, "_</pre>
        raw2D_", sep="")
       fout <- paste(file.prefix,formatC(1, width=2,</pre>
        flag="0"), ".png", sep="")
       dev2bitmap(file=fout, width=pwid, height=phei,
         res=reso)
       fout <- paste(file.prefix,formatC(1, width=2,</pre>
        flag="0"), ".pdf", sep="")
       dev.print(pdf, file=fout)
       fout <- paste(file.prefix,formatC(1,width=2,</pre>
        flag="0"), ".eps", sep="")
       dev.print(postscript,file=fout)
     }
     # 2D contours
     xylim <- c(-1,1) * xylim.fudge * (r2[1]*1e3) #</pre>
      plot limits [mm]
720 # --- 2D CONTOUR PLOTS ---#
```

```
721 if(MAKE.PLOTS){
```

```
# plot 2D contour, bare
722
                                                         748
       dev.new()
723
                                                         749
       par(oma=rep(0,4), mar=rep(0,4))
724
       if(BKGR.BLACK) par(fg="white", col="white",
725
        col.axis="white", col.lab="white", col.main="
        white", col.sub="white")
                                                         751
       filled.contour.nokey(Xc, Yc, Zphys, xlim=xylim
726
         , ylim=xylim, asp=1,levels=seq(Zlim[1], Zlim 752
        [2], length.out=Nc), col=palette,frame.plot=
        FALSE)
                                                         753
       # save to file
727
       file.prefix <- paste(out.dir, fname.id, "2</pre>
728
                                                         754
        Dbare_", sep="")
                                                         755
       fout <- paste(file.prefix,formatC(1, width=2,</pre>
729
        flag="0"), ".png", sep="")
                                                         756
       dev.print(png, file=fout, width=pwpx, height=
                                                              }
                                                        757
730
        phpx)
                                                         758
       #dev2bitmap(file=fout, width=pwid, height=phei759
731
        , res=reso)
                                                         760
       fout <- paste(file.prefix,formatC(1, width=2,</pre>
                                                        761
732
        flag="0"), ".pdf", sep="")
                                                         762
733
       dev.print(pdf, file=fout)
                                                         763
       fout <- paste(file.prefix,formatC(l,width=2,</pre>
734
                                                         764
        flag="0"), ".eps", sep="")
       dev.print(postscript,file=fout)
735
     }
736
                                                         765
     if(MAKE.PLOTS){
737
       # plot 2D contour, legend.
738
       wiland <- 7; heland <- 6
739
                                                         766
       dev.new(width=wiland, height=heland)
740
                                                         767
       par(oma=rep(0,4), mar=c(4.2,3.2,4.2,0.5), mgp=
741
        c(1.5, 0.5, 0))
                                                         768
       if(BKGR.BLACK) par(fg="white", col="white",
742
        col.axis="white", col.lab="white", col.main="
                                                        769
        white", col.sub="white")
       tick <- diff(range(Xc)) * 0.015; tcol <- "</pre>
743
                                                         770
        white"
       plot.ticks <- function() {axis(1); axis(2);</pre>
744
                                                        771
        lines(c(Xcm,Xcm), c(min(Yc), min(Yc)+tick),
        col=tcol);lines(c(Xcm,Xcm), c(Ycm-tick, Ycm+ 772
        tick), col=tcol);lines(c(Xcm,Xcm), c(max(Yc)-773
        tick,max(Yc)), col=tcol);lines(c(min(Xc),min(
        Xc)+tick), c(Ycm,Ycm), col=tcol);lines(c(Xcm-774
        tick,Xcm+tick), c(Ycm,Ycm), col=tcol);lines(c775
                                                              }
        (max(Xc)-tick,max(Xc)), c(Ycm,Ycm), col=tcol)776
        }
                                                         777
       plot.title <- function() {title(main=bquote(</pre>
745
                                                        778
        paste(I[fil], " = ", .(format(IH[1], nsmall 779
        =2)), " A ","B = ", .(format(B[1]*10, nsmall 780
        =1)), " kG ", "V = ", .(format(V[1]/1e3,
        nsmall=2)), " kV ",I[e], " = ", .(format(I[1]781
        *1e3, nsmall=0)), " mA")),xlab="X [mm]",ylab=
        "Y [mm]")}
       filled.contour(Xc, Yc, Zphys, xlim=xylim, ylim782
746
        =xylim, asp=1,levels=seq(Zlim[1], Zlim[2],
        length.out=Nc), col=palette,plot.title=plot. 783
        title(),plot.axes=plot.ticks(),key.title=
        title(main=expression(paste(j, " [", A/cm<sup>2</sup>,
        "]"))),key.axes=axis(4))
       if(RUBBER.STAMP){mtext(paste(fns[1], "\n","
747
        Analyzed on", date(), " ",R.version.string),
        side = 1, line = 3, adj = 1, font = 3, cex =
```

```
# save to file
 file.prefix <- paste(out.dir, fname.id, "2D_",</pre>
    sep="")
 fout <- paste(file.prefix,formatC(1, width=2,</pre>
  flag="0"), ".png", sep="")
  #dev.print(png, file=fout, width=wiland*reso,
   height=heland*reso)
 dev2bitmap(file=fout, width=wiland, height=
  heland, res=reso)
  fout <- paste(file.prefix,formatC(1, width=2,</pre>
   flag="0"), ".pdf", sep="")
  dev.print(pdf, file=fout)
  fout <- paste(file.prefix,formatC(l,width=2,</pre>
   flag="0"), ".eps", sep="")
  dev.print(postscript,file=fout)
  # --- 2D CENTERED CONTOUR PLOTS ---#
if(MAKE.PLOTS){
  # plot 2D contour, bare, centered
  dev.new()
  par(oma=rep(0,4), mar=rep(0,4))
  if(BKGR.BLACK) par(fg="white", col="white",
   col.axis="white", col.lab="white", col.main="
  white",col.sub="white")
  filled.contour.nokey(Xc-Xcorr, Yc-Ycorr, Zphys
   , asp=1,levels=seq(Zlim[1], Zlim[2], length.
   out=Nc), col=palette,frame.plot=FALSE)
  # save to file
  file.prefix <- paste(out.dir, fname.id, "2</pre>
  Dcentbare_", sep="")
  fout <- paste(file.prefix,formatC(1, width=2,</pre>
  flag="0"), ".png", sep="")
  dev.print(png, file=fout, width=pwpx, height=
  phpx)
  #dev2bitmap(file=fout, width=pwid, height=phei
   , res=reso)
 fout <- paste(file.prefix,formatC(1, width=2,</pre>
   flag="0"), ".pdf", sep="")
 dev.print(pdf, file=fout)
  fout <- paste(file.prefix,formatC(l,width=2,</pre>
   flag="0"), ".eps", sep="")
  dev.print(postscript,file=fout)
if(MAKE.PLOTS){
  # plot 2D contour, legend, centered
 wiland <- 7; heland <- 6
 dev.new(width=wiland, height=heland)
 par(oma=rep(0,4), mar=c(4.2,3.2,4.2,0.5), mgp=
   c(1.5, 0.5, 0))
  if(BKGR.BLACK) par(fg="white", col="white",
   col.axis="white", col.lab="white", col.main="
   white", col.sub="white")
  tick <- diff(range(Xc)) * 0.015; tcol <- "</pre>
   white"
  plot.ticks <- function() {axis(1); axis(2);</pre>
   lines(c(Xcm,Xcm), c(min(Yc), min(Yc)+tick),
   col=tcol);lines(c(Xcm,Xcm), c(Ycm-tick, Ycm+
   tick), col=tcol);lines(c(Xcm,Xcm), c(max(Yc)-
   tick,max(Yc)), col=tcol);lines(c(min(Xc),min(
   Xc)+tick), c(Ycm,Ycm), col=tcol);lines(c(Xcm-
   tick,Xcm+tick), c(Ycm,Ycm), col=tcol);lines(c
   (max(Xc)-tick,max(Xc)), c(Ycm,Ycm), col=tcol)
```

.7)}

```
}
                                                         814
       plot.title <- function() {title(main=bquote(</pre>
784
        paste(I[fil], " = ", .(format(IH[1], nsmall
                                                        815
        =2)), " A ", "B = ", .(format(B[1]*10, nsmall
        =1)), " kG ", "V = ", .(format(V[1]/1e3,
                                                         816
        nsmall=2)), " kV ",I[e], " = ", .(format(I[1]
        *1e3, nsmall=0)), " mA")),xlab="X [mm]",ylab=817
        "Y [mm]")}
                                                         818
785
       filled.contour(Xc-Xcorr, Yc-Ycorr, Zphys,
                                                         819
        levels=seq(Zlim[1], Zlim[2], length.out=Nc),
                                                        820
        col=pplot.title=plot.tplot.axes=plot.tkey.
                                                         821
        title=title(main=expression(paste(j, " [", A/822
        cm^2, key.axes=axis(4))
                                                         823
       if(RUBBER.STAMP){
                                                         824
786
       mtext(paste(fns[1], "\n", "Analyzed on", date()825
787
         , " ",R.version.string),side = 1, line = 3,
        adj = 1, font = 3, cex = .7
                                                         826
       }
788
       # save to file
789
       file.prefix <- paste(out.dir, fname.id, "2</pre>
                                                         827
790
        Dcent_", sep="")
791
       fout <- paste(file.prefix,formatC(1, width=2,</pre>
        flag="0"), ".png", sep="")
       #dev.print(png, file=fout, width=wiland*reso,
792
        height=heland*reso)
       dev2bitmap(file=fout, width=wiland, height=
793
        heland, res=reso)
       fout <- paste(file.prefix,formatC(1, width=2, 828</pre>
794
        flag="0"), ".pdf", sep="")
       dev.print(pdf, file=fout)
795
       fout <- paste(file.prefix,formatC(1,width=2,</pre>
796
        flag="0"), ".eps", sep="")
       dev.print(postscript,file=fout)
797
                                                         829
     }
798
799
      # --- 3D PLOTS ---#
800
     if(MAKE.PLOTS){
801
                                                         830
       # plot 3D
802
                                                         831
       rgl.open()
803
       #par3d(windowRect=c(0, 0, pwpxRGL, phpxRGL),
804
                                                         832
        family="serif", cex=2.5)
       par3d(windowRect=c(0, 0, pwpxRGL, phpxRGL),
                                                         833
805
        family="serif")
       bg3d(pbgr)
806
                                                         834
       view3d(fov=45, userMatrix=matrix(c(0.7,-
807
        0.7,0,0,0.54,0.54,0.57,0,-0.4,-
                                                         835
        0.4,0.86,0,0,0,0,1), nrow=4, ncol=4, byrow=
        TRUE))
                                                         836
       if(D3.AXIS){persp3d(Xc, Yc, Zphys, col=mcol,
                                                         837
808
        lit=FALSE,xlab="X [mm]", ylab="Y [mm]", zlab=
        "j [A/cm<sup>2</sup>]",axes=TRUE,box=FALSE)}
                                                         838
       else{persp3d(Xc, Yc, Zphys, col=mcol, lit=
                                                         839
809
        FALSE, xlab="", ylab="", zlab="", axes=TRUE)}840
       grid3d(c("x+","y+"), at = NULL,, lwd = 1, lty 841
810
        = 1, n = 5)
                                                         842
        #bbox3d(color=c("#333377","black"), emission 843
811
         ="#333377", specular="#33337FF", shininess=5, 844
        alpha=0.8, xlab="X [mm]", ylab="Y[mm]", zlab="845
        j[A/cm^2]")
                                                         846
        #axes3d(edges="bbox", labels = TRUE, tick =
812
         TRUE, nticks = 5, main="Profile", xlab="X [mm
         ]", ylab="Y[mm]", zlab="j[A/cm^2]")
                                                         847
        #box3d()
813
```

```
#title3d(main = "Profile", sub = NULL, xlab =
   "X", ylab = "Y", zlab = "Z", line = NA)
 file.prefix <- paste(out.dir, fname.id, "3D_",</pre>
   sep="")
 fout <- paste(file.prefix,formatC(1, width=2,</pre>
  flag="0"), ".png", sep="")
 rgl.snapshot(fout)
}
if(MAKE.PLOTS){
  # 2D vacuum measurement
  if(VACUUM){
   #Plim <- c(1e-8, 1.5e-7)
   Plim <- range(Zvac, na.rm=TRUE)</pre>
   dev.new(width=wiland, height=heland)
   par(oma=c(0, 0, 0, 1), mar=c
     (4.2,3.2,4.2,0.5), mgp=c(1.5, 0.5, 0))
   if(BKGR.BLACK) par(fg="white", col="white",
    col.axis="white", col.lab="white", col.main=
     "white", col. sub="white")
   plot.title <- function() {title(main=bquote(</pre>
    paste(I[fil], " = ", .(format(IH[1], nsmall
    =2)), " A ", "B = ", .(format(B[1]*10,
    nsmall=1)), " kG ","V = ", .(format(V[1]/1
    e3, nsmall=2)), " kV ",I[e], " = ", .(
    format(I[1]*1e3, nsmall=0)), " mA")),xlab="
    H corrector setting [A]",ylab="V corrector
    setting [A]")}
   filled.contour(X, Y, Zvac, asp=1,levels=seq(
    Plim[1], Plim[2], length.out=Nc), col=
    palette,plot.title=plot.title(),key.title=
    title(main=expression(paste(P, " [mbar]")))
    )
   if(RUBBER.STAMP){mtext(paste(fns[1], "\n","
    Analyzed on", date(), " ",R.version.string)
     ,side = 1, line = 3, adj = 1,font = 3, cex
    = .7)
   # save to file
   file.prefix <- paste(out.dir, fname.id, "_</pre>
    vac2D_", sep="")
   fout <- paste(file.prefix,formatC(1, width</pre>
    =2, flag="0"), ".png", sep="")
   #dev.print(png, file=fout, width=wiland*reso
     , height=heland*reso)
   dev2bitmap(file=fout, width=wiland, height=
    heland, res=reso)
   fout <- paste(file.prefix,formatC(1, width</pre>
    =2, flag="0"), ".pdf", sep="")
   dev.print(pdf, file=fout)
   fout <- paste(file.prefix,formatC(1, width</pre>
    =2, flag="0"), ".eps", sep="")
   dev.print(postscript, file=fout)
  }
}
if(EXPORT.CALIBRATED.PROFILE){
 message("Exporting calibrated profile...")
 Xmat <- Xc %o% rep(1, Ny)</pre>
 Ymat <- rep(1, Nx) %o% Yc
 Zmat <- Zphys
 CalProf <- data.frame(x.mm = as.vector(Xmat),y</pre>
   .mm = as.vector(Ymat), j.Acm2 = as.vector(
  Zphys))
 fout <- paste(out.dir, fname.id, "_",formatC(1</pre>
   , width=2, flag="0"), "_calprof.txt", sep="")
```

```
write.table(CalProf, file=fout, row.names=
848
                                                            894
         FALSE, col.names=TRUE)
        system(paste("gzip", fout))
849
                                                            895
      7
850
                                                            896
      if(EXPORT.RADIAL){
851
                                                            897
        message("Calculating average radial profile...898
852
         ")
        Xmat <- (Xc-Xcorr) %o% rep(1, Ny)</pre>
853
                                                            899
        Ymat <- rep(1, Nx) %o% (Yc-Ycorr)
854
                                                            900
        r <- as.vector(sqrt(Xmat<sup>2</sup> + Ymat<sup>2</sup>))
855
                                                            901
        rstep <- max(Xcstep, Ycstep)/4</pre>
856
                                                            902
        rf <- cut(r, breaks=seq(min(r), min(r2[1]*</pre>
857
         1500, max(r)), by=rstep), ordered=TRUE)
                                                            903
        j <- as.vector(Zphys)</pre>
                                                            904
858
        rm <- tapply(r, rf, mean)</pre>
859
        jm <- tapply(j, rf, mean)</pre>
860
                                                            905
        rpro <- data.frame(r.mm=rm, j.au=jm)</pre>
                                                            906
861
        message("Plotting radial profile...")
862
        dev.new()
863
                                                            907
864
        par(oma=rep(0,4), mar=c(3.2, 3.2, 0.5, 0.5),
                                                            908
         mgp=c(1.5, 0.5, 0))
865
        plot(rm, jm, xlab="Radial distance [mm]", ylab909
         ="Current density [arb. units]",pch="+",col="
         blue")
                                                            910
        file.prefix <- paste(radial.dir, fname.id, "_"911</pre>
866
         , formatC(1, width=2, flag="0"), "_radial",
         sep="")
                                                            912
        fout <- paste(file.prefix, ".pdf", sep="")</pre>
867
        dev.print(pdf, file=fout)
868
        message("Exporting radial profile to text file913
869
         ...")
        fout <- paste(file.prefix, ".txt", sep="")</pre>
870
                                                            914
        write.table(rpro, file=fout, row.names=FALSE,
871
                                                            915
         col.names=TRUE)
                                                            916
      7
872
                                                            917
873
                                                            918
      # downsample a large array for plotting
874
                                                            919
       purposes
                                                            920
      # function returns ordered vector of indices
                                                            921
875
      idwnsmpl <- function(x=runif(100), n=10){i <-</pre>
                                                            922
876
       sort(sample(x=1:length(x), size=n));return(i)} 923
877
      # export particle coordinates
878
                                                            924
      if(EXPORT.PARTICLES){
879
                                                            925
        message("Exporting particle coordinates for
880
                                                            926
         Warp field solver...")
                                                            927
        Znorm < - Z/sum(Z)
881
                                                            928
        Np <- round(Nmax*Znorm)
882
                                                            929
        cat("Max # of particles per bin:", max(Np), "\930
883
         n")
                                                            931
        cat("Empty bins:", length(Np[Np==0]), "\n")
884
                                                            932
        xpart <- numeric(Nmax); xpart <- NA</pre>
885
                                                            933
        ypart <- numeric(Nmax); ypart <- NA</pre>
886
        k <- 1
887
                                                            934
        for(i in 1:Nx){
888
                                                            935
         for(j in 1:Ny){
889
          if(Np[i,j]>=1 & k <= Nmax){</pre>
890
                                                            936
           for(m in 1:Np[i,j]){
891
            # generate coordinates uniformly within
892
                                                            937
             each bin
            xpart[k] <- Xc[i] + runif(1, -0.5*Xcstep, 938</pre>
893
             0.5*Xcstep)
```

```
ypart[k] <- Yc[j] + runif(1, -0.5*Ycstep,</pre>
      0.5*Ycstep)
     k <- k+1
   }}
 }}
  cut <- (is.na(xpart)==FALSE) & (is.na(ypart)==</pre>
  FALSE)
 xpart <- xpart[cut]; ypart <- ypart[cut]</pre>
 part_coord <- data.frame(xpart, ypart)</pre>
  dev.new()
 par(oma=rep(0,4), mar=c(3.2, 3.2, 0.5, 0.5),
  mgp=c(1.5, 0.5, 0))
 N.plot.max <- 4096
  if(length(xpart) > N.plot.max) ii <- idwnsmpl(</pre>
   x=xpart, n=N.plot.max)
  else ii <- 1:length(xpart)</pre>
 plot(xpart[ii], ypart[ii], pch=".", asp=1,xlab
   ="X coord. [mm]", ylab="Y coord. [mm]")
 message("Saving Particle Plot")
  file.prefix <- paste(out.dir, fname.id, "_",</pre>
   sep="")
  fout <- paste(file.prefix,formatC(1, width=2,</pre>
   flag="0"), "_particles.eps", sep="")
  dev.print(postscript, file=fout)
 message("Writing coordinates of ", length(
  xpart), " particles...")
  fnpart <- paste(file.prefix,formatC(1, width</pre>
   =2, flag="0"), "_", length(xpart), "_
   particles.txt", sep="")
  write.table(part_coord, file=fnpart, row.names
   =FALSE, col.names=FALSE)
  system(paste("gzip", fnpart))
}
```

```
•
```

```
# polar decomposition
if(POLAR.DECOMPOSITION){
 dev.new(width=9, height=7)
  source("../Polar/polar_decomposition.R")
 nn <- 1:32
                 #Was 128
 mm < - 0:32
                #Was 48
 PD <- pol_dec(Xc=Xc-Xcorr,Yc=Yc-Ycorr,Z=Z,nn=
  nn,mm=mm,a=30,BOUNDARY=0)
 P_mat <- PD$P
  #rownames(P_mat)<-1:32</pre>
  #colnames(P_mat)<-0:32</pre>
 nModes <- rowSums(Mod(P_mat))</pre>
 mModes <- colSums(Mod(P_mat))</pre>
 I mat <- PD$I
```

plots

```
par(oma=c(0,0,3,0))
a <- layout(matrix(c(1,4,2,5,3,6),2,3),widths=
c(0.9,0.9,0.9))
layout.show(a)
image(Xc-Xcorr,Yc-Ycorr,Z,asp=1,col=palette,
main="Measurement")
image(Xc-Xcorr,Yc-Ycorr,Mod(I_mat),asp=1,col=
palette,main="Reconstruction")
image(Xc-Xcorr,Yc-Ycorr,abs(Mod(I_mat)-Z),asp
=1,col=gray(0:64/64), main="Errors")
ticks <- c(0.00001,
0.0001,0.0001,0.001,0.01,0.10)
```

939	<pre>image.plot(nn,mm,log10(Mod(P_mat)), col=heat. 949</pre>	<pre>save(P_mat, file=fncoe)</pre>
	<pre>colors(64),main="Mode amplitudes", xlab=" 950</pre>	# save log
	Radial mode n",ylab="Azimuthal mode m",nlevel951	<pre>fnlog <- paste(file.prefix,formatC(l,width=2,</pre>
	=64, axis.args=list(at=log10(ticks), labels=	flag="0"), ".txt", sep="")
	ticks)) 952	<pre>cmnd <- paste("mv pol_dec_log.txt", fnlog)</pre>
940	<pre>plot(nn,nModes,type="o",log="y",pch=16, main="953</pre>	<pre>system(cmnd)</pre>
	Radial modes",xlab="Radial mode n", ylab=" 954	# save figures
	Amplitude") 955	<pre>fnfig <- paste(file.prefix,formatC(l,width=2,</pre>
941	<pre>plot(mm,mModes,type="o",log="y",pch=16, main="</pre>	flag="0"),"_polar", ".pdf", sep="")
	Azimuthal modes", xlab="Azimuthal mode m", 956	<pre>cmnd <- paste("cp polar.pdf", fnfig)</pre>
	ylab="Amplitude") 957	system(cmnd)
942	<pre>mtext(fname.id, outer=TRUE) 958</pre>	}
943	<pre>dev.print(pdf,file="polar.pdf",width=9,height 959</pre>	}
	=7,pointsize=10,family="Times") 960	<pre>if(MAKE.PLOTS) dev.off()</pre>
944	layout(1) 961	<pre>quartz.options(reset=TRUE)</pre>
945	<pre>file.prefix <- paste(polar.dir, fname.id, "_",962</pre>	pdf.options(reset=TRUE)
	sep="") 963	ps.options(reset=TRUE)
946	# save coefficients 964	
947	<pre>fncoe <- paste(file.prefix, 965</pre>	$cat("\nStop: ", date(),"\n")$
948	<pre>formatC(1,width=2, flag="0"), ".RData", sep="966</pre>	sink()
	")	

C.2 Transverse Field Measurement

```
95
2 # NAME: Fieldsolver_single.py
3 # AUTHOR: Vince Moens, Giulio Stancari
4 # LAST MODIFIED: 29.07.2013
                                                   96
5 # PURPOSE: Calculation of charge density,
                                                   97
   electric potential and electric fields of the
   measured electron beam profiles
8 from warp import *
9 from numpy.linalg import *
10 import matplotlib.mlab as mlab
11 import matplotlib.pyplot as plt
12 import math as math
13
14 # Options
15 printfields = true
16 particleplot = true
17 chargedensityplot = true
18 potentialplot = true
19 electricfieldplot = true
20 electricfieldlineplot = true
21 zoomedplots = true
22 makepsfile = false
                                                   101
23 findmaxcenter = true
24 security_check = true
25 LHC = true
26
27 # Loading Files
28 \text{ fns} = \begin{bmatrix} 1 \\ 1 \end{bmatrix}
29 # new profiles acquired with ACL script
   "../Profile/Results/Chart_Colors/
30
     HG1b_121218_9p25A_3-3-3
     kG_500V_51mA_b_57_8102_particles.txt", #0
```

. . .

```
95 "../Profile/Results/Chart_Colors/
HG1b_130521_9p25A_06-24-06
kG_4kV_1368mA_136_8192_particles.txt" #63
96 ]
```

```
98 Bgun=[3, 3, 3, 3, 3, 3, 3, 3, 3, 3, 3, 3, 1, 1, 1, 2.5,
2, 2.5, 2.5, 2.5, 2.5, 2, 2, 2, 2, 1.5, 1.5, 1.5,
3, 2.5, 3, 3, 3, 3, 3, 1, 1, 0.8, 0.8, 0.8, 0.8,
0.8, 0.8, 0.8, 0.8, 0.8, 0.8, 0.8, 1, 1, 1,
0.6, 0.4, 0.4, 1, 1, 1, 0.6, 1, 0.4, 1, 0.4,
0.4, 0.4, 0.6, 0.6, 0.6]
```

```
103 Current=[0.051, 2.940, 1.100, 0.388, 0.140,
1.924, 0.694, 2.394, 0.440, 0.027, 0.118, 0.398,
0.726, 1.770, 0.928, 1.484, 0.412, 0.865,
1.460, 0.213, 0.448, 0.744, 0.090, 0.390, 0.192,
0.582, 0.051, 0.050, 0.050, 0.050, 0.050,
```

```
0.050, 0.070, 0.070, 0.073, 0.526, 1.450, 0.195, 157 # --- beam
     3.7, 3.1, 0.026, 0.009, 0.944, 1.99, 2.53, 158 top.ekin = Voltage[n]
    1.022, 0.196, 2.754, 0.073, 0.195, 0.072, 0.073, 159 if security_check:
     0.534, 1.44, 0.196, 1.98, 0.488, 3.88, 0.125, 160
                                                           if not ((str(int(top.ekin))[0]+"kV" in item) or
    0.332, 0.658, 0.526, 0.924, 1.368]
                                                               (str(int(top.ekin)) in item)):
104 n=63
                                                              print("Voltage is wrong!")
                                                       161
105 if(n in [24]):LHC = false
                                                       162
                                                              sys.exit()
106 item=fns[n]
                                                       163 top.vbeam = 0
107 print("Plotting results...")
                                                       164 top.ibeam = -Current[n]
108 palette("ImageJ_Fire.gp")
                                                       165 if security_check:
109 Bgun=Bgun[n] *0.1
                                                       166 if str(int(top.ibeam*-1e3)) not in item:
110 Bmain=Bmain[n]*0.1
                                                              print("Current is wrong!")
                                                       167
111 Bcoll=Bcoll[n]*0.1
                                                       168
                                                              sys.exit()
112 print("Bgun = "+str(Bgun)+" T; Bmain = "+str(
                                                       169 top.lrelativ = true
    Bmain)+" T; Bcoll = "+str(Bcoll)+"_T")
                                                       170 top.relativity = true
                                                       171 top.derivqty()
113
114 # --- Set comment lines, user's name.
                                                       172 totalweight = top.ibeam * dz / abs(elec.zion) /
115 if item.startswith("../Profile/Results/
                                                           top.echarge / top.vbeam
    Chart_Colors/"): top.runid="Results/H1"+ item
                                                       173
    [38:]
                                                       174 # --- initialize
116 top.pline2 = "Hollow electron beam -1-in cathode"
                                                       175 print("Initializing...")
117 top.pline1 = "Field calculation from measured
                                                       176 package("wxy")
    current density profile"
                                                       177 generate()
118 top.runmaker = "V. Moens"
                                                       178 installconductor(driftpipe)
119 setup(makepsfile)
                                                       179
120 winon()
                                                       180 # --- read particle positions
                                                       181 print("Reading particle positions...")
121
122 # --- Define the electron speices
                                                       182 posi = fromfile(item, sep=' ')
123 elec = Species(type=Electron)
                                                       183 npart = len(posi)/2
                                                       184 posi = reshape(posi, (npart,2))
124
125 # --- lattice
                                                       185
126 print("Defining lattice...")
                                                       186 # --- Set particle weight for the beam
127 rpipe = 0.03 # test stand: 3 cm; TEL2: 3.5 cm
                                                       187 elec.sw = totalweight/npart
128 syslen = 2.86
                                                       188 elec.nps = 0
129 top.drftzs[0] = 0
                                                       189
130 top.drftze[0] = syslen
                                                       190 # --- add particles according to measured current
131 top.drftap[0] = rpipe
                                                           density
132 print("Beam pipe radius = "+str(rpipe)+" m")
                                                       191 print("Calculating charge density according to
                                                           particle distribution...")
133
134 # --- field solver
                                                       192 print("Macroparticle weight = %f" % elec.sw)
135 print("Setting up field solver...")
                                                       193 xinit = posi[:,0]*math.sqrt(float(Bcoll)/float(
136 w3d.solvergeom = w3d.XYgeom
                                                           Bmain)) * mm
137 \text{ fudge} = 1.05
                                                       194 yinit = posi[:,1]*math.sqrt(float(Bcoll)/float(
                                                           Bmain)) * mm
138 w3d.xmmax = rpipe*fudge
139 w3d.xmmin = -rpipe*fudge
                                                       195 zinit = zeros(npart)
                                                       196 vxinit = zeros(npart)
140 w3d.ymmax = rpipe*fudge
141 w3d.ymmin = -rpipe*fudge
                                                       197 vyinit = zeros(npart)
142 w3d.zmmax = syslen
                                                       198 vzinit = zeros(npart) + top.vbeam
143 w3d.zmmin = 0
                                                       199 elec.addpart(xinit,yinit,zinit,vxinit,vyinit,
144 \text{ w3d.nx} = 256
                                                           vzinit)
145 \text{ w3d.ny} = 256
                                                       200
146 w3d.nz = 1
                                                       201 # --- deposit particle charge density
147 dx = (w3d.xmmax -w3d.xmmin) / w3d.nx
                                                       202 loadrho()
148 dy = (w3d.ymmax -w3d.ymmin) / w3d.ny
                                                       203
149 dz = (w3d.zmmax -w3d.zmmin) / w3d.nz
                                                       204 # --- find corresponding potentials and fields
150 w3d.l2symtry = false
                                                       205 print("Solving for fields...")
151 w3d.l4symtry = false
                                                       206 fieldsolxy(-1)
152
                                                       207
153 # --- Conductors
                                                       208 # X vs. Y
154 driftpipe = ZCylinderOut(condid = 1, radius=0.03, 209 if particleplot:
     zlower=0, zupper=syslen, voltage=0)
                                                       210 ppxy()
155 scraper=ParticleScraper(driftpipe)
                                                       211
                                                            fma()
                                                       212 # charge density in x-y plane
156
```

```
Appendix C. Codes
```

```
213 if chargedensityplot:
                                                        266
214
     pcrhoxy(lframe=1, cellarray=1, contours=None)
                                                        267
215
     fma()
                                                        268
216 # potential
                                                        269
217 if potentialplot:
     pcphixy(lframe=1, cellarray=1)
218
                                                        271
219
     fma()
                                                        272
220 # electric field components and strength
                                                        273
221 if electricfieldplot:
                                                        274
     pcselfexy(lframe=1, comp='x', cellarray=1)
222
223
     fma()
                                                        275
     pcselfexy(lframe=1, comp='y', cellarray=1)
224
                                                        276
     fma()
225
                                                        277
226
     Ef = sqrt(w3d.selfe[0,:,:,0]**2 + w3d.selfe
227
      [1,:,:,0]**2 + w3d.selfe[2,:,:,0]**2)
     ppgeneric(grid=Ef,x=w3d.xmesh,y=w3d.ymesh,
228
      contours=10,cellarray=1,titles=0)
     ptitles("Electric field strength in x-y plane",
229
      "X","Y")
     fma()
230
231
232 if electricfieldlineplot:
     plg(Ef[w3d.nx/2,:]/1e3, w3d.ymesh*1e3, color="
233
      red")
     plg(Ef[:,w3d.ny/2]/1e3, w3d.xmesh*1e3, color="
234
      blue")
     ptitles("E-field vs X (blue) and Y (red)","X or 288 Gauss=Gauss/sum(Gauss)
235
       Y (mm)","E-field (kV/m)")
     fma()
236
237
238 # zoomed-in plots
239 if zoomedplots:
     if particleplot:
240
       pcrhoxy(lframe=1, cellarray=1, contours=None) 292 Efcw_RMS=std(Efcw)
241
       limits(-0.01, 0.01, -0.01, 0.01)
242
       fma()
243
     if potentialplot:
244
       pcphixy(lframe=1, cellarray=1, contours=None) 296
245
       limits(-0.01, 0.01, -0.01, 0.01)
246
       fma()
247
     if electricfieldplot:
248
       ppgeneric(grid=Ef,x=w3d.xmesh,y=w3d.ymesh,
249
        contours=None,cellarray=1,titles=0)
       limits(-0.01, 0.01, -0.01, 0.01)
250
       ptitles("Electric field strength in x-y plane"303 q=1
251
        ,"X","Y")
       fma()
252
253
254 # --- write results
255 if printfields:
     print("Writing results to file...")
256
     roi = 0.01 # region of interest in m, +/-
257
     ispan = int(round(roi/dx))
258
     print("Half span in grid points = "+repr(ispan)
259
      )
     i1 = w3d.nx/2 - ispan
260
     i2 = w3d.nx/2 + ispan
261
     name = "Results/H1"+item[38:-4]+"_fields.txt"
262
     out = open(name, 'wb')
263
264
     line = "Xmesh_mm\tYmesh_mm\tRho_Cm3\tPhi_V\
      tEx_Vm\tEy_Vm\n"
     out.write(line)
265
```

```
for i in range(i1, i2):
         for j in range(i1, i2):
             xx = repr(w3d.xmesh[i]*1e3)
                 = repr(w3d.ymesh[j]*1e3)
             VV
             rho = repr(w3d.rho[i,j,0])
             pot = repr(w3d.phi[i,j,0])
             ex = repr(w3d.selfe[0,i,j,0])
             ey = repr(w3d.selfe[1,i,j,0])
             line = xx+"\t"+yy+"\t"+rho+"\t"+pot+"\t"
              +ex+"\t"+ey+"\n"
             out.write(line)
     out.close()
278 ## -- Calculating the emittance Groth in the
     center --##
279 print "\n Calculating Emittance Growth in the
    center n"
_{280} intervall = 0.8
281 mean=0
282 sigma=4.69e-4
283 print "Sigma = "+ str(sigma)
284 Cath_radius=6.75e-3
285 print "Cathode Radius = "+ str(Cath_radius)
286 R = array([[sqrt(w3d.ymesh[i]**2+w3d.xmesh[j]**2)
     for i in range(w3d.ymesh.size)] for j in range(
    w3d.xmesh.size)])
287 Gauss=mlab.normpdf(R,mean,sigma)
289 Efw=array([[Ef[i,j]*Gauss[i,j] for i in range(w3d
     .ymesh.size)] for j in range(w3d.xmesh.size)])
290 Efcw=Efw[R <= 0.00675*intervall*math.sqrt(float(
    Bcoll)/float(Bmain))]
291 Efc=Ef[R <= 0.00675*intervall*math.sqrt(float(</pre>
    Bcoll)/float(Bmain))]
293 Efc_RMS=std(Efc)
294 Efcw_mean=average(Efcw)
295 Efc_mean=average(Efc)
297 #Calculating emittance growth at center
298 m0=938.272e6
299 print "Proton Rest mass = "+str(m0)+" eV"
300 if (LHC): Etot=4e12
301 else: Etot=980e9
302 print "Proton Energy = "+str(Etot)+" eV"
304 print "Elementary Charge is = "+str(q)+" electron
     charge"
305 Er_center=Efcw_RMS
306 Er_center_4=Efcw_RMS*(6.75e-3*math.sqrt(float(
    Bcoll)/float(Bmain))/(4*sigma))
307 Er_center_6=Efcw_RMS*(6.75e-3*math.sqrt(float(
    Bcoll)/float(Bmain)) /(6*sigma))
308 print "Maximum electric field in center is "+str(
    Er_center)+" V/m"
309 L=2.65
310 print "Length of interaction region is "+str(L)+"
     m"
311 gamma=Etot/m0
312 print "gamma is "+str(Etot/m0)
313 if(LHC): beta=250
```

```
314 else: beta=150
```

```
315 print "beta is "+str(beta)
```

316 if (LHC): R_freq=1.12e4 339 print "emittance growth rate is "+str(gr_center_6)+" 1/s at 6 sigma" 317 else: R_freq=4.7e4 340 print "\n" 318 t_rev=1/R_freq 319 print "revolution time is "+str(t_rev) 341 print str(item[38:-4]) 320 if (LHC): $_{342}$ print "\n Calculating Emittance Growth overall \n Emit=3.75e-6 321 print "Emittrance of LHC is "+str(Emit) 343 322 344 name = "Results/H1"+item[38:-4]+"_Emax_Gaussian. 323 else: 324 Emit=2.5e-6 txt" print "Emittrance of TEV is "+str(Emit) 345 print "Saving data to text file: "+str(name)+"\n" 325 326 kick_center=math.atan(Etot*q*Er_center*L/(Etot**2 346 out = open(name, 'wb') 347 titleline = item[38:-4]+"\tEmax [kV/m]\tKick []\ -m0**2)) 327 kick_center_4=math.atan(Etot*q*Er_center_4*L/(tEmittance Growth [m/turn] \tEmittance Growth Etot**2-m0**2)) Rate [1/s]\n" 348 out.write(titleline) 328 kick_center_6=math.atan(Etot*q*Er_center_6*L/(Etot**2-m0**2)) 349 line = "Center:\t"+str(Er_center)+"\t"+str(329 print "The kick is "+str(kick_center) kick_center)+"\t"+str(delepsilon_center)+"\t"+ 330 delepsilon_center=beta*kick_center**2*gamma $str(gr_center) + "\n"$ 331 delepsilon_center_4=beta*kick_center_4**2*gamma 350 out.write(line) 332 delepsilon_center_6=beta*kick_center_6**2*gamma 351 line = "Center_6s:\t"+str(Er_center_6)+"\t"+str(333 gr_center=delepsilon_center/(t_rev*Emit) kick_center_6)+"\t"+str(delepsilon_center_6)+"\t 334 gr_center_4=delepsilon_center_4/(t_rev*Emit) "+str(gr_center_6)+"n" 352 out.write(line) 335 gr_center_6=delepsilon_center_6/(t_rev*Emit) 336 print "Emittance growth is "+str(353 line = "Center_4s:\t"+str(Er_center_4)+"\t"+str(delepsilon_center)+" m/turn" kick_center_4)+"\t"+str(delepsilon_center_4)+"\t 337 print "emittance growth rate is "+str(gr_center)+ "+str(gr_center_4)+"n" " 1/s at cathode radius" 354 out.write(line) 338 print "emittance growth rate is "+str(gr_center_4 355 out.close())+" 1/s at 4 sigma" Attachments/Fieldsolver single app.py

C.3 Numerical WARP simulations

```
22
2 # NAME: tbench_complex.py
3 # AUTHOR: Vince Moens, Giulio Stancari
4 # LAST MODIFIED: 05.08.2013
5 # PURPOSE: 3D Simulations in WARP using gun or
                                        26
  profile injection
8 # NOTES ON EMITTANCE TYPES
9 # GUN: The code automatically injects particles
  from the cathode conductor using Child Langmuir
                                        91
  law.
10 # PROFILE: Profiles measured in the test bench
  are injected into the lattice. The gun is
                                        92
  ommitted fromt the lattice. Injection takes
                                        93
  place at the end of that anode.
11
13 # >>> Package Loading <<< #
15
16 from warp import * # warp code
17 from datetime import *
18
20 # >>> File Loading <<< #
```

```
23 # --- Profiles ---#
```

```
24 fns = [
```

```
25 # new profiles acquired with ACL script
```

% "../../HG1b/Profile/Results/Chart_Colors/ HG1b_121218_9p25A_3-3-3 kG_500V_51mA_b_57_8102_particles.txt", #0

. . .

```
91 "../../HG1b/Profile/Results/Chart_Colors/
HG1b_130521_9p25A_06-24-06
kG_4kV_1368mA_136_8192_particles.txt" #63
```

94 Bgun=[3, 3, 3, 3, 3, 3, 3, 3, 3, 3, 3, 3, 1, 1, 1, 2.5, 2, 2.5, 2.5, 2.5, 2.5, 2, 2, 2, 2, 1.5, 1.5, 1.5, 3, 2.5, 3, 3, 3, 3, 3, 1, 1, 0.8, 0.8, 0.8, 0.8, 0.8, 0.8, 0.8, 0.8, 0.8, 0.8, 0.8, 1, 1, 1, 0.6, 0.4, 0.4, 1, 1, 1, 0.6, 1, 0.4, 1, 0.4, 0.4, 0.4, 0.6, 0.6, 0.6]

```
96 Bcoll=[3, 3, 3, 3, 3, 3, 3, 3, 3, 3, 3, 1, 1, 1,
                                                   134 os.system("cp tbench_complex.py* ../Results/"+
    2.5, 2, 2.5, 2.5, 2.5, 2.5, 2, 2, 2, 1.5, 1.5,
                                                       date+"/tbench_"+date+time+"_test.py")
    1.5, 3, 2.5, 3, 3, 3, 3, 3, 3, 3, 2, 0.8, 0.8, 0.8, 135 top.runid = machine_type+"_"+date+time+"_"+
     machine_injtype+"_"+file_ending
    1, 0.6, 0.4, 0.4, 1, 1, 1, 0.6, 1, 0.4, 1, 0.4, 136 if machine_type == "tbench": top.pline2 = "
    0.4, 0.4, 0.6, 0.6, 0.6]
                                                        Electron Lens Test Bench"
                                                    137 else: top.pline2 = "Tevatron Electron Lens 2"
97
98 Voltage=[500, 8000, 4000, 2000, 1000, 6000, 3000, 138 if machine_emittype == 1: top.pline1 = "Constant-
     7000, 8000, 8000, 889, 2000, 3000, 5556, 3556,
                                                        injection_" + machine_injtype
    5000, 2083, 3472, 5000, 1333, 2222, 3111, 750,
                                                    139 elif machine_emittype == 2: top.pline1 = "Child-
    2000, 1250, 3000, 500, 500, 500, 500, 500, 500,
                                                       Langmuir_" + machine_injtype
    500, 500, 500, 2000, 4000, 1000, 8000, 7000,
                                                    140 else: top.pline1 = "other injection method"
    250, 125, 3000, 5000, 6000, 3125, 1000, 6250,
                                                    141 top.runmaker = "V. Moens"
    500, 1000, 500, 500, 2000, 4000, 1000, 5000,
                                                    142
    2000, 8000, 750, 1500, 2500, 2000, 3000, 4000]
                                                   99 Current=[0.051, 2.940, 1.100, 0.388, 0.140,
                                                   144 # >>> Variables <<< #
    0.726, 1.770, 0.928, 1.484, 0.412, 0.865,
                                                   146 # --- Machine Parameters ---#
    1.460, 0.213, 0.448, 0.744, 0.090, 0.390, 0.192, 147 machine_zstart = .0e0
     0.582, 0.051, 0.050, 0.050, 0.050, 0.050,
                                                   148 machine_syslen = 2.86
    0.050, 0.070, 0.070, 0.073, 0.526, 1.450, 0.195, 149 machine_zplat = machine_syslen
     3.7, 3.1, 0.026, 0.009, 0.944, 1.99, 2.53,
                                                   150 machine_piperad = 3*cm
    1.022, 0.196, 2.754, 0.073, 0.195, 0.072, 0.073, 151 zfinal = machine_zstart + machine_syslen
     0.534, 1.44, 0.196, 1.98, 0.488, 3.88, 0.125,
                                                   152
    0.332, 0.658, 0.526, 0.924, 1.368]
                                                   153 # --- Electron Gun ---#
                                                   154 # - Cathode
100
101 # --- Selecting profile ---#
                                                   155 Cathode zstart = -29.25*mm
102 \text{ item} = 51
                                                   156 Cathode_zend = 0.0*mm
103 file_ending = "test"
                                                   157 Cathode_radi = 6.75*mm
104 Bmain = float(Bmain[item])/10
                                                   158 Cathode_rado = 12.7*mm
105 print("Magetic Field in Main Solenoid: %g " %
                                                   159 Cathode_radcurvb = 10*mm
    Bmain)
                                                    160 Cahtode_radcurvs = 0.5*mm
106 Bgun = float(Bgun[item])/10
                                                   161 Cathode_voltage = Cathode_Potential
107 Bcoll = float(Bcoll[item])/10
                                                   162 # - Anode
108 Cathode_Potential = -Voltage[item]
                                                   163 Anode_zstart = 9.48*mm
109 print("Cathode Potential: %g " %
                                                   164 Anode_z1 = Anode_zstart + 1.5*mm
    Cathode_Potential)
                                                   165 Anode_z2 = Anode_zstart + 3.5*mm
110 Current = Current[item]
                                                   166 Anode_z3 = Anode_z1 + 9*mm
111 print("Current: %g " % Current)
                                                   167 \text{ Anode}_{z4} = \text{Anode}_{z3} + 11.25 \text{*mm}
112 npart = -50*Cathode_Potential/500
                                                   168 Anode_z5 = Anode_z4 + 5.625*mm
                                                   169 Anode_zend = Anode_z5 + 58.5*mm
113
170 Anode_ri = 14.25*mm
115 # >>> Options <<< #
                                                    171 Anode_ro = Anode_ri+5.33*mm
172 Anode_r1 = Anode_ri
117 machine_injtype = "profile"
                                                   173 Anode_r2 = Anode_ro
118 print("Injection type is:"+machine_injtype)
                                                   174 Anode_r3 = Anode_ri
119 if (machine_injtype=="gun"): machine_emittype = 2 175 Anode_r4 = Anode_ri + 0.675*mm
120 elif (machine_injtype=="profile"):
                                                   176 Anode_radtipi = Anode_ri + 1.5*mm
    machine_emittype = 1
                                                   177 Anode_radtipo = Anode_ro -3.5*mm
121 else:
                                                   178 Anode_r5 = Anode_ri + 5.625*mm
122 print("Wrong machine_injtype!!")
                                                   179 Anode_rendi = Anode_r5
                                                   180 Anode_rendo = Anode_rendi + 1.35*mm
123
    quit()
124 machine_type = "tbench"
                                                   181 Anode_radcurvb = 3.5*mm
                                                   182 Anode_radcurvs = -1.5*mm
125
183 Anode_voltage = 0.0e0
127 # >>> Headers <<< #
                                                   184 # - Electrode F
185 ElectrodeF_zstart = Cathode_zstart
129 now=datetime.now()
                                                   186 ElectrodeF_zend = 0.98*mm
130 date=now.strftime("%y%m%d")
                                                    187 ElectrodeF_z1 = ElectrodeF_zend -0.5*mm
131 time=now.strftime("%H%M")
                                                    188 ElectrodeF_z2 = ElectrodeF_zend -1.4*mm
132 if not os.path.exists("../Results/"+date+"/"):
                                                    189 ElectrodeF_ri = 13.1*mm
      os.makedirs("../Results/"+date+"/")
                                                    190 ElectrodeF_ro = ElectrodeF_ri + 1.9*mm
133
                                                    191 ElectrodeF_r1 = ElectrodeF_ri + 0.5*mm
```

```
192 ElectrodeF_radcurvs = -0.5*mm
193 ElectrodeF_radcurvb = 1.4*mm
194 ElectrodeF_voltage = Cathode_Potential
195 # - Electrode C
196 ElectrodeC_zstart = Cathode_zstart
197 ElectrodeC_zend = 1.97*mm
198 ElectrodeC_ri = 20.5*mm
199 ElectrodeC_ro = 22.0*mm
200 ElectrodeC_radcurv = 0.75*mm
201 ElectrodeC_z1 = ElectrodeC_zend-0.75*mm
202 ElectrodeC_voltage = Cathode_Potential
203 # - Gun drift pipe
204 Gun_pipe_zstart = 84.375*mm
205 Gun_pipe_zend = 178.875*mm
206 Gun_pipe_ri = 36*mm #Should this not be 3 cm?
207 Gun_pipe_ro = 33.75*mm
208 Gun_pipe_voltage = 0.0
209
210 # ---- Solenoids ---#
211 # - Gun Solenoid
212 tbench_solenoid_gun_zstart = -13*cm
213 tbench_solenoid_gun_zend = 37*cm
214 tbench_solenoid_gun_radi = 28*cm
215 tbench_solenoid_gun_rado =
    tbench_solenoid_gun_radi+5.433*cm
216 tbench_solenoid_gun_b = Bgun
217 tbench_solenoid_gun_voltage = 0.0
218 # - Main Solenoid
219 tbench_solenoid_main_zstart = 0.60
220 tbench_solenoid_main_zend = 2.52
221 tbench_solenoid_main_radi = 0.20
222 tbench_solenoid_main_rado =
    tbench_solenoid_main_radi+14.48*cm
223 tbench_solenoid_main_b = Bmain
224 tbench_solenoid_main_voltage = 0.0
225 # - Collector Solenoid
226 tbench_solenoid_col_zstart = 2.67
227 tbench_solenoid_col_zend = 3.17
228 tbench_solenoid_col_radi = 28*cm
229 tbench_solenoid_col_rado =
    tbench_solenoid_col_radi + 5.433*cm
230 tbench_solenoid_col_b = Bcoll
231 tbench_solenoid_col_voltage = 0.0
232
233 # --- Drift Spaces ---#
234 # - First Drift
235 tbench_drift1_zstart = 37*cm
236 tbench_drift1_zend = 0.60
237 tbench_drift1_ap = machine_piperad
238 # - Second Drift
239 tbench_drift2_zstart = 2.52
240 tbench_drift2_zend = 2.67
241 tbench_drift2_ap = machine_piperad
242
243 # ---Beam size & position ---#
244 beama0 = 17.5e0*mm
_{245} beamb0 = 17.5e0*mm
_{246} beamap0 = .0e0*mm
_{247} beambp0 = .0e0*mm
248 beamx0 = .0e0*mm
_{249} beamy0 = .0e0*mm
250 \text{ beamxpO} = .0e0*mm
251 beamyp0 = .0e0*mm
```

```
252
253 # --- Beam inject parameters ---#
254 beamxinject = .0e0*mm
255 beamyinject = .0e0*mm
256 beamxpinject = .0e0*mm
257 beamypinject = .0e0*mm
258 beamainject = 17.5*mm
259 beambinject = 17.5*mm
260 beamapinject = .0e0*mm
261 beambpinject = .0e0*mm
262 beamainjmin = 6.75*mm
263 beambinjmin = 6.75*mm
264 beamzinject = machine_zstart
265
267 # >>> Script <<< #
269
270 #-----#
271 # Invoke setup routine #
272 #-----#
273 setup(makepsfile=0)
274 # winon()
275 palette("ImageJ_Fire.gp")
276
277 #-----#
278 # Particle Loading #
279 #-----#
280 # We only need to load the particles if the
    inject type is profle
281 if (machine_injtype == "profile"):
282 print("Reading particle positions...")
283 posi = fromfile(fns[item], sep=' ')
_{284} npart = len(posi)/2
285 posi = reshape(posi, (npart,2))
286 print("Calculating charge density according to
     particle distribution...")
287 print("Number of particles = %e" % npart)
288
289 #-----#
290 # Particle Properties #
291 #-----#
292 # --- Particle parameters ---#
293 electron_Iz= -Current294 cyc_freq=echarge*Bmain/emass
                     = pi/(2*cyc_freq)
295 timestep
                    = .0e0
= -Cathode_Potential
296 electron_vz
297 electron_ekin
298 electron_q
                     = -1.e0
                      = .0e0
299 vthz
300 lrelativity
                      = true
301 relativity
                       = true
302 sw=int((-electron_Iz*timestep/echarge)/npart)
303 elec = Species(type=Electron,color=red,weight=sw)
304 elec.ibeam = electron_Iz
305 print ("Beam Current: %g" % elec.ibeam)
306 elec.zion = electron_q
307 print ("Particle charge: %g" % elec.zion)
308 top.dt
           = timestep
309 print ("Cyclotron Frequency: %g" % cyc_freq)
310 print ("Timestep: %g" % top.dt)
311 elec.vbeam
              = electron_vz
```

312 print ("Particle velocity: %g" % elec.vbeam)

```
313 elec.ekin
                   = electron_ekin
                                                              w3d.uxt[:] = vthz
                                                     374
                                                              w3d.uyt[:] = vthz
314 elec.aion
                   = top.emass/top.amu
                                                     375
                                                              w3d.uzt[:] = elec.vbeam
315 top.derivqty()
                                                     376
                  = lrelativity
                                                           installuserparticlesinjection(
316 elec.lrelativ
                                                     377
317 elec.relativity = relativity
                                                           hollow_cathode_source)
318 elec.vthz
                  = vthz
                                                     378
                                                     379 #-----#
319 nsteps = 1.3*machine_syslen/elec.vbeam/timestep
320 print("The number of time steps is: %f" % nsteps) 380 # Lattice #
321
                                                     381 #-----#
322 #-----#
                                                      382 # - Gun Solenoid
323 # Beam Design #
                                                      383 addnewsolenoid(zi=tbench_solenoid_gun_zstart, zf=
324 #-----#
                                                          tbench_solenoid_gun_zend, ri=
325 # - size
                                                          tbench_solenoid_gun_radi, ro=
326 elec.a0 = Cathode_rado
                                                          tbench_solenoid_gun_rado, maxbz=
327 elec.b0 = Cathode_rado
                                                          tbench_solenoid_gun_b)
328 elec.ap0 = beamap0
                                                     384 # - Drift before main solenoid
329 elec.bp0 = beambp0
                                                     385 addnewdrft(zs=0.37, ze=0.60, ap=machine_piperad)
330 # - centroid
                                                     386 # - Main Solenoid
331 \text{ elec.x0} = \text{beamx0}
                                                     387 addnewsolenoid(zi=tbench_solenoid_main_zstart, zf
332 elec.xp0 = beamxp0
                                                          =tbench_solenoid_main_zend, ri=
333 elec.y0 = beamy0
                                                          tbench_solenoid_main_radi, ro=
334 elec.yp0 = beamyp0
                                                          tbench_solenoid_main_rado, maxbz=
335
                                                          tbench_solenoid_main_b)
336 #-----#
                                                     388 # - Drift after main solenoid
337 # Injection #
                                                     389 addnewdrft(zs=2.52, ze=2.67, ap=machine_piperad)
338 #-----#
                                                     390 # - Collector Solenoid
339 # --- Beam Injection ---#
                                                     391 addnewsolenoid(zi=tbench_solenoid_col_zstart, zf=
340 elec.npmax = npart
                                                          tbench_solenoid_col_zend, ri=
341 top.inject = machine_emittype
                                                          tbench_solenoid_col_radi, ro=
342 top.zinject[0] = beamzinject
                                                          tbench_solenoid_col_rado, maxbz=
343 if (machine_injtype=="gun"):
                                                          tbench_solenoid_col_b)
     elec.npinject = int(npart**2*sw*elec.sq/(elec. 392
344
      ibeam*timestep*nsteps))
                                                      393 # >>> Set input parameters describing the 3d
     print("number of particles injected per time
                                                          simulation.
345
                                                     394 w3d.nx = 32
     step: %g" % elec.npinject)
     top.xinject[0] = beamxinject
                                                     395 \text{ w3d.ny} = 32
346
     top.yinject[0] = beamyinject
                                                     396 \text{ w3d.nz} = 256
347
     top.xpinject[0] = beamxpinject
                                                     397 top.prwall = machine_piperad
348
     top.ypinject[0] = beamypinject
                                                     398
349
                                                     399 # >>> Set to finite beam.
     top.ainject[0] = elec.a0
350
     top.binject[0] = elec.b0
                                                     400 w3d.xmmin = -machine_piperad
351
     top.ainjmin[0] = Cathode_radi
                                                     401 w3d.xmmax = machine_piperad
352
     top.binjmin[0] = Cathode_radi
                                                     402 w3d.ymmin = -machine_piperad
353
     top.apinject[0] = beamapinject
                                                     403 w3d.ymmax = machine_piperad
354
     top.bpinject[0] = beambpinject
                                                     404 w3d.zmmin = machine_zstart
355
     top.vzinject[0,0] = 0.0
                                                     405 w3d.zmmax = machine_syslen
356
     top.vinject[0] = -500.0
                                                     406 dx = (w3d.xmmax-w3d.xmmin) / w3d.nx
357
                                                     407 dy = (w3d.ymmax-w3d.ymmin) / w3d.ny
358
359 # --- Profile Injection ---#
                                                     408 dz = (w3d.zmmax-w3d.zmmin) / w3d.nz
360 if machine_injtype == 'profile':
                                                     409
     xinit = posi[:,0] * mm #+ Xoff*mm
                                                     410 # >>> Set up some diagnostic windows.
361
     yinit = posi[:,1] * mm #+ Yoff*mm
                                                     411 top.xwindows[:,1] = [-5.e-2,5.e-2]
362
     zinit = zeros(npart)
                                                     412 top.zwindows[:,1] = [machine_zstart,2*elec.vbeam*
363
     vxinit = zeros(npart)
                                                          top.dt]
364
                                                     413 top.zwindows[:,2] = [machine_syslen/2-elec.vbeam*
     vyinit = zeros(npart)
365
     vzinit = zeros(npart) + elec.vbeam
                                                          top.dt, machine_syslen/2+elec.vbeam*top.dt]
366
                                                     414 top.zwindows[:,3] = [machine_syslen-2*elec.vbeam*
     def hollow_cathode_source():
367
     if w3d.inj_js == elec.jslist[0]:
                                                          top.dt, machine_syslen]
368
        w3d.npgrp = npart
                                                     415
369
        gchange('Setpwork3d')
                                                     416 # >>> Time histories
370
                                                     417 elec.nhist = int(nsteps/10)
371
        w3d.xt[:] = xinit
         w3d.yt[:] = yinit
                                                     418 top.ifzmmnt = 2
372
         w3d.zt[:] = top.zinject
                                                     419 top.itmomnts[0:3]=[0,nsteps,elec.nhist]
373
```

gun_cathode_radi=[None,Cathode_radcurvb,None, 476 None] gun_cathode=ZSrfrv(rsrf=gun_cathode_r,zsrf= 477 gun_cathode_z,rad=gun_cathode_radi,voltage= Cathode_voltage,xcent=.0e0,ycent=.0e0,zcent=.0 e0) 478 # - Anode 479gun_anode_r = [Anode_r1, Anode_r3, Anode_r4, Anode_r5, Anode_rendi, Anode_rendo, Anode_rendo, Anode_r2, Anode_r2, Anode_radtipo, Anode_radtipi] gun_anode_z = [Anode_z1, Anode_z3, 480 Anode_z4, Anode_z5, Anode_zend, Anode_zend, Anode_z5,Anode_z4,Anode_z2,Anode_zstart, Anode_zstart] gun_anode_radi = [None, None, None, None, None, 481 None, None, None, Anode_radcurvb, None, Anode_radcurvs] gun_anode = ZSrfrv(rsrf=gun_anode_r, 482 zsrf=gun_anode_z,rad=gun_anode_radi, voltage= Anode_voltage,xcent=.0e0,ycent=.0e0,zcent=.0e0) 483 # - Electrode F gun_electrodef_r = [ElectrodeF_r1, 484 ElectrodeF_ro,ElectrodeF_ro,ElectrodeF_ri, ElectrodeF_ri] gun_electrodef_z = [ElectrodeF zend. 485 ElectrodeF_z2,ElectrodeF_zstart, ElectrodeF_zstart,ElectrodeF_z1] gun_electrodef_radi = [ElectrodeF_radcurvb,None 486 ,None,None,ElectrodeF_radcurvs] gun_electrodef = ZSrfrv(rsrf= 487 gun_electrodef_r,zsrf=gun_electrodef_z,rad= gun_electrodef_radi,voltage=ElectrodeF_voltage ,xcent=0,ycent=0,zcent=.0e0) 488 # - Electrode C gun_electrodeC_r = [ElectrodeC ro. 489 ElectrodeC_ro,ElectrodeC_ri,ElectrodeC_ri] gun_electrodeC_z = [ElectrodeC_z1, 490 ElectrodeC_zstart, ElectrodeC_zstart, ElectrodeC_z1] gun_electrodeC_radi = [None,None,None, 491 ElectrodeC_radcurv] = ZSrfrv(rsrf= gun_electrodeC 492 gun_electrodeC_r,zsrf=gun_electrodeC_z,rad= gun_electrodeC_radi,voltage=ElectrodeC_voltage ,xcent=0,ycent=0,zcent=.0e0) gun_conductors=[gun_driftpipe,gun_driftpipe, 493 gun_cathode,gun_anode,gun_electrodef, gun_electrodeC] 494 installconductor(gun_conductors) 495 496 # --- Lattice --- # 497 pipe = ZCylinderOut(radius=machine_piperad,zlower =Gun_pipe_zstart,zupper=machine_syslen,voltage =0.,xcent=0,ycent=0,zcent=0)) 498 lattice_conductors=[pipe] 499 installconductor(lattice_conductors) 500 fieldsolve() 501 502 #-----# 503 # Plotting Lattice # 504 #-----#

```
505 # --- Plotting Envelope Function
                                                        540 fma()
506 penv(color="fg",marks=0,marker=None,msize=1.0,
                                                        541
    lframe=0,titles=1,ascale=None,bscale=None,zscale 542 # --- Plotting Electric Field
     =None)
507 fma()
508
                                                        545
509 # --- Plotting Potential
                                                        546
510 if (machine_injtype=="gun"):
                                                        547
     gun_driftpipe.draw(filled=190,color="fg")
511
                                                        548
     gun_cathode.draw(filled=160,color='fg')
512
                                                        549
     gun_anode.draw(filled=100,color='fg')
513
     gun_electrodef.draw(filled=150,color='fg')
514
                                                        550
     gun_electrodeC.draw(filled=150,color='fg')
515
516 pfzr(fullplane=1,plotsg=1,cond=1,fill=1,plotphi
                                                        551
     =1,plotrho=0,plotselfe=0,comp='z',titles=1)
   limits(Cathode_zstart,Gun_pipe_zend,-1.2*
                                                        553
517
    machine_piperad,1.2*machine_piperad)
   fma()
518
519
520 # --- Plotting Potential
                                                        557
521 if (machine_injtype=="gun"):
                                                        558
522
     gun_driftpipe.draw(filled=190,color="fg")
                                                        559
523
     gun_cathode.draw(filled=160,color='fg')
                                                        560
     gun_anode.draw(filled=100,color='fg')
524
                                                        561
     gun_electrodef.draw(filled=150,color='fg')
525
                                                        562
     gun_electrodeC.draw(filled=150,color='fg')
526
527 pfzr(fullplane=1,plotsg=1,cond=1,fill=1,plotphi
                                                        563
     =0,plotrho=1,plotselfe=0,comp='E',titles=1)
                                                        564
528 limits(Cathode_zstart,Gun_pipe_zend,-1.2*
                                                        565
    machine_piperad,1.2*machine_piperad)
                                                        566
529 fma()
                                                        567
530
                                                        568
531 # --- Plotting Potential
                                                        569
532 if (machine_injtype=="gun"):
                                                        570
     gun_driftpipe.draw(filled=190,color="fg")
                                                        571
533
     gun_cathode.draw(filled=160,color='fg')
                                                        572
     gun_anode.draw(filled=100,color='fg')
     gun_electrodef.draw(filled=150,color='fg')
536
     gun_electrodeC.draw(filled=150,color='fg')
537
538 pfzr(fullplane=1,plotsg=1,cond=1,fill=1,plotphi
     =0,plotrho=0,plotselfe=1,comp='E',titles=1)
539 limits(Cathode_zstart,Gun_pipe_zend,-1.2*
     machine_piperad,1.2*machine_piperad)
```

```
543 if (machine_injtype=="gun"):
     gun_driftpipe.draw(filled=190,color="fg")
     gun_cathode.draw(filled=160,color='fg')
     gun_anode.draw(filled=100,color='fg')
     gun_electrodef.draw(filled=150,color='fg')
     gun_electrodeC.draw(filled=150,color='fg')
  pfzr(fullplane=1,plotsg=1,cond=1,fill=1,plotphi
    =0,plotrho=0,plotselfe=1,comp='z',titles=1)
   limits(Cathode_zstart,Gun_pipe_zend,-1.2*
    machine_piperad,1.2*machine_piperad)
   fma()
   ## --- REPETITIVE PLOTS
554 def myplots():
     # --- Plotting Electric Field
     if (machine_injtype=="gun"):
       gun_driftpipe.draw(filled=190,color="fg")
       gun_cathode.draw(filled=160,color='fg')
       gun_anode.draw(filled=100,color='fg')
       gun_electrodef.draw(filled=150,color='fg')
       gun_electrodeC.draw(filled=150,color='fg')
     limits(Cathode_zstart,machine_syslen,-1.2*
      machine_piperad,1.2*machine_piperad)
     ppzx(color="density",ncolor=30)
     fma()
     pptrace(filled=1,particles=0,contours=30)
     fma()
   installplalways(myplots)
   # >>> run time steps and dump the results.
   step(nsteps)
   # step(50)
573 os.system("mv "+machine_type+"_"+date+time+"_"+
    machine_injtype+"_"+file_ending+"* ../Results/"+
```

date+"/")

Attachments/tbench complex app.py

Appendix D

Dimensions

D.1 Electron Guns

Hollow Electron Gun 1 Inch

Cathode	[mm]
Length	29.25
Inner Radius	6.75
Outer Radius	12.7
Radius of Curvature	10
Anode	[mm]
Distance to Cathode	9.48
Length	85.2
Inner radius	14.25
Outer Radius	$22 \mathrm{d}$
Anode radius of Curvature	3.5
Electrode F	[mm]
Length	74
Inner Radius	13.1
Outer Radius	15
Radius of Curvature	1.4
Electrode C	[mm]
Length	49.5
Inner Radius	20.5
Outer Radius	22
Radius of Curvature	0.75

Hollow Electron Gun 0.6 Inch

Cathode	[mm]
Length	e
Inner Radius	9
Outer Radius	15.24
Radius of Curvature	10
Anode	[mm]
Distance to Cathode	9.48
Length	90
Aperture	18
Outer Radius	40
Electrode F	[mm]
Length	38.9
Inner Radius	8.05
Outer Radius	11.1
Electrode C	[mm]
Length	20.5
Inner Radius	17.0
Outer Radius	20.5

D.2 Electron Lenses

Tevatron Electron Lens 2

Bend Solenoids	[mm]
Width	90
Inner Radius	193
Outer Radius	265
Main Solenoid	[mm]
Length	2688.5
Inner Radius	200
Outer Radius	344.8
Collector and Gun Solenoid	[mm]
Length	500
Inner Radius	280
Outer Radius	334.3
Path lengths	[mm]
Main Solenoid - Nearest Bend	82.7
Inter Bend Distance	52.9
Cathode Surface - First Bend	281.6

Tevatron Electron Lens Test Stand

Main Solenoid	[mm]
Length	1920
Inner Radius	100
Collector and Gun Solenoid	[mm]
Length	500
Inner Radius	140
Drift Pipe	[mm]
Length	2860
Inner Radius	30
Path lengths	[mm]
Main Solenoid - Nearest Bend	82.7
Inter Bend Distance	52.9
Cathode Surface - First Bend	281.6

Appendix E

Maps

E.1 Map of LHC Collimators



Figure E.1: The super-symmetry of the LHC. IP1, IP2, IP5 and IP8 are the experimental halls ALICE, ATLAS, CMS and LHC-b respectively. IP3 and IP7 are the momentum and betatron cleaning sites and IP6 contains the beam dumps.

E.2 Map of Tevatron



Figure E.2: Map of the tevatron and its various sectors