

Particle and Ray Tracing Codes (PARMILA & TURTLE Introduction)

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Presentation Outline

Particle & Ray Tracing Codes

1. Beam Dynamics Codes Used in LINAC Design
2. Introduction to TURTLE
3. Introduction to PARMILA - For Simulation of Transfer Lines
 ⇒ Other Applications of PARMILA Later in Course
4. Space Charge Modeling in PARMILA
5. Using PARMILA & TURTLE to Study Some Beamlines
 ⇒ You will use the Simulation Lab computers in the classroom

1. Beam Dynamics Codes Used in LINAC Design

- Very Large Number of Programs Used in Accelerator Community
 - More Than 200 Codes Listed in the Los Alamos Accelerator Code Group (LAACG) Compendium of Accelerator Codes (~1990)
 - Optics Codes Account for About 1/2
 - The LAACG Compendium Represents a Lower Bound!
- Restrict Comments to Codes of Interest for LINAC Design & Low β

"Design Codes" vs "Simulation Codes"
"Envelope Codes" vs "Particle Codes"

- TRACE 3-D and TRANSPORT are "Design Codes"
 - Parameter Variation and Constraint Fitting
 - Fast Running by Using Beam Envelopes (No Particles)
- To Evaluate a Design One Needs a "Particle" or "Ray Tracing" Code
 - Simulates a Group of Macroparticles Transiting the Beamlne
 - "Simulation Codes" Used for Performance Evaluation of Design
- Three "Simulation Codes" on **Simulation Lab computers**
 ⇒ **PARMILA, PARMTEQ, TURTLE**
- **PARMILA & PARMTEQ** are both "Design Codes" & "Simulation Codes"
 - Can Produce Structure Layouts & Designs - no envelope calculations

1. Beam Dynamics Codes Used in LINAC Design (con't)

- **Particle Codes**
 - Space Charge is a Major Consideration in the Design of a Code
 - How Fields Are Modeled is a Major Consideration
 - 2-D versus 3-D Simulation is Also a Major Consideration
- **Several Advanced Codes in Varied Use for LINAC Simulation**
- **One Example is WARP (LBNL, LLNL, NRL)**
 - PIC (Particle-in-Cell) Code
 - Detailed Description of External Fields (magnets, cavities,...)
 - Parallel (& Serial) Processing Employed
 - Applied to Heavy Ion Fusion (HIF) and Intense Electron Beams
- **Several Other Advanced Codes (PIC and EM) in DOE SciDAC Program**
 - SciDAC: Scientific Discovery through Advanced Computing
 - Capable of Simulating $\sim 10^{19}$ Particles / Bunch through Detailed Fields
 - Requires Massively Parallel Computer Clusters on the "Petascale"
- **Use of These Advanced Codes can be Complex**
 - Assistance of Code Developers Typically Required
 - ⇒ Not as "User Friendly" as the Codes in the Simulation Lab!

1. Beam Dynamics Codes Used in LINAC Design (cont'd)

- For completeness we should mention a few Electron Linac codes
- PARMELA
 - "Phase And Radial Motion for Electron Linear Accelerators"
 - Los Alamos National Laboratory (LANL) code
 - Patterned after PARMILA, Includes Space Charge
 - Used Primarily for Low Energy Electron Linac Design
- LIAR
 - Linear Accelerator Research code
 - Stanford Linear Accelerator Center (SLAC)
 - Used Primarily for High Energy Electron Linacs
 - Includes the DIMAD code for Tracking Calculations
- DIMAD - a TRANSPORT spin-off with Tracking Capabilities
 - SLAC & Thomas Jefferson National Accelerator Facility (TJNAF)
- Several Specialized Codes for Electron Guns (Ion Sources)
- Some of the Codes in the Simulation Lab are Often Used for Electrons:
TRACE 3-D, SUPERFISH, TRANSPORT, TURTLE

1. Beam Dynamics Codes Used in LINAC Design (cont'd)

Some Electron Linac Codes

Code Name	Version(s)	Website(s)	Ref.
PARMELA*	LANL	http://laacg1.lanl.gov	[E-1]
LIAR*	SLAC	http://www-project.slac.stanford.edu/lc/local/AccelPhysics/Accel_Physics_index.htm	[E-2]
DIMAD*	SLAC	http://www-project.slac.stanford.edu/lc/local/AccelPhysics/Accel_Physics_index.htm	[E-3]
TRACE 3-D	LANL, PBO Lab Module†	http://laacg1.lanl.gov http://www.ghga.com/accelsoft	[2]
SUPERFISH	LANL	http://laacg1.lanl.gov	[8]
TRANSPORT	PBO Lab Module (FNAL)	http://www.ghga.com/accelsoft http://fermitools.fnal.gov	[1]
TURTLE	PBO Lab Module (FNAL)	http://www.ghga.com/accelsoft http://fermitools.fnal.gov	[3]

*These electron codes are *not* being used in this USPAS course

† Includes electrostatic elements

References in Preceding Table

- [E-1] L. M. Young, "PARMELA," Los Alamos National Laboratory Report No. LA-UR-96-1835 (1996).
- [E-2] R. Assmann, C Adolphsen, K. Bane, P. Emma, L. Hendrickson, F. Ostiguy, T. Raubenheimer, A. Seryi, R. Siemann, G. Stupakov, P. Tenenbaum, K. Thompson and F. Zimmerman "LIAR - A Computer Program for the Modeling and Simulation of High Performance Linacs," Stanford Linear Accelerator Center, SLAC/AP-103, 3 February 2003, 131 pages (2003).
- [E-3] R. V. Servranckx, K. L. Brown, L. Schachinger, D. Douglas and P. G. Tenenbaum, "User's Guide to the Program DIMAD," Stanford Linear Accelerator Center, 20 April 2001, 57 pp (2001); R. V. Servranckx, K. L. Brown, L. Schachinger, D. Douglas and P. G. Tenenbaum, "User's Guide to the Program DIMAD," Stanford Linear Accelerator Center, 14 January 2004, 48 pp (2004).
- [1] D. C. Carey, K. L. Brown and F. Rothacker, "Third-Order TRANSPORT with MAD Input - A Computer Program for Designing Charged Particle Beam Transport Systems," SLAC-R-530, 316 pp (1998).
- [2] K. R. Crandall and D. P. Rusthoi, "TRACE 3-D Documentation," Third Edition, Los Alamos National Laboratory Report No. LA-UR-97-886, 106 pages (1997).
- [3] D. C. Carey, "TURTLE with MAD Input (Trace Unlimited Rays Through Lumped Elements), a Computer Program for Simulating Charged Particle Beam Transport Systems, and DECAY-TURTLE Including Decay Calculations," Fermilab-Pub-99/232, 196 pp (1999).
- [8] M. T. Menzel and H. K. Stokes, "User's Guide for the POISSON/SUPERFISH Group of Codes," Los Alamos National Laboratory Report No. LA-UR-87-115, ~250 pages (1987); The Los Alamos Accelerator Code Group, "Reference Manual for the POISSON/SUPERFISH Group of Codes," Los Alamos National Laboratory Report No. LA-UR-87-126, ~400 pages (1987).

2. Introduction to TURTLE

- **Acronym: "Trace Unlimited Rays Through Lumped Elements"**
- **Companion Code to TRANSPORT**
- **"Simulation Code" Rather Than a "Design Code"**
 - Calculates the 6-D coordinates of macroparticles through beamline
 - Same 6-D coordinates as TRANSPORT: $(q_i) = (x, x', y, y', l, \delta)$
 - Can routinely run 1,000 - 100,000+ particles on a PC
 - No parameter variation, fitting, matching, ...
- **Can examine effects not modeled in TRANSPORT**
 - Three distributions in standard TURTLE:
Gaussian, Rectangular, Uniform
 - Alternatively can import initial particle distribution from a file
 - Beam loss on apertures and slits
 - Non elliptical beams - either as input, or created in a beamline
 - Can add phase space "cuts" to distributions along beamline
- **Output Used by Some Radiation Transport Codes - Shielding**
- **No Space Charge**

2. Introduction to TURTLE (continued)

- **TURTLE is complimentary program to TRANSPORT**
 - Both use same description of beamline and elements
 - 1st, 2nd, 3rd order calculated in both
 - Use the same 6-D phase space coordinates
 - ⇒ **Use TRANSPORT to get a design; use TURTLE to evaluate it**
- A special version, called DEACY-TURTLE, for radioactive beams
 - Tracks particles that have a finite lifetime
 - Primary particles (parents) decay randomly during beamline passage according to their lifetime
 - Tracks the decay product particles (daughters) thru beamline
 - User specifies parent lifetime, masses and charges of daughters
 - Transparency of apertures and slits can be particle dependent
- PBO-Lab provides TURTLE support for the same "non-TRANSPORT" optics elements that PBO-Lab makes available for TRANSPORT
 - TRACE 3-D elements: Doublet, Triplet, RF Gap, Thin Lens
 - Extended fringe-field elements: Magnetic Quad, ElectroStatic Quad

2. Introduction to TURTLE (continued)

- Generates Data for 2 Types of Plots
 - Scatter Plots of Phase Space Data (2-D Plot) e.g. x' versus x
 - Histograms of Data (1-D Plot) e.g. # particle is at bin location x
- ⇒ **Plot Specifications for TURTLE set in PBO Lab Marker Piece**
- Any of the 6-D Phase Space Variables Can be Used x, x', y, y', l, δ
 - 36 Possible Scatter Plots
 - 6 Possible Histograms
- Native Plots are Text Based - "Line Printer" (without a post processor)
- Multiple Locations in Beam Line - User Specified Plots at Each
- Apertures (on magnets) Can Be Ignored or Included (Particle Loss)
- DECY-TURTLE Version Has Above Capabilities for Decay Particles
 - But DECY-TURTLE **does not support all 3rd order modeling**

⇒ **Next week (time permitting) will use TURTLE to simulate the use of octupoles in a high energy beam transport (HEBT) line to create a nearly uniform spot density from a Gaussian beam**

Before discussing PARMILA, will look at some typical TURTLE output plots

2. Introduction to TURTLE (continued)

Typical Scatter Plot (TURTLE native output)

```

+          -0.100   -0.060   -0.020    0.020    0.060    0.100
+          TOTALS
+*****-----*-----*-----*-----*-----*-----*-----*-----*-----*
-0.003 TO -0.003 I  1   1 1131 12           B           I   22
-0.003 TO -0.003 I           11 2241 1           711          I   21
-0.003 TO -0.002 I           1 1 33 2           6           I   16
-0.002 TO -0.002 I           113 2321          6           I   19
-0.002 TO -0.002 I           1 255433          E1          I   38
-0.002 TO -0.002 I           1327589443          1B          I   58
-0.002 TO -0.002 I           2 2 4562711          I           I   48
-0.002 TO -0.001 I           4247945521          E2          I   59
-0.001 TO -0.001 I           1 295BDE8751          I31          I   99
-0.001 TO -0.001 I           416P$USD711          9B 1          I   260
-0.001 TO -0.001 I           8B$$$$$XI1          R3  1          I 1182
-0.001 TO -0.001 I           12D$$F$$$S$8          V2           I   950
-0.001 TO 0.000 I           14Q$  1X$$5          X121          I   727
0.000 TO 0.000 I           22R$  3$$E          $7           I   665
0.000 TO 0.000 I           3V$   $$$          $91          I   638
0.000 TO 0.000 I           1B$   $$$          $R211         I   616
0.000 TO 0.000 I           13L   N$$5          $W  1          I   636
0.000 TO 0.001 I           1 T   9$$$          $Q721         I   753
0.001 TO 0.001 I           13P   B$$$UA$$U2          1 I   929
0.001 TO 0.001 I           T   15W$$$$$62 11          I   1194
0.001 TO 0.001 I           1 E6   47EP$$J854          I   262
0.001 TO 0.001 I           2J   1ACAI9651 212          I   98
0.001 TO 0.002 I           J   322647823          I   56
0.002 TO 0.002 I           K   27587623          1           I   61
0.002 TO 0.002 I           D3   1124731132 2 1          I   44
0.002 TO 0.002 I           5B   1 351331222          I   39
0.002 TO 0.002 I           1A   43312          I   24
0.002 TO 0.003 I           9   22 4112 1   1          I   23
0.003 TO 0.003 I           54   1 11 5 111          I   20
0.003 TO 0.003 I           14   1 11 3   1          I   12
+*****-----*-----*-----*-----*-----*-----*-----*-----*-----*
I
I
I
I   1754343443444443334471
I   12555118093180211299858384211
TOTALS I 00110002247295757521144381220762055313246322020001 I   9569
0   TOTAL NUMBER OF ENTRIES =      10000 INCLUDING UNDERFLOW AND OVERFLOW AS FOLLOWS
0
0   LEFT          RIGHT
ABOVE       38           173            3
                  1           9569            0
BELOW        8            174            34
SUM OF SQUARES =      1777943.
CENTER =      0.000   RMS HALF WIDTH =      0.032
CENTER =      0.000   RMS HALF WIDTH =      0.002
CORRELATION =   0.2913
ONO 4 TWO DIMENSIONAL PLOT OF
HORIZONTAL AXIS Y IN M      14.608 M   FROM THE START
VERTICAL AXIS Y' IN R      14.608 M   FROM THE START

```

[MS Word font: Courier New 7 pt., normal spacing]

2. Introduction to TURTLE (continued)

Typical Histogram (TURTLE native output)

```

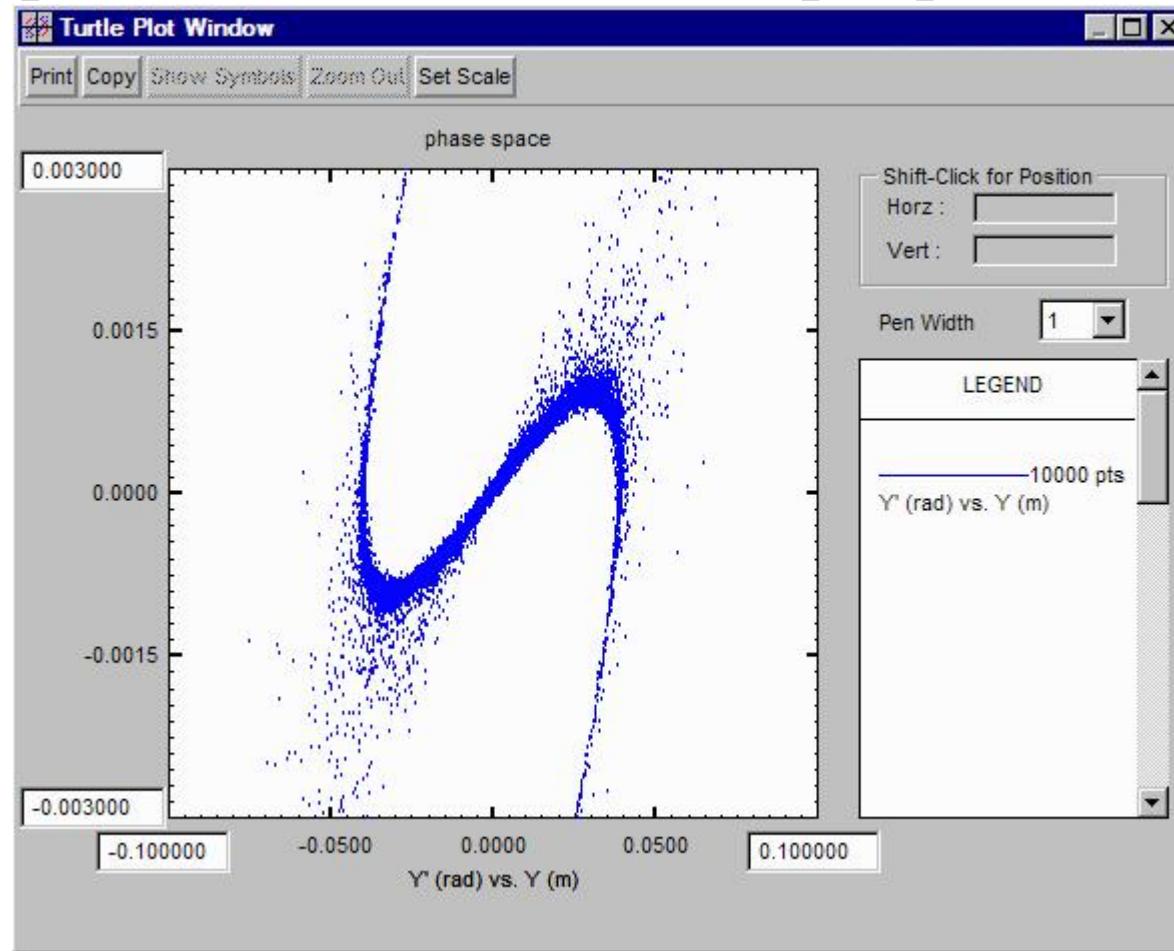
DISTRIBUTION OF      Y IN M      14.608 M   FROM THE START
0
INTERVAL           SCALE FACTOR: 100 X'S EQUAL    908 RAYS
0LESS THAN -0.100    47    XXXXX
-0.100 TO -0.095     3
-0.095 TO -0.090     8
-0.090 TO -0.085     7
-0.085 TO -0.080     3
-0.080 TO -0.075    11  X
-0.075 TO -0.070     8
-0.070 TO -0.065    10  X
-0.065 TO -0.060    15  X
-0.060 TO -0.055    13  X
-0.055 TO -0.050    29  XXX
-0.050 TO -0.045    55  XXXXX
-0.045 TO -0.040   184  XXXXXXXXXXXXXXXXXXXXXXXX
-0.040 TO -0.035   908  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
-0.035 TO -0.030   591  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
-0.030 TO -0.025   491  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
-0.025 TO -0.020   510  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
-0.020 TO -0.015   529  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
-0.015 TO -0.010   527  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
-0.010 TO -0.005   514  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
-0.005 TO 0.000    520  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
0.000 TO 0.005    548  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
0.005 TO 0.010    508  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
0.010 TO 0.015    535  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
0.015 TO 0.020    515  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
0.020 TO 0.025    524  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
0.025 TO 0.030    507  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
0.030 TO 0.035    607  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
0.035 TO 0.040    859  XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX
0.040 TO 0.045    205  XXXXXXXXXXXXXXXXXXXXXXXXX
0.045 TO 0.050    47   XXXXX
0.050 TO 0.055    28   XXX
0.055 TO 0.060    24   XX
0.060 TO 0.065    17   X
0.065 TO 0.070     9
0.070 TO 0.075    19   XX
0.075 TO 0.080     6
0.080 TO 0.085     9
0.085 TO 0.090     4
0.090 TO 0.095     4
0.095 TO 0.100     5
0GREATER THAN 0.100 37   XXXX
0      TOTAL NUMBER OF ENTRIES = 10000      INCLUDING UNDERFLOW AND OVERFLOW
0      CENTER = 0.000      RMS HALF WIDTH = 0.032

```

[MS Word font: Courier New 7 pt., normal spacing]

2. Introduction to TURTLE (continued)

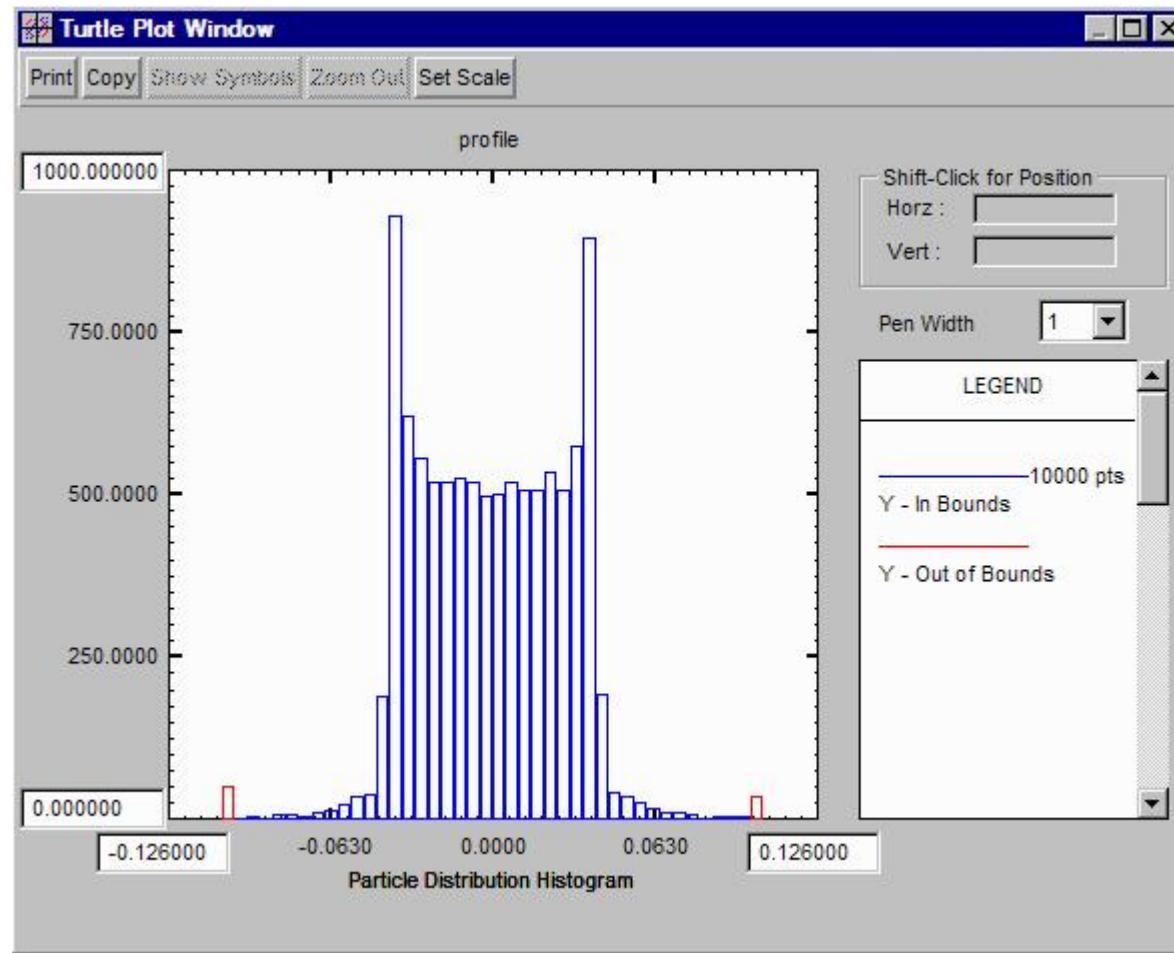
Typical Scatter Plot (PBO Lab postprocessor plot)



- PBO-Lab uses same (default) color assignments for all codes
- Provide for "Auto-Scaling" of Plots - Useful for Initial Setup
- Can Adjust Scales After a Run- Unlimited "Zoom" Capability

2. Introduction to TURTLE (continued)

Typical Histogram (PBO Lab postprocessor plot)



- PBO-Lab plots provide data visualization & document preparation tools
- Print, Copy, scale adjustment, ...
- Need to look at TURTLE (text) output for selected quantitative data

3. Introduction to PARMILA - For Simulation of Transfer Lines

- PARMILA is Both a Simulation Code and a Design Code ("Dual" Modes)
 - Much More than Just a Companion to TRACE 3-D
 - First Version Written by Don Swenson (~1963) for DTL Modeling
 - "PARMILA 1" (~1990) Transport Lines, Matching Sections, DTLs
 - "PARMILA 2" (~1998) Several New RF Structures / Capabilities
 - Coupled Cavity Linac (CCL)
 - Coupled Cavity DTL (CCDTL)
 - Superconducting Cavity (SC) Linac (2 types)
 - Funnels, Choppers, Improved Space Charge, ...
- Simulation Mode - Can be used for Transfer Lines, Matching Sections
 - Solves (Numerically “Integrates”) the Particle Equations of Motion
 - Conceptually Similar to Differential Matrix Model of Components
 - Advances Individual Macroparticles in Steps
 - Space Charge Calculated & Applied at Certain Steps - "Scheff"
 - Particles Removed (& Logged) Once Outside an
 - Inherently a Nonlinear Code (especially space charge)
⇒ Use TRACE 3-D to get a design; use PARMILA to evaluate it
- Design Mode - Used for DTLs, CCLs, Superconducting Structures
 - Will Discuss This in Detail Later in Course
 - For Now, Just a Brief Overview

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

- **Design Mode**
 - PARMILA's Most Important Capabilities - Few Other Codes Have This
- **How Does the Design Mode Work for a DTL?**
 - User Specifies Certain "Goals" for the DTL Design
 - e.g. Final Output Energy (but also other options for DTL goals)
 - User Sets Some Initial Values for Various Parameters
 - e.g. Initial Quadrupole Strength at DTL Entrance
 - Options to Specify "Ramping" or Changes with Distance
 - e.g. Ramp Accelerating Gradient
 - ... Other DTL Criteria/Options that will be Discussed Later
- **When PARMILA is Run in Design Mode:**
 - A DTL LINAC Design is Generated (Subroutines GENLINx)
 - The Design is Output to Several Text Files
- **With PARMILA 2 Several Other Capabilities**
 - Design of CCL, CCDTL, and Superconducting Structures
 - A TRACE 3-D Equivalent Lattice (of DTL, CCL, ...) is Output to a File
 - Other Files Generated for use in Graphics Displays
 - Can Run an "End-to-End" Simulation of Complete LINAC (after RFQ)

⇒ **Design Mode is Similar to an "Expert System"**

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

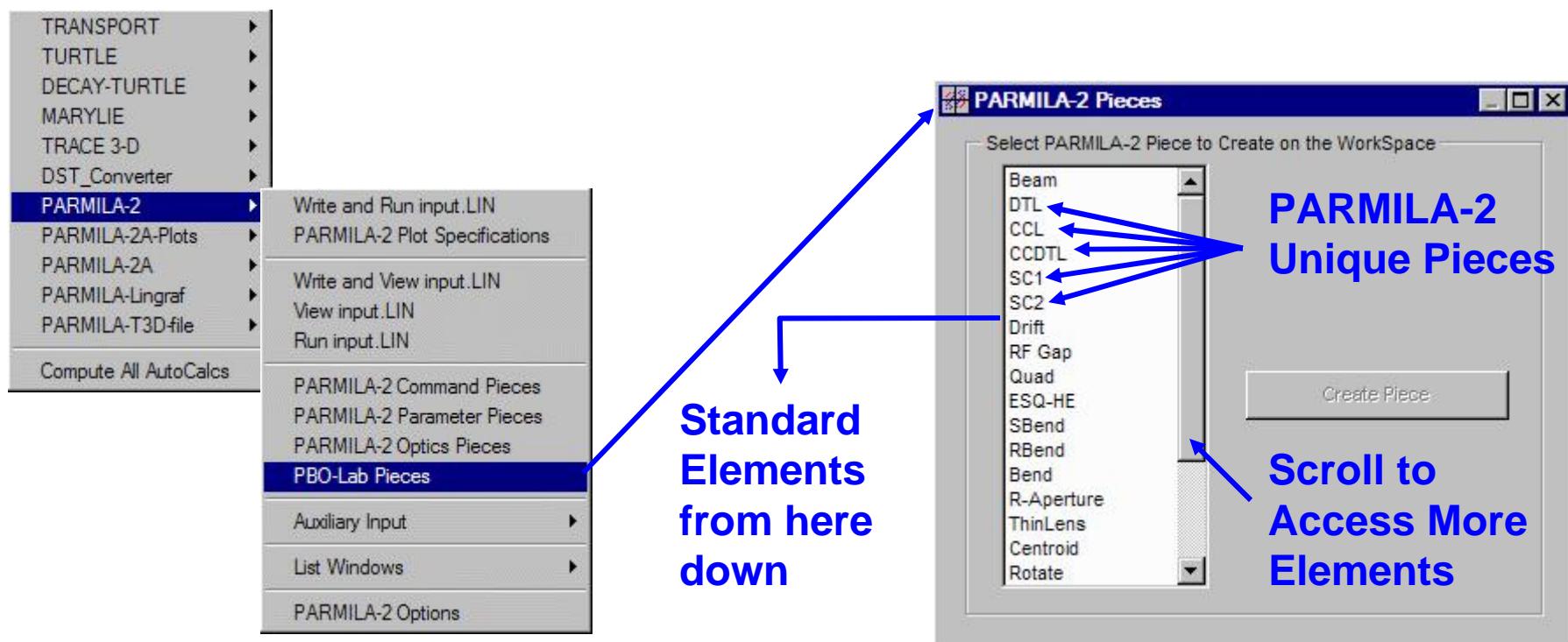
Optical Elements in PARMILA-2

- Standard Beamline Optical Elements (~12)
 - Eight are Essentially Same as TRACE 3-D / TRANSPORT Elements:
Drift, Quad, Bend (with Edges), **Solenoid***,
RF Gap, Thin Lens, Rotate (Displace & Tilt elements)
 - Steering Magnet ("Kicker") - Similar to TRANSPORT, TURTLE
 - Element Displacement - Similar to TRACE 3-D, TRANSPORT, TURTLE
 - Apertures (Circular, Rectangular) - Similar to TURTLE
- ***Solenoid in PARMILA version 2.36 does not work correctly**
⇒ **Use PBO Lab PARMILA-2 Module - it has a fix**
- Magnetic Quadrupole with Higher-Order Multipoles ("Mpoles")
- A Few Additional "Element" Type Items (i.e. Location Specific)
 - Funnel: Doubles RF Frequency and Beam Current (power only)
 - Chopper: Reduces Average Beam Current (power only)
 - Chopper: Reduces Average Beam Current (power only)
- Documentation (LA-UR-98-4478) Says Electrostatic (ES) Elements:
 - **ES Quadrupole, ES Deflector, DC Column**
 - **However, these are not in PARMILA version 2.36**

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

Simulating Transfer Lines in PARMILA-2 with PBO Lab

- For a PBO Lab Model of Standard Beamline Optical Elements
⇒ The Beamlne on the Model Space is Already "PARMILA-2 Complete"
- What "Standard Beamline Optical Elements" are supported?
- Get List from PARMILA-2 Menu "PBO Lab Pieces":



3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

Simulating Transfer Lines in PARMILA-2 with PBO Lab

- PBO Lab will write PARMILA-2 formatted element parameters (input.LIN file)
- PBO Lab also writes **automatically** other input lines **required** by PARMILA-2
 - Will look at these "other" required input lines shortly
 - Just note that PBO Lab provides **default** data for required input lines
- Use PARMILA-2 Command "Write and View 'input.LIN' " to see the results
- Use PARMILA-2 Command "Write and Run 'input.LIN' " to execute
⇒ **PBO Lab will produce a default "Simulation" for the transfer line**
- Error messages may be generated by PARMILA-2 if defaults not adequate
- Note: PARMILA-2 input file format **significantly different** from PARMILA-1
⇒ **PARMILA-2 cannot read PARMILA-1 input files**
- Let's look at a transfer line (Example B with RF gaps) **input file** for PARMILA-2

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

Transfer Line Input File ("input.LIN") for PARMILA-2

```

run 11 <----- Generally Starts File
title <----- Title indicates next line is not to be executed
Example B.pbol run on 01/15/11 07:04:56 <----- PBO Lab automatically writes a Title with date & time
; *** PARMILA OASIS Module Version 2.2.1.4 for PBO Lab 3.0 *** <----- ; means a comment follows (until next line)
; linac: W0, Fbunch, Ibeam, Mc^2, Nq <----- PBO Lab has a liberal use of comments
linac 2.00000000 80.00000000 100.00000000 1875.00000000 1.00000000 <----- linac = Global Parameters:
                                                               Initial Beam Energy MeV
                                                               Radio-Frequency (MHz)
                                                               Current (mA)
                                                               Particle Mass (MeV)
                                                               Particle Charge (|e|)

; ***** End of the input.lin HEADER SECTION ***** <----- PBO Lab extensive
; *** User Auxillary HEADER INPUT will appear here *** <----- use of comments!
; ***** Start of the input.lin LATTICE SECTION *****

; Initial beam parameters data format:
; type, number of macroparticles, alpha-x, beta-x (cm/rad), emittance-x (pi-cm-rad, boundary)
; line 2: alpha-y, beta-y (cm/rad), emittance-y (pi-cm-rad, boundary)
; line 3: alpha-z, beta-z (cm/rad), emittance-z (pi-cm-rad, boundary)
; line 4: centroids: x (cm), xp (rad), y(cm), yp (rad), del-phi (degrees), del-E (MeV)
;; input 0 1000 3.11400000 75.36000000 0.00600000
; input -0 1000 3.11400000 75.36000000 0.00600000 <----- number of macroparticles (PBO Lab Global)

input -8 1000 3.11400000 75.36000000 0.00600000 <----- input: (4 lines) Initial Twiss Parameters for X, Y, Z
      -2.62020000 57.89000000 0.00600000 <----- -8 = Uniform in 3-D Spatial Coords. (-) Fit to Twiss:
      -0.12100000 109.22502084 0.01201661 <----- Alpha, Beta, Emittance (Attention to UNITS, "0.0"!)
      0.00000000 0.00000000 0.00000000 0.000 0.00000000 0.500000 <----- Beam Centroids (see comments)
                                                               Random number seed

; ReadDist 0.000000 part_rfq.dst <----- Reads Input Beam Data from "part_rfq.dst" File (IF Uncommented)

```

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

Transfer Line Input File ("input.LIN") for PARMILA-2 (con't)

- Setting up the "input" (initial beam) line can be challenging
- PARMILA-2 has **numerous distribution possibilities** (not quite limitless!)
 - Nine basic "types" (0-8) with **numerous** variants
 - Basic type & variants set using a **signed 3-digit Type code (lmn)**
 - The last digit (n) **defines the Basic "type"** described in Documentation
 - The first two digits (lm) are **not commonly used** (no entry, only n)
 - Minus ($-lmn$) **adjusts random distributions to achieve exact (Twiss) values**
 - Units for longitudinal Twiss & emittance **differ for different "type"**
 - Documentation (LA-UR-98-4478) not always "clear"
- PBO Lab Beam Piece is useful for some commonly used distributions
 - PBO Lab: 6-D Equivalent Uniform
 - "-8" PARMILA-2 Type Code
 - Uniform in Spatial Coordinates
 - and** Uniform in Divergences
 - Emittances are **Boundary Values**
 - PBO Lab: 4-D Equivalent Uniform (KV) "-6" PARMILA-2 Type Code
 - PBO Lab: Gaussian (5-Sigma) "-52" PARMILA-2 Type Code
- PBO Lab "input" line "type" can be **overridden with PARMILA-2 Options**
- PBO Lab **PARMILA-2 Options** used to read distribution from file, set file name



3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

Transfer Line Input File ("input.LIN") for PARMILA-2 (con't)

PBO Lab Beam Piece writes several additional lines required for simulations:

```
STRUCTURE 0 0 80.00000000 ← STRUCTURE defines a new PARMILA-2 Linac Section (DTL, ...)  

TRANSPORT ← TRANSPORT specifies the type of PARMILA-2 Structure  

Drift 0.00000000 20000.00000000 0 1 ← Drift is a zero length drift (not required, but useful for output)  

Bore 20000.00000000 ← Bore sets an radial aperture (in cm) for removing particles  

Scheff 0.05000000 0.05000000 9 9 0 1 1 ← Scheff necessary for simulation (more about this later)  

PrtBeam 99.00000000 ← PrtBeam writes out 99% rms beam data ("Beam.out" file)  

OUTPUT 2 1 1 1000 1 ← OUTPUT specifies how often to output data ("part_dtl.dst" file)  

; End of PBO Lab Beam Piece data, including STRUCTURE 0 for writing input beam to Parmila.out
```

PBO Lab also has a PARMILA-2 Command Piece named
 "Start a TRANSPORT STRUCTURE"

Needed for matching sections between complex Structures (DTL, CCL, SCL)

Not necessary for a simple transfer line simulation, but this is an example

This Command Piece writes 4 data lines (& 1 comment line):

```
BEGIN ; starts beam dynamics calculation for prior STRUCTURE  

;***** Start of New STRUCTURE - PBO Lab PARMILA-2 Module *****  

STRUCTURE 1 0 80.00000000 80.00000000 ← 1 = Structure ID (unique, increasing)  

TRANSPORT 0.00000000 ; TRANS ← 0 = Specifies the space charge routine  

QuadCalc 1 1 1 ← 80.0 = Radiofrequency (MHz) for Section  

                           80.0 = SF Data Radiofrequency (not used here)  

                           0.0 = Starting RF Phase (degrees) for initial beam distribution  

                           QuadCalc = Activates quadrupole simulation options:  

                           Fringe Fields  

                           Third Order  

                           Conserve Energy
```

- PBO Lab writes certain default "input.LIN" lines so PARMILA-2 will execute

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

Transfer Line Input File ("input.LIN") for PARMILA-2 (con't)

PBO Lab writes optics elements in format and units needed by PARMILA -2:

Drift 17.30000000 20000.00000000 1 1 ; Comment: D4

This example has optics elements for:

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

Drift

Quad 9.60000000 1.00000000 1 -2650.00000000 0 ; Comment: Q5

Quad

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

Cavity (RF Gap)

Drift 3.97500000 20000.00000000 1 1 ; Comment: D6

Commented (;) lines are Piece Parameters
that user has elected to not use.
This example has quadrupole *rolls* and
displacements (;Displace) turned off.

Cavity 0.0 20000. 1 0.18150000 80.00000000 -40.00000000 ; Comment: G7

Drift 3.97500000 20000.00000000 1 1 ; Comment: D8

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

Suggestion: Use PBO Lab Piece Comment
field to help identify elements in "input.LIN"

Quad 9.60000000 1.00000000 1 2650.00000000 0 ; Comment: Q9

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

Drift 4.27400000 20000.00000000 1 1 ; Comment: D10 ;

Cavity 0.0 20000. 1 0.18150000 80.00000000 -40.00000000 ; Comment: G11

Drift 4.27400000 20000.00000000 1 1 ; Comment: D12

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

Third entry (0 or 1) controls output to
"Parmila.plt" binary file used by Lingraf.
Suggestion: Use PBO Lab PARMILA-2
Option "OutputFlag" to turn ALL on (or off).

Quad 9.60000000 1.00000000 1 -2650.00000000 0 ; Comment: Q13

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

Drift 4.57400000 20000.00000000 1 1 ; Comment: D14

Cavity 0.0 20000. 1 0.18150000 80.00000000 -40.00000000 ; Comment: G15

Drift 4.57400000 20000.00000000 1 1 ; Comment: D16

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

Quad 9.60000000 1.00000000 1 -2650.00000000 0 ; Comment: Q17

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

Drift 4.87300000 20000.00000000 1 1 ; Comment: D18

Cavity 0.0 20000. 1 0.18150000 80.00000000 -40.00000000 ; Comment: G19

Drift 4.87300000 20000.00000000 1 1 ; Comment: D20

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

Quad 4.80000000 1.00000000 1 -2650.00000000 0 ; Comment: Q21

;Displace 0.0 0.0 1 0.0 0.0 0.00000000

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

Transfer Line Input File ("input.LIN") for PARMILA-2 (con't)

PBO Lab has custom PARMILA -2 Command Pieces and Parameter Pieces to specify space charge details, cutoff parameters, output details, etc.

```
Bore 2000.00000000
; SC: delta-R delta-L N-rad N-long N-adj N-beta-lambda Remesh
Scheff 0.05000000 0.05000000 9 9 0 1 1
PrtBeam 99.00000000
; OUTPUT 1 1 1 2 3 4 5 6 7 8
OUTPUT 2 1 1 1000 1
; OUTPUT 3 1 1 2 3 4
; OUTPUT 4 1 1 2 3 4
```

This group is similar to the "default" lines PBO Lab writes for "STRUCTURE 0" however user sets the parameter values

PBO Lab automatically writes needed simulation "begin" and "end" lines

```
begin
end
```

- **SCHEFF ("Space Charge Effects") line needed for PARMILA-2 simulations**
- SCHEFF, together with optional lines, control space charge calculations
- Will look at the space charge models and options available, but first look at some typical PARMILA-2 outputs for the above transfer line

3. Introduction to PARMILA - For Simulation of Transfer Lines

PARMILA-2 Outputs - Example B with RF Gaps

Main Output File: "Parmila.out"

Los Alamos National Laboratory Parmila - Ion Linac Design Code
Program Parmila written by Harunori Takeda

This program is provided as a service to the accelerator community by the Los Alamos Accelerator Code Group (LAACG).
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Program Parmila 2.35 released 7-25-2005

Parmila starting at 01-16-2011 14:28:25
Parmila input file: input.lin
LANL.INI file: C:\OASIS\LANL.INI 8-11-2010 16:59:46
First random number (using system clock) for linac error studies = 0.251825749465751880

Lines that start with "LIN:" are from file input.lin:

LIN: run 1 1

[Reiteration of parts of input file](#)

```
*****
Example B.pbol run on 01/16/11 14:28:23
*****
```

LIN: linac 2.00000000 80.00000000 100.00000000 1875.00000000 1.00000000
Maximum numbers of Particles: 90000
Cells: 10000
Tanks: 25
Structures: 50
Space-charge mesh size: 80 (radial), 200 (longitudinal)
LIN: input -8 1000 3.11400000 75.36000000 0.00600000
LIN: -2.62020000 57.89000000 0.00600000
LIN: -0.12100000 109.22502084 0.01201661
LIN: 0.00000000 0.00000000 0.00000000 0.000 0.00000000 0.500000
• • •
LIN: PrtBeam 99.00000000
LIN: OUTPUT 2 1 1 1000 1
LIN: BEGIN ; starts beam dynamics calculation for prior STRUCTURE
Starting beam dynamics... Current = 100.0 mA
Space charge calculations use the Scheff algorithm.

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

PARMILA-2 Outputs - Example B with RF Gaps (con't)

Main Output File: "Parmila.out" (con't)

Starting Transport section

```
Wsync =      2.0000 MeV, 1000 good particles
Beam centroid: x (cm)  xp (mr)  y (cm)  yp (mr)  phi (deg)  W (MeV)
          0.00000  0.00000  0.00000  0.00000      0.000    2.000
Phis =      0.0000 deg
Transport finished element 1: 1. Drift
```

Longitudinal normalized emittance is in deg MeV, beta(n) is in deg/MeV.

Emittance, Twiss parameters, beam sizes, energy and phases are at the end of the element.

element ngood	plane	emittance (cm-mrad),(deg-MeV)	alpha	beta(u)	rms(u)	max	Design	Average	Reference	Average	
		100%	90%	rms(n)	(cm/mrad), (deg/MeV)	x or y (cm)	x or y (cm)	Phase (deg)	Phase (deg)	Energy (MeV)	Energy (MeV)
1	x-xp	0.30472	0.22549	0.05531	3.05259	0.07252	0.2947	0.6763	0.0000	0.0000	2.0000
	y-yp	0.30766	0.22182	0.05542	-2.64926	0.05899	0.2660	0.6100			
	phi-w	1.13070	0.81924	0.20009	-0.13512	570.366	10.6828	0.0000			

Transport length differs from n*beta*lambda by 0.0000 deg, equivalent to 0.0000 cm.

End of structure 1 Run 1

```
=====
LIN: STRUCTURE 1 0 80.00000000 80.00000000
LIN: TRANSPORT 0.00000000 ; TRANS
LIN: QuadCalc 1 1 1
LIN: Drift 17.30000000 20000.00000000 1 1 ; Comment: D4
LIN: Quad 9.60000000 1.00000000 1 -2650.00000000 0 ; Comment: Q5
LIN: Drift 3.97500000 20000.00000000 1 1 ; Comment: D6
LIN: Cavity 0.0 20000. 1 0.18150000 80.00000000 -40.00000000 ; Comment: G7
LIN: Drift 3.97500000 20000.00000000 1 1 ; Comment: D8
• • •
LIN: Drift 4.87300000 20000.00000000 1 1 ; Comment: D18
LIN: Cavity 0.0 20000. 1 0.18150000 80.00000000 -40.00000000 ; Comment: G19
LIN: Drift 4.87300000 20000.00000000 1 1 ; Comment: D20
LIN: Quad 4.80000000 1.00000000 1 -2650.00000000 0 ; Comment: Q21
LIN: Bore 2000.00000000
LIN: Scheff 0.05000000 0.05000000 9 9 0 1 1
```

PBO Lab Structure "Zero"

Provides initial beam calculation summary

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

PARMILA-2 Outputs - Example B with RF Gaps (con't)

Main Output File: "Parmila.out" (con't)

Beam distribution parameters:

rms(n)	100% ellipse	alfa	beta(u)
(cm-mr)	(cm-mr)	(cm/rad), (deg/MeV)	
1	0.0553	0.3047	3.0526 72.5210
2	0.0554	0.3077	-2.6493 58.9937
3	0.2001	1.1307	-0.1351 570.3655

LIN: PrtBeam 99.00000000

LIN: OUTPUT 2 1 1 1000 1

LIN: begin

Starting beam dynamics... Current = 100.0 mA

Space charge calculations use the Scheff algorithm.

Standard PARMILA-2 initial beam summary

Starting Transport section

Beam simulation summary

Wsync = 2.0000 MeV, 1000 good particles

Beam centroid: x (cm)	xp (mr)	y (cm)	yp (mr)	phi (deg)	W (MeV)
0.00000	0.00000	0.00000	0.00000	0.000	2.000

Phis = 0.0000 deg

Transport finished element 1: 1. Drift

Emittance, Twiss parameters, beam sizes, energy and phases are at the end of the element.

element ngood	plane	emittance (cm-mrad), (deg-MeV)	alpha	beta(u)	rms(u)	max	Design	Average	Reference	Average	
		100%	90%	rms(n)	(cm/mrad), (deg/MeV)	x or y (cm)	x or y (cm)	Phase (deg)	Phase (deg)	Energy (MeV)	
1	1000	x-xp y-yp phi-w	0.34947 0.42452 1.49577	0.23100 0.24132 0.83265	0.05630 0.05799 0.20511	0.50220 -6.56283 0.75988	0.01481 0.20755 611.154	0.1343 0.5104 11.1962	0.3010 1.1460 0.0000	-0.0023 2.0000	2.0000 2.0001

Transport finished element 2: 3. Quadrupole magnet

2	960	x-xp y-yp phi-w	0.35399 0.53637 1.71989	0.23080 0.23958 0.87469	0.05677 0.05750 0.21239	-2.19065 8.72791 1.26147	0.02729 0.14375 691.166	0.1831 0.4230 12.1160	0.4352 0.8385 0.0000	0.0000 0.1122 2.0000	0.1122 2.0000 1.9989
---	-----	-----------------------	-------------------------------	-------------------------------	-------------------------------	--------------------------------	-------------------------------	-----------------------------	----------------------------	----------------------------	----------------------------

Transport finished element 3: 1. Drift

3	960	x-xp y-yp phi-w	0.35532 0.55389 1.77224	0.23025 0.24556 0.87824	0.05667 0.05837 0.21334	-3.18785 6.35442 1.47655	0.04867 0.08226 744.456	0.2443 0.3224 12.6025	0.5923 0.6423 0.0000	0.0000 0.1121 0.1121	0.1121 2.0000 1.9989
---	-----	-----------------------	-------------------------------	-------------------------------	-------------------------------	--------------------------------	-------------------------------	-----------------------------	----------------------------	----------------------------	----------------------------

RF gap Wsync = 2.1390 MeV

Beam centroid: x (cm)	xp (mr)	y (cm)	yp (mr)	phi (deg)	W (MeV)
-0.00154	-0.13805	-0.00071	-0.02453	0.112	2.136

Transport finished element 4: 2. RF Cavity

• • •

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

PARMILA-2 Outputs - Example B with RF Gaps (con't)

Main Output File: "Parmila.out" (con't)

• • •

Transport finished element 17: 1. Drift

17	910	x-xp	0.60290	0.22709	0.05735	4.48549	0.07977	0.2960	0.5930	0.0000	0.3259	2.5561	2.5458
		y-yp	0.81246	0.26269	0.06238	-10.02267	0.21431	0.5059	1.3492				
		phi-w	2.46901	1.01229	0.23985	-1.47309	504.815	11.0036	0.0000				

Transport finished element 18: 3. Quadrupole magnet

18	829	x-xp	0.63402	0.22914	0.05804	1.22366	0.05471	0.2466	0.5064	0.0000	0.0951	2.5561	2.5466
		y-yp	0.35664	0.20040	0.05098	-1.17775	0.22395	0.4675	0.9924				
		phi-w	2.26173	0.98204	0.22782	-1.31140	494.279	10.6115	0.0000				

Transport length differs from n*beta*lambda by -34.0009 deg, equivalent to -1.8462 cm.

Number of particles surviving

Beam dynamics RF phase shift

End of structure 2 Run 1

=====

LIN: end

Beam centroid: x (cm)	xp (mr)	y (cm)	yp (mr)	phi (deg)	W (MeV)
0.00562	-0.36958	0.00727	0.05663	0.157	2.546

Parmila finished at 01-16-2011 14:28:26

PARMILA-2 run summary

To eliminate SFdata output, insert NoSFTableOut before any SFdata table.

Total time 0 min 1.69 sec

Scheff time 0 min 0.05 sec, S/T= 2.96 %

- "Parmila.out" for **PARMILA-2 transport simulations** provides RMS beam data
- Other **text file outputs**: "Quad.out" "Design.out" "Beam.out" "Tr3din.t3d"
 - ⇒ **Use PBO Lab menu: View -> PARMILA-2 -> Auxiliary Files**
- Special (**Lingraf**) plot data written in **binary** format to "Parmila.plt"
 - ⇒ **Use PBO Lab PARMILA-Lingraf Module**
- Final particle distribution data written in **binary** format to "part_dtl.dst"

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

PARMILA-2 Outputs - Example B with RF Gaps (con't)

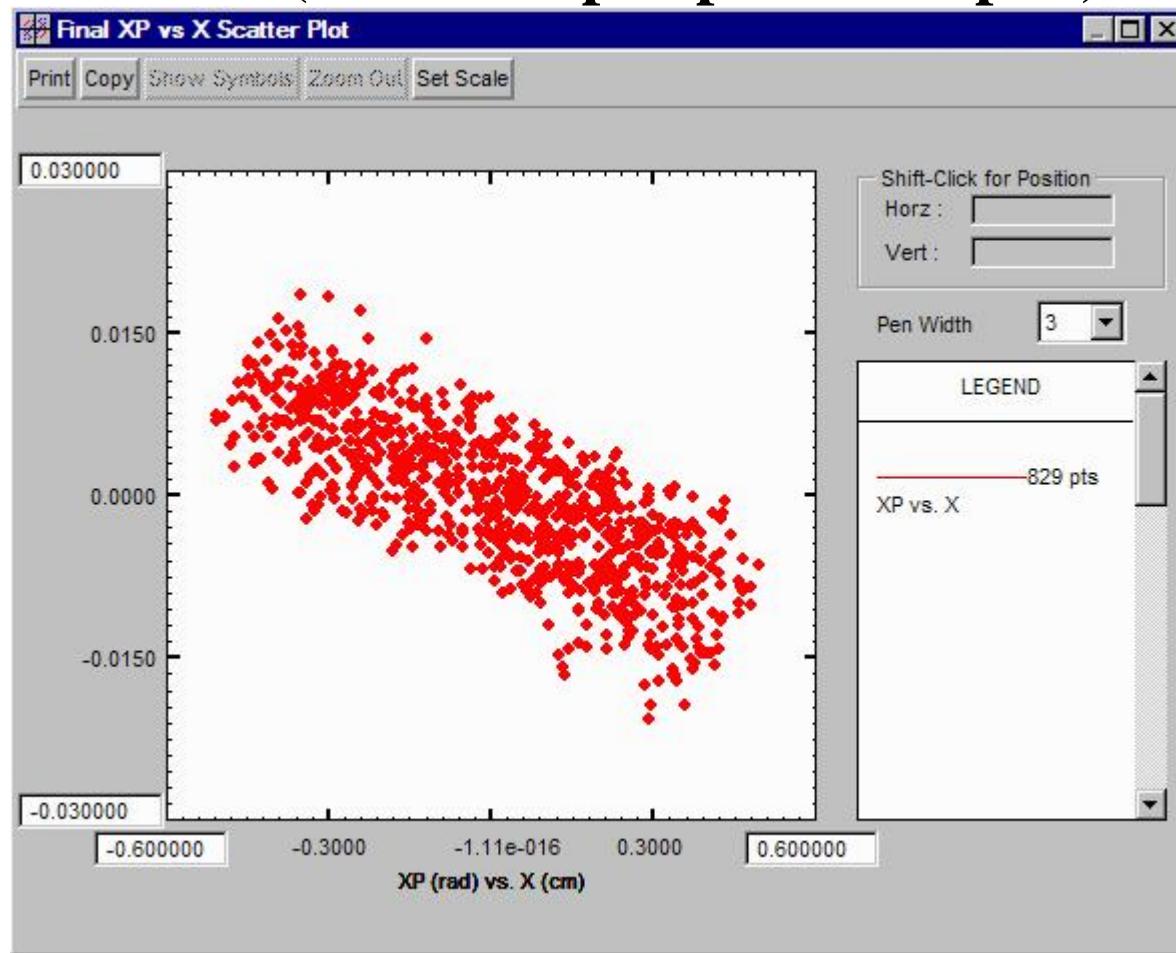
Binary Output File: "part_dtl.dst"

- Program "readdst.exe"
 - Converts *binary* file "part_dtl.dst" to *formatted text* file "part_dtl.TXT"
- From PBO Lab use: Commands->PARMILA Plots->Write and Run 'input.temp'
 - Use the Windows file open dialog to select "part_dtl.dst"
 - Formatted text file "part_dtl.TXT" will be opened at completion
 - Three default phase space plots will be displayed for x-x', y-y', W-phi
 - Other plots of the data can be selected by the user (histogram, scatter)
- First few lines from "part_dtl.TXT" illustrated below

```
Parmila data from "Example B.pbol run on 01/16/11 14:28:23" run at time 01-16-2011 14:28:25
Structure number      =          2
Cell or element number =          0
Design particle energy = 2.000000    MeV
Number of particles   =        829
Beam current          = 82.90000   (frequency scaled equivalent for every rf bucket filled)
RF Frequency          = 80.00000   MHz
Bunch Freq             = 80.00000   MHz
Chopper fraction       = 1.000000
The input file particle coordinates were written in double precision.
      x(cm)           xpr(=dx/ds)      y(cm)           ypr(=dy/ds)      phi(radian)      W(MeV)
-0.2740483529     8.5920352496E-03  0.5239293919   2.4765092954E-03 -1.9268399094E-02  2.529683666
-0.1522036535     2.6527460134E-03  -7.0805802018E-02  1.0753200838E-03  -0.2567370817   2.453311165
-0.1422078128     2.5051232246E-03  -0.3606680071   1.3703372955E-03   0.1262587515   2.577495844
  0.2895415372    -8.4868751524E-03  -0.3359137013   -7.3630744864E-04  -0.2435264125   2.489523835
-0.3386279447     7.8457234313E-03  -0.3981546281   6.2457500245E-04   7.7077898501E-02  2.581053081
```

3. Introduction to PARMILA - For Simulation of Transfer Lines (continued)

Typical Scatter Plot (PBO Lab postprocessor plot) PARMILA-2



- PBO-Lab PARMILA Plots has same basic features as plots for TURTLE
- Apertures (of quadrupoles) responsible for beam loss in this example

4. Space Charge Modeling in PARMILA

Space Charge in PARMILA-2

- PARMILA-2 has two basic space charge calculation algorithms
 - SCHEFF is a 2-D method that is the **default**
 - PICNIC is a 3-D that is an **option**
 - To invoke **either** algorithm the "input.LIN" file **must have a SCHEFF line**
- SCHEFF ("Space Charge EFFects")
 - Particles first transformed to beam rest frame (**from lab frame**)
 - r-z mesh centered on the beam rest frame (rectangular mesh of r & z)
 - Data on the SCHEFF line is used to define mesh parameter
 - At specific steps for each element (# depends upon element):
 - Charge is allocated to points on the mesh to determine effective field
 - Space charge impulse force is applied to particles using that field
 - Particles then transformed back to lab frame
- PICNIC (written by Nicolas Pichoff) computes the fields differently
 - Computes RMS beam size for all 3 dimensions (x_{rms} , y_{rms} , z_{rms})
 - Uses a 3-D mesh given by $3.5 \times$ RMS sizes ($3.5x_{rms}$, $3.5y_{rms}$, $3.5z_{rms}$)
 - For particle outside the core, the force for a Gaussian fit to core used
- Both use the number of mesh points parameters defined by Scheff line

4. Space Charge Modeling in PARMILA

Space Charge in PARMILA-2

- SCHEFF line has form

Scheff	ΔR_{sc}	ΔZ_{sc}	N_R	N_Z	N_{bunch}	$N_{\beta\lambda}$	Remesh
e.g. Scheff	0.05000	0.05000	9	9	0	1	1

- The mesh intervals will be dynamically recalculated with SCHEFF method
 - According the Remesh value on the Scheff line
 - The dynamical mesh intervals of SCHEFF can be overridden using a line
MeshFact 2.0 3.0 \Rightarrow mesh intervals changed to: $2.0 \times \Delta R_{sc}$ $2.0 \times \Delta Z_{sc}$
- SCHEFF includes space charge from adjacent bunches according to N_{bunch}
 - According the Remesh value on the Scheff line
 - Adjacent bunches are take to be $N_{\beta\lambda} \times \beta\lambda$ from bunch being computed
- PICNIC is invoked with line:
Use3DPicnic
- The number of mesh points for PICNIC are given by
 $N_x = 2 \times N_R$ $N_y = 2 \times N_R$ $N_z = 2 \times N_z$
- SCHEFF can be re-invoked using line
UseSCHEFF
- Note that loss particles can impact the dynamic meshing for either method

5. Using PARMILA & TURTLE to Study Some Beamlines
⇒ **Use the Simulation Lab computers in the classroom**

- Compare PARMILA & TRACE 3-D at Low (~0) Current
- Explore the use of Lingraf to display data
- Compare PARMILA & TURTLE at Low (~0) Current
- Use PARMILA to Study High Current Transport
- Compare SCHEFF & PICNIC Space Charge Calculations
- Compare PARMILA & TRACE 3-D at High Current