

PyHEADTAIL examples

All our thanks to
D. Amorim, Kevin Li and Michael Schenck
CERN BE/ABP-HSC

and

Martial Fol (AZIMUTEC), Marie Gauthier and Coline Morin (ESI)

Reference: <http://kli.web.cern.ch/kli/>

Agenda

- Goals
- Plan
- Introduction to PyHEADTAIL
- Structure of ipython files
- Tracking examples
- Setting up the environment
 - Virtual box
 - Ubuntu
 - Anaconda
 - PyHEADTAIL

Goals

- Run macroparticle simulations with a state-of-the-art open source tracking code (PyHEADTAIL)
- Simulate several case-studies related to the course
- Play with beam parameters and observe the impact on the longitudinal beam dynamics

Plan

- 1h15 towards the end of the course in the computer room
- A virtual box with examples is prepared in the computer room
- Following popular demand, it is also possible to set up your own simulation environment on your own PC.
- The detailed procedure and examples are on the Indico site.
- Note: we should expect incompatibilities linked to open source codes!

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Introduction to PyHEADTAIL

- Open source macroparticle tracking code developed at CERN:
- Download link: <https://github.com/PyCOMPLETE/PyHEADTAIL>
- Primary use: tracking simulation of collective effects in synchrotron accelerators
 - Transverse and longitudinal beam dynamics (with feedback)
 - Electron/ion cloud
 - Impedances
 - Space charge
- Reference:
[Introduction to PyHEADTAIL: USPAS course](#) by Kevin Li et al (2015)
- Not the only code of his kind!
 - **BLOND**: longitudinal dynamics simulation code <https://blond.web.cern.ch/>
 - **HEADTAIL**: the father of PyHEADTAIL! G. Rumolo et al (reference)
 - **elegant**: 6D tracking code developed at Argonne National Lab ([link](#))
 - **mbtrack** and **sbtrack**: R. Nagaoka et al “Studies of Collective Effects in SOLEIL and DIAMOND Using the Multiparticle Tracking Codes sbtrack and mbtrack”, PAC09, Vancouver, May 2009.
 - ORBIT (<http://web.ornl.gov/~jzh/JHolmes/ORBIT.html>), pyORBIT (<http://sourceforge.net/projects/py-orbit/>)
 - And so many others!

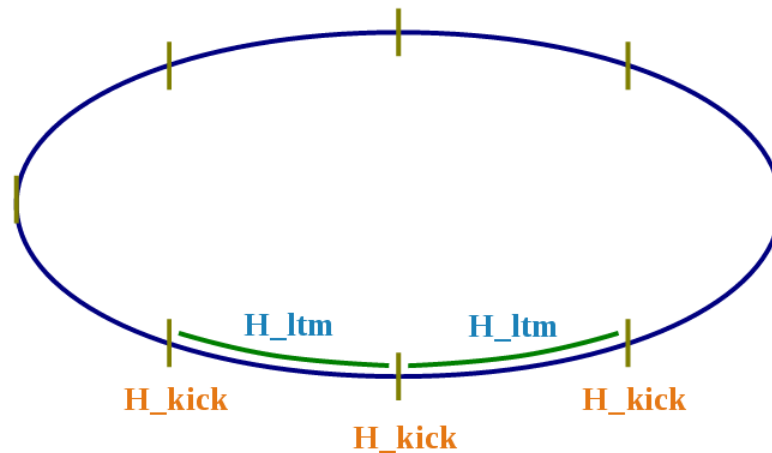
introduction to PyHEADTAIL

Courtesy Kevin Li
USPAS 2015

How does PyHEADTAIL work?



- PyHEADTAIL is a macroparticle tracking code designed specifically to simulate collective effects in circular accelerators



- H_ltm: linear transfer map
 - Chromaticity
 - Amplitude detuning
 - ...
- H_kick: collective interaction
 - Wakefields
 - Electron cloud
 - Feedback
 - Space-charge
 - ...



introduction to PyHEADTAIL



A real world example

Courtesy
Kevin Li
USPAS 2015



- Load Python packages and modules

```
1 from __future__ import division
2 import cProfile, itertools, sys, time, timeit
3
4 from scipy.constants import c, e, m_p
5
6 from cobra_functions import stats, random
7 from beams.beams import *
8 from monitors.monitors import *
9 from spacecharge.spacecharge import *
10 from trackers.transverse_tracker import *
11 from trackers.longitudinal_tracker import *
12
```

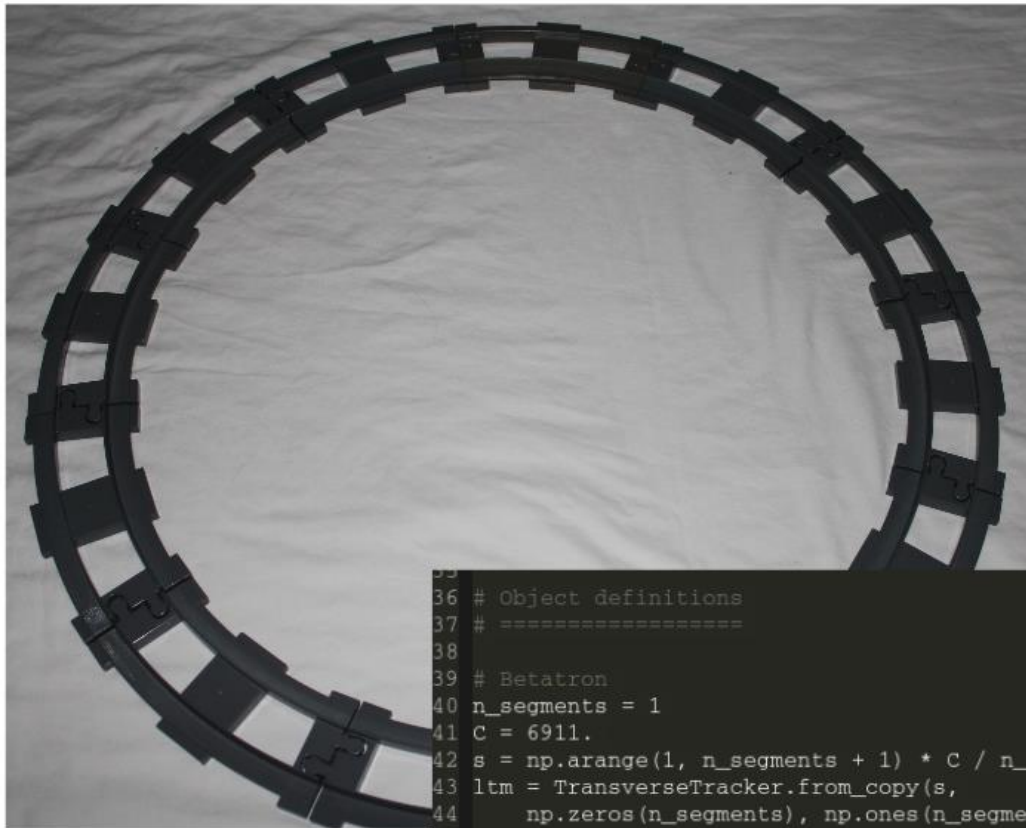


introduction to PyHEADTAIL



A real world example

Courtesy
Kevin Li
USPAS 2015



- Load Python packages and modules
- Build linear periodic transfer maps

```
35  
36 # Object definitions  
37 # =====  
38  
39 # Betatron  
40 n_segments = 1  
41 C = 6911.  
42 s = np.arange(1, n_segments + 1) * C / n_segments  
43 ltm = TransverseTracker.from_copy(s,  
44     np.zeros(n_segments), np.ones(n_segments) * beta_x, np.zeros(n_segments),  
45     np.zeros(n_segments), np.ones(n_segments) * beta_y, np.zeros(n_segments),  
46     Qx, 0, 0, Qy, 0, 0)  
47
```

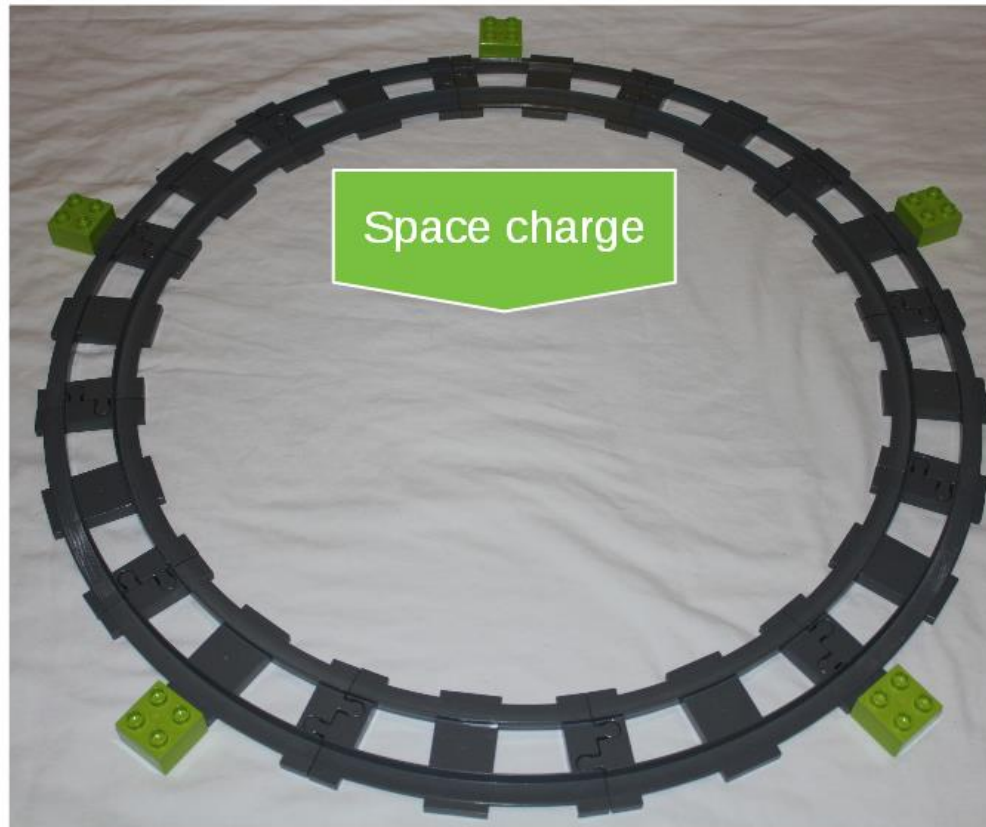


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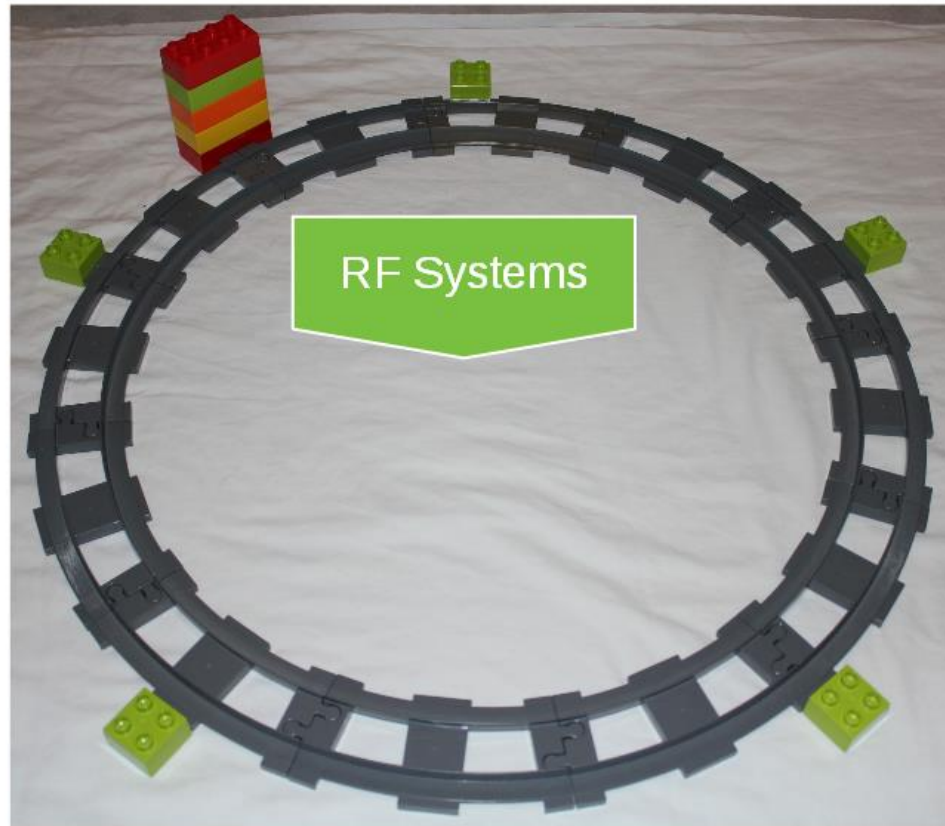
- Load Python packages and modules
- Build linear periodic transfer maps
- Add (collective) kick elements

introduction to PyHEADTAIL



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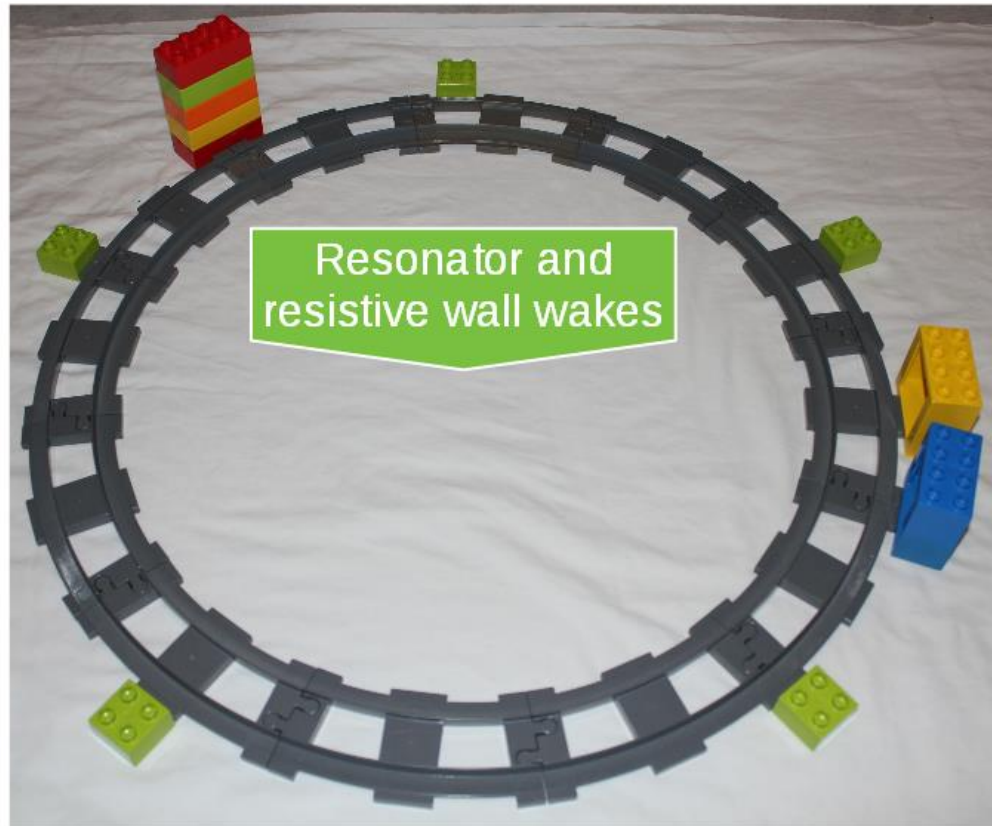
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introduction to PyHEADTAIL



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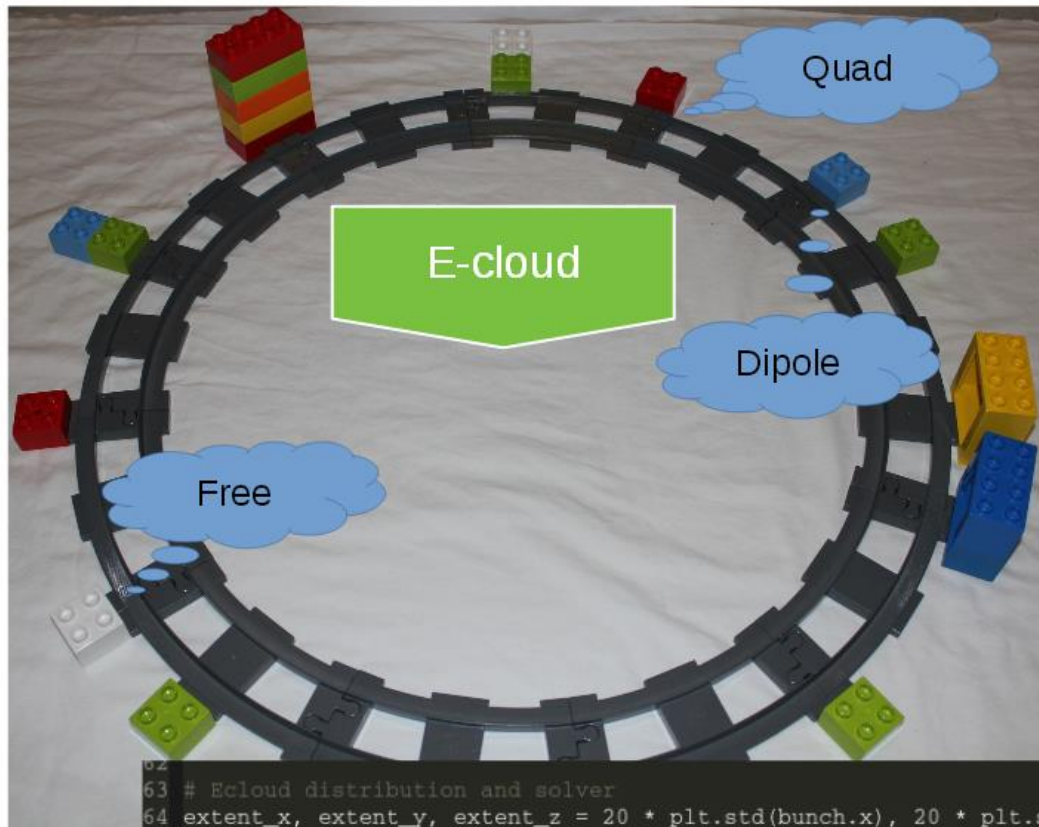
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introduction to PyHEADTAIL



A real world example

Courtesy
Kevin Li
USPAS 2015



- Load Python packages and modules
- Build linear periodic transfer maps
- Add (collective) kick elements

```
62  
63 # Ecloud distribution and solver  
64 extent_x, extent_y, extent_z = 20 * plt.std(bunch.x), 20 * plt.std(bunch.y), C / n_segments  
65 cloud = Cloud(100000, lcell, extent_x, extent_y, extent_z)  
66 ecloud = SpaceCharge(cloud, 'cloud', extent_x, extent_y, 128, 128, slices)  
67
```

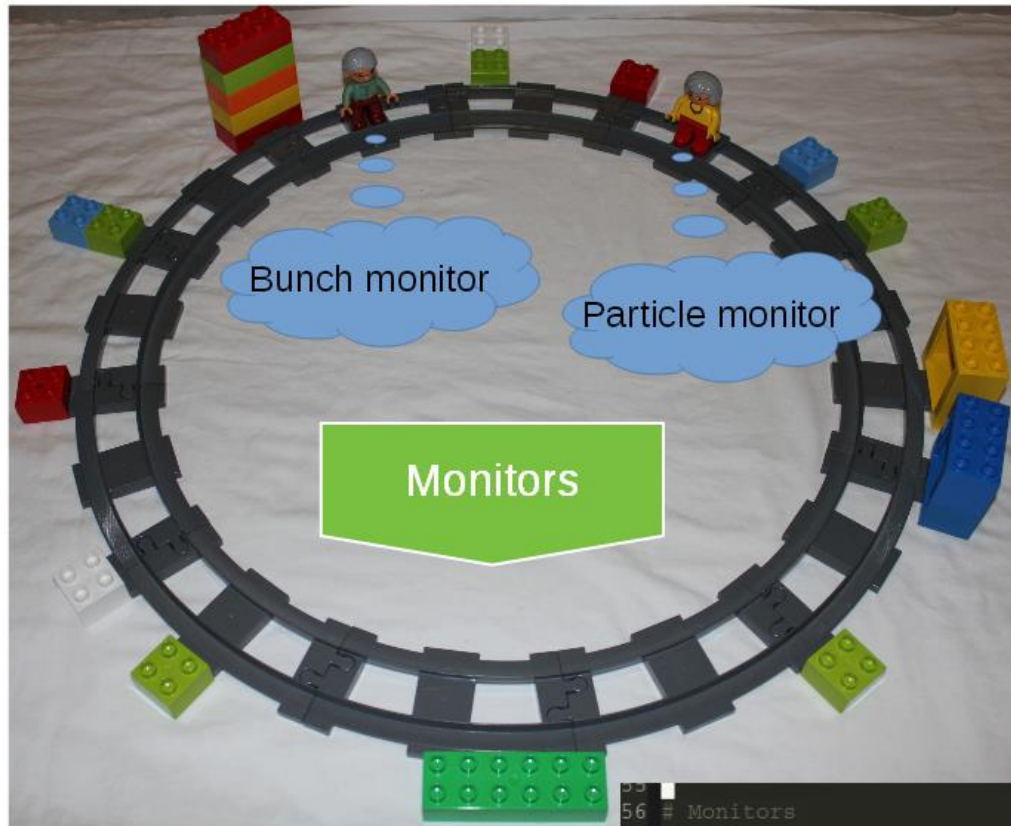


introduction to PyHEADTAIL



A real world example

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USPAS 2015



- Load Python packages and modules
- Build linear periodic transfer maps
- Add (collective) kick elements

```
56 # Monitors
57 bunchmonitor = BunchMonitor('bunch', n_turns, slices)
58 particlemonitor = ParticleMonitor('particles', n_turns, slices)
```

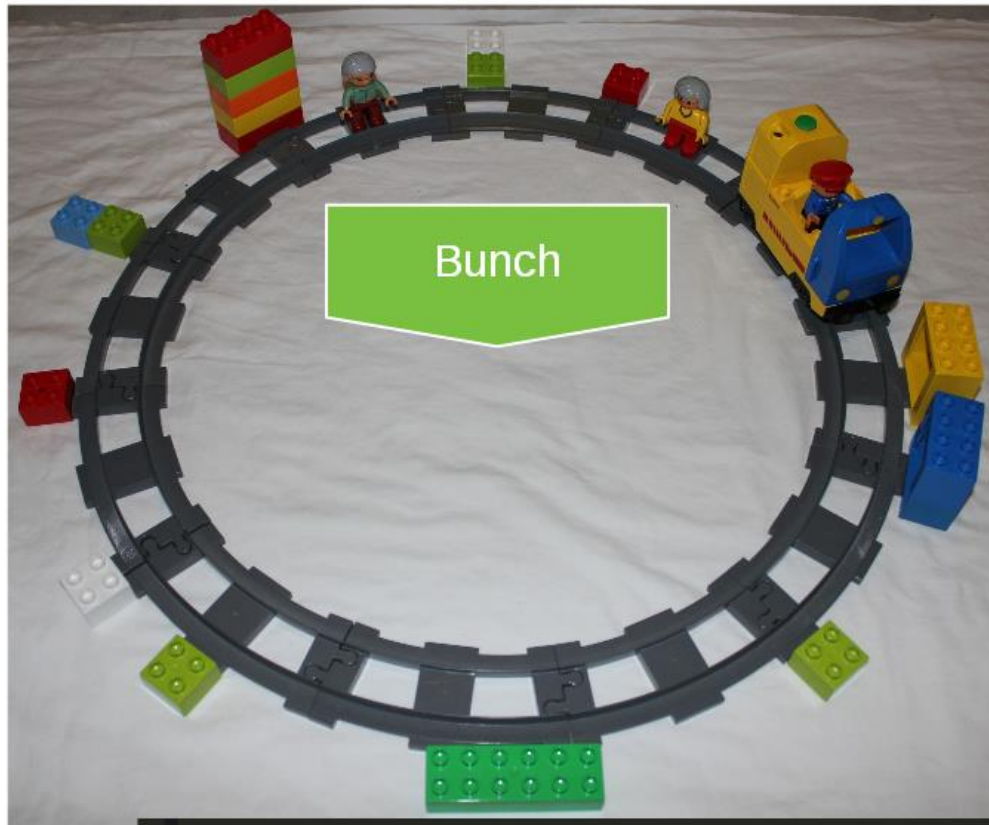


introduction to PyHEADTAIL



A real world example

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Kevin Li
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- Load Python packages and modules
- Build linear periodic transfer maps
- Add (collective) kick elements
- Place beam

```
60 # Bunch  
61 bunch = Bunch(500000, e, gamma, 1.15e11, m_p, 0, beta_x, epsn_x, 0, beta_y, epsn_y, beta_z, sigma_z)  
62
```

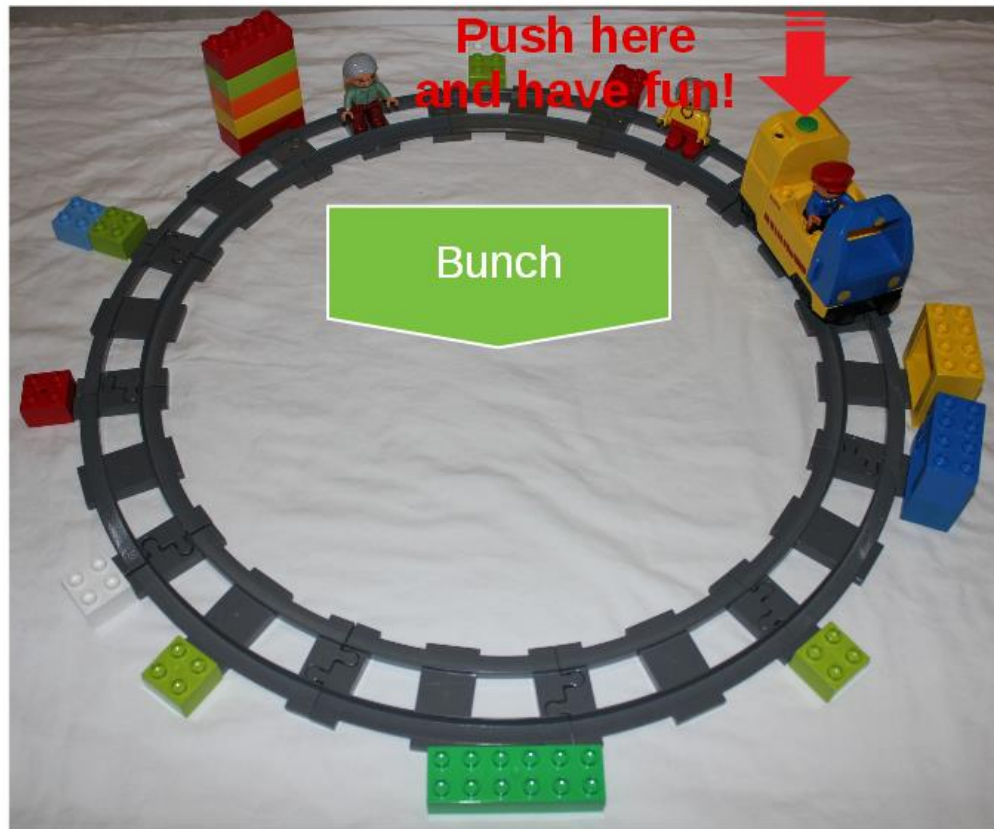


introduction to PyHEADTAIL



A real world example

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USPAS 2015



- Load Python packages and modules
- Build linear periodic transfer maps
- Add (collective) kick elements
- Place beam

```
86 for i in range(n_turns):  
87     for m in map_:  
88         m.track(bunch)  
89
```



In our case: only RF systems

JUAS

Tracking

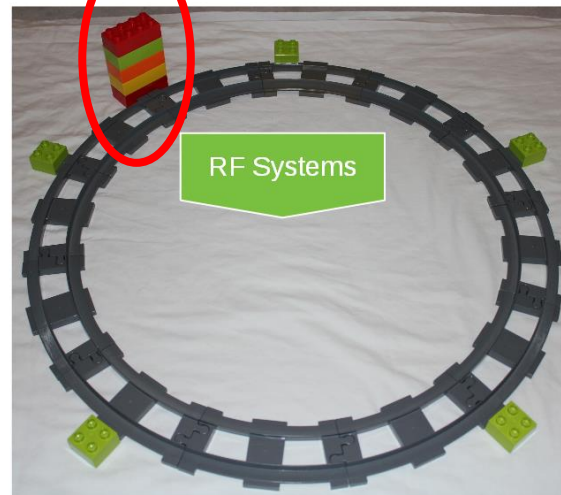
- ◆ The motion of the particles can be tracked turn by turn using the recurrence relation (between turn n and turn $n+1$)

$$\Delta E_{n+1} = \Delta E_n + e \hat{V}_{RF} [\sin \phi_n - \sin \phi_s]$$

$$\phi_{n+1} = \phi_n - \frac{2 \pi h \eta}{\beta_s^2 E_s} \Delta E_{n+1}$$

JUAS - Jan 2016 - E.Métral

A real world example



- Load Python packages and modules
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Structure of the ipython file

Import needed libraries

```
In [ ]: import sys
sys.path.append("../..")
import numpy as np
from scipy.constants import m_p, c, e
from PyHEADTAIL.particles.particles import Particles
import PyHEADTAIL.particles.generators as generators
from PyHEADTAIL.trackers.transverse_tracking import TransverseMap
from PyHEADTAIL.trackers.simple_long_tracking import RFSystems, LinearMap
import PyHEADTAIL.cobra_functions.stats as st
import matplotlib.pyplot as plt
```

Set beam and machine parameters

```
In [ ]: # general simulation parameters
n_particles = 1000
n_segments = 1

# machine parameters
circumference = 1000*np.pi
inj_alpha_x = 0e-1.2
inj_alpha_y = 0e15
inj_beta_x = 16.*5.9 # in [m]
inj_beta_y = 16.*5.7 # in [m]
Qx = 6.25
Qy = 6.25
gamma_tr = 6.1
alpha_c_array = [gamma_tr**2]
V_rf = 20e3 # in [V]
harmonic = 5
phi_offset = 0 # measured from aligned focussing phase (0 or pi)
# pipe radius = 5e-2
Bdot=0 # in T/s
bending_radius=70 # in m

# beam parameters
Ekin = 1.4e9 # in [eV]
intensity = 8e12
epsn_x = 5e-6 # in [m*rad]
epsn_y = 5e-6 # in [m*rad]
#epsn_z = 1. # 4pi*sig_z*sig_dp (*p0/e) in [eVs]

# calculations
gamma = 1 + e * Ekin / (m_p * c**2)
beta = np.sqrt(1 - gamma**(-2))
print('beta: ' + str(beta))
eta = alpha_c_array[0] - gamma**(-2)
print('eta: ' + str(eta))
if eta < 0:
    phi_offset = np.pi - phi_offset
Etot = gamma * m_p * c**2 / e
p0 = np.sqrt(gamma**2 - 1) * m_p * c
Qs = np.sqrt(np.abs(eta) * V_rf / (2 * np.pi * beta**2 * Etot))
print('Qs: ' + str(Qs))
beta_z = np.abs(eta) * circumference / (2 * np.pi * Qs)
print('beta_z: ' + str(beta_z))
turn_period = circumference / (beta * c)
p_increment_0 = e*bending_radius*Bdot*turn_period
sigma_z_0 = 230e-9/4*beta*c/10

# BETATRON
# Loop on number of segments and create the TransverseSegmentMap
# For each segment:
s = np.arange(0, n_segments + 1) * circumference / n_segments
alpha_x = inj_alpha_x * np.ones(n_segments)
beta_x = inj_beta_x * np.ones(n_segments)
D_x = np.zeros(n_segments)
alpha_y = inj_alpha_y * np.ones(n_segments)
beta_y = inj_beta_y * np.ones(n_segments)
D_y = np.zeros(n_segments)

# Define RF systems
rfsystems = RFSystems(circumference, [harmonic], [V_rf], [phi_offset],
                    alpha_c_array, gamma, p_increment=p_increment_0)
# Generate the particle distribution
bunch = generators.ParticleGenerator(macroparticlenumber=n_particles,
                                   intensity=intensity, charge=e, mass=m_p,
                                   circumference=circumference, gamma=gamma,
                                   distribution_x=generators.gaussian2D(epsn_x, alpha_x=inj_alpha_x),
                                   distribution_y=generators.gaussian2D(epsn_y, alpha_y=inj_alpha_y),
                                   distribution_z=generators.RF_bucket_distribution(rfsystems)
                                   ).generate()
```

Define RF system

Generate a matched distribution corresponding to these parameters

Loop over number of turns

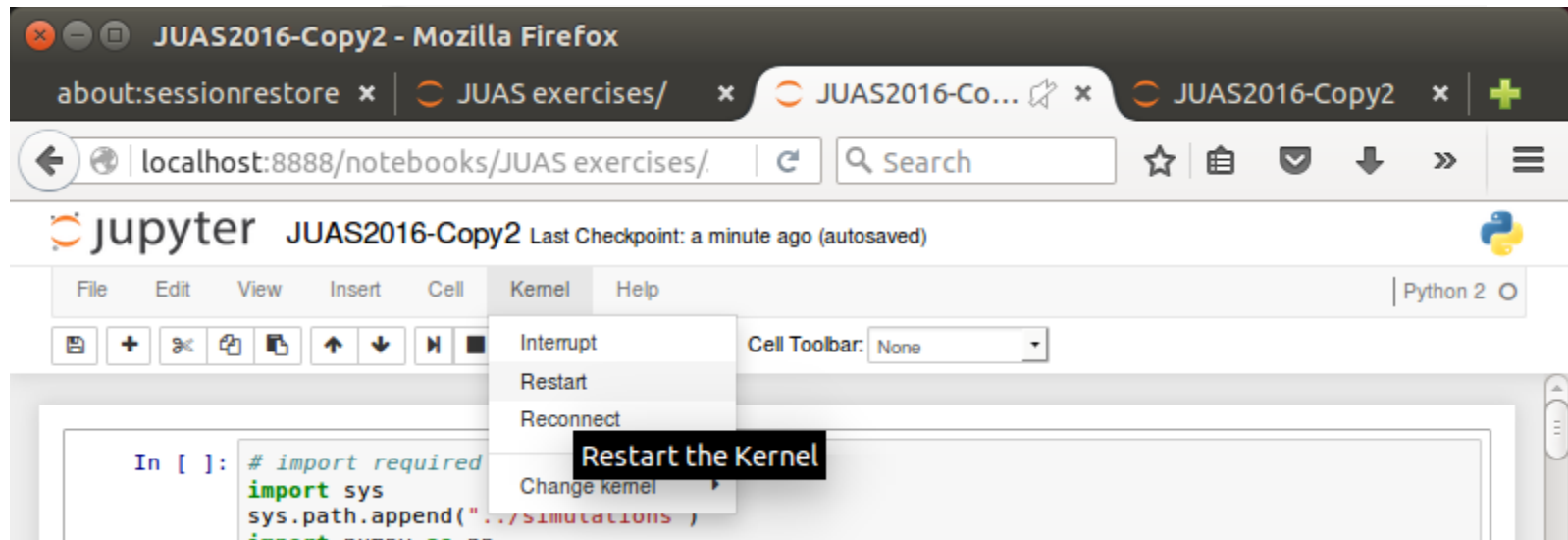
```
In [ ]: # plot phase space
plt.close()
plt.ion()
fig = plt.figure(1)
for i in np.arange(0, 50, 1):
    # track the particles
    rfsystems.track(bunch)
    if i%10 == 0:
        # monitor the particles
        bucket = rfsystems.get_bucket(gamma=bunch.gamma)
        # plot the RF bucket envelope
        z = np.linspace("bucket.interval", num=100)
        dp = np.linspace(-0.005, 0.005, num=100)
        Z2, DPP = np.meshgrid(z, dp)
        HH = bucket.hamiltonian(Z2, DPP)
        plt.contour(Z2, DPP, HH, levels=[0], colors='magenta')

    # plot the particles in phase space
    plt.plot(bunch.z, bunch.dp, 'o')
    plt.xlabel('z in m')
    plt.ylabel('Delta p/p')
    plt.show()
    plt.pause(0.1)
    plt.cla()
plt.ioff()
```

track
monitor
plot

ipython cheat sheet

- To restart the kernel at any point: “Kernel” → “Restart”



- To only interrupt the kernel: “Kernel” → “Interrupt”
- Shift-enter: execute current cell and move to next cell
- Ctrl-enter: execute current cell and stay on current cell

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Tracking example: injection energy

- ◆ nTOF bunch in the CERN PS (near transition)

Average machine radius: R [m]	100
Bending dipole radius: ρ [m]	70
\dot{B} [T/s]	0
\hat{V}_{RF} [kV]	200
h	8
α_p	0.027
Longitudinal (total) emittance: ϵ_L [eVs]	2
Number of protons/bunch: N_b [1E10 p/b]	800
Norm. rms. transverse emittance: $\epsilon_{x,y}^*$ [μm]	5
Trans. average betatron function: $\beta_{x,y}$ [m]	16
Beam pipe [cm \times cm]	3.5 \times 7
Trans. tunes: $Q_{x,y}$	6.25

20 kV at injection

$\Rightarrow \gamma_t \approx 6.1$

Tracking example: injection energy

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Beam pipe [cm \times cm]	3.5 \times 7
Trans. tunes: $Q_{x,y}$	6.25

$$E_{kin} = 1.4e9 \text{ \# in [eV]}$$

$$\text{circumference} = 100 * 2 * \pi$$

$$\text{bending_radius} = 70 \text{ \# in m}$$

$$\text{Bdot} = 0 \text{ \# in T/s}$$

20 kV at injection

$$V_{rf} = 20e3 \text{ \# in [V]}$$

$\Rightarrow \gamma_t \approx 6.1$

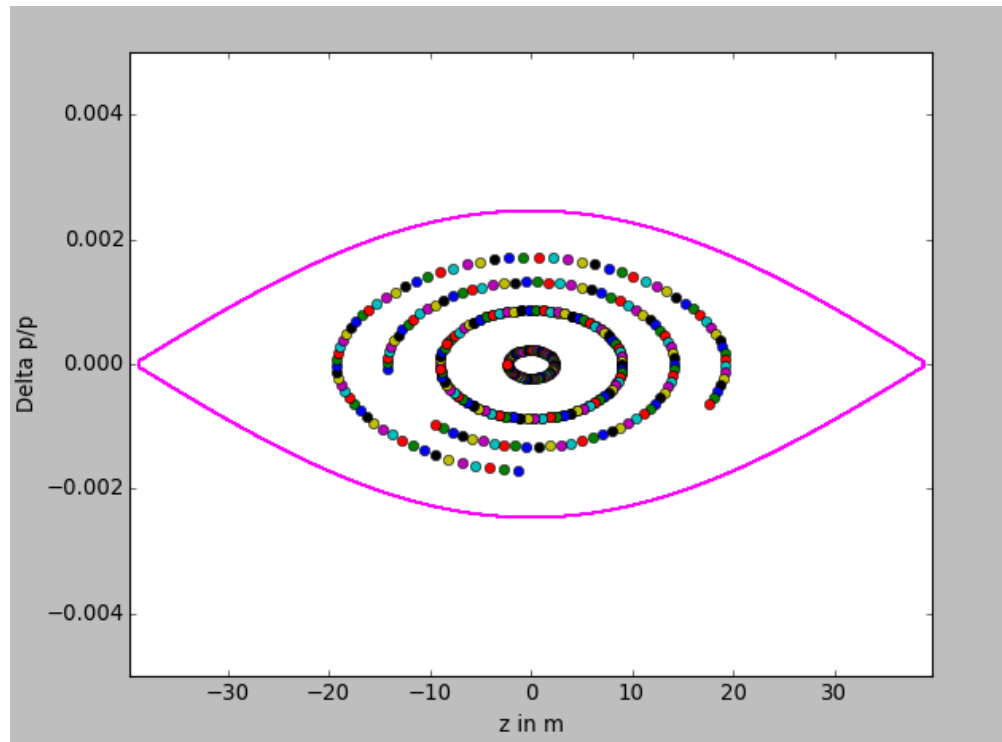
$$\text{harmonic} = 8$$

$$\text{gamma_tr} = 6.1$$

$$\text{sigma_z_0} = 230e-9 / 4 * \text{beta} * c$$

Let's run that example

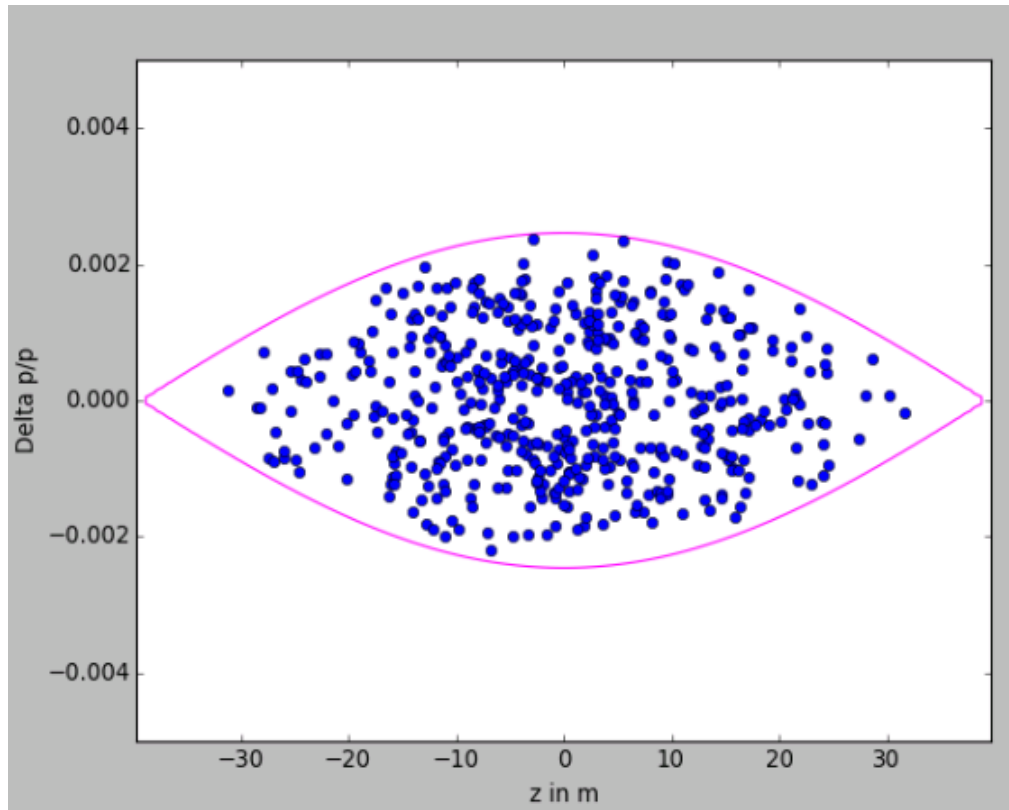
- Use a small number of particles ($n_{\text{particles}} \sim 10$) to observe the synchrotron motion at large and small amplitudes



→ Can you estimate the synchrotron tune Q_s ? Is it consistent with the course?

Synchrotron motion

- Use a larger number of particles (~ 500) and reset the graph every time (“plt.cla()” uncommented)



- Are the particles rotating in the correct direction?
- Plot the phase space as in the course (in ϕ [degrees], ΔE [MeV])
- Is the maximum energy consistent with the course?

Top energy

- What parameters should be changed for top energy (still no acceleration and $h=8$)?
- What changes in the plot compared to injection energy?

Agenda

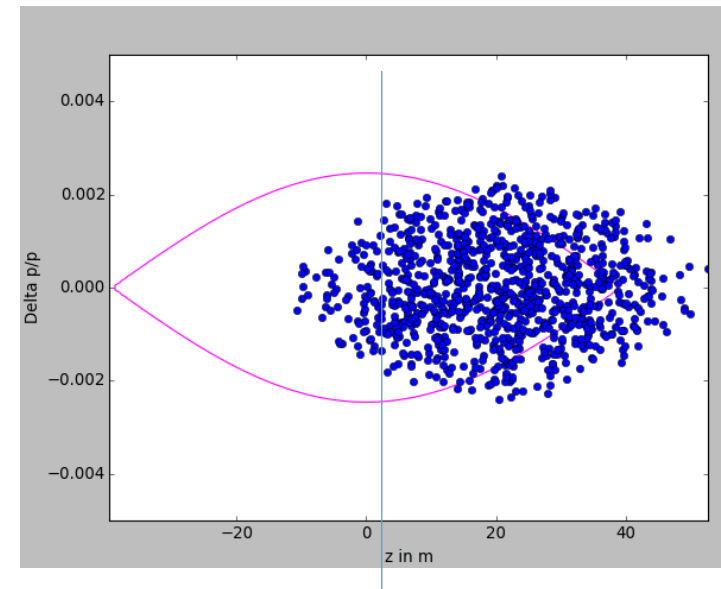
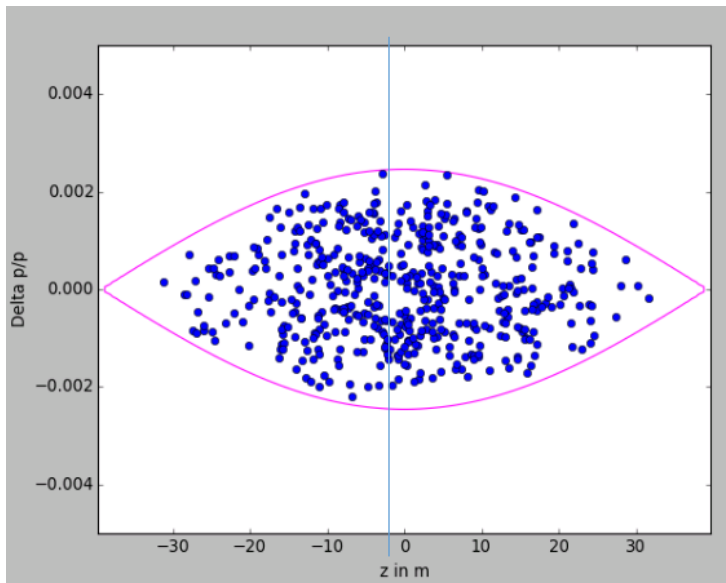
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 - Impact of voltage mismatch
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Impact of phase mismatch

- Generate ~ 1000 particles
- Add e.g. 5 or 20 m to all particles:
- Track particles for ~ 5000 turns
- Plot the distribution every 100 turns:

```
bunch.z = bunch.z+5
```

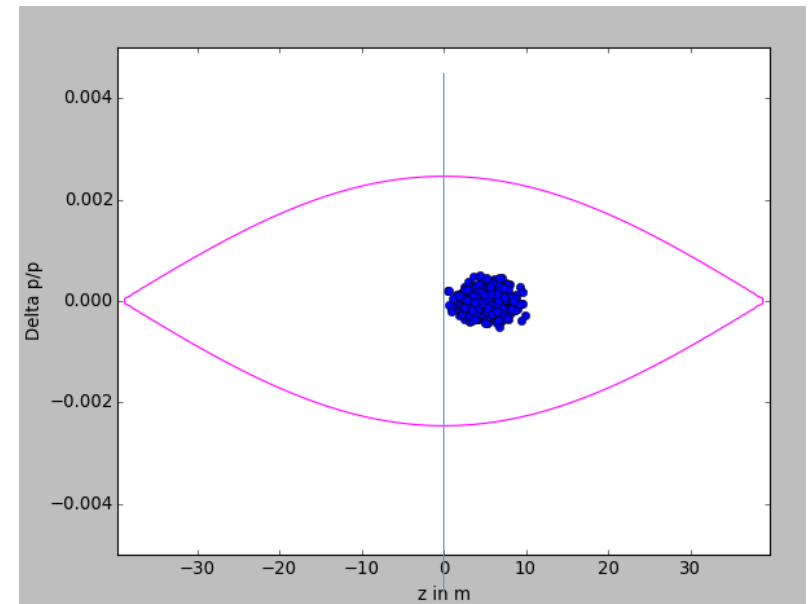
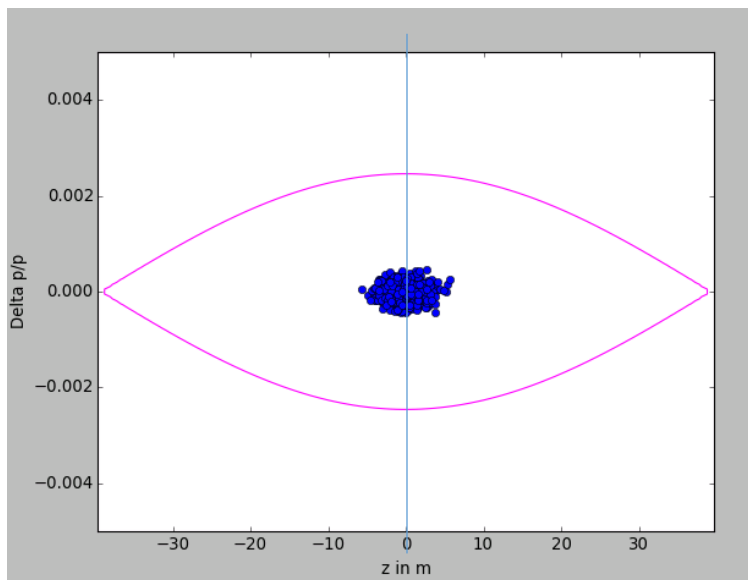
```
if i%100 == 0:
```



→ What will happen?

Impact of phase mismatch

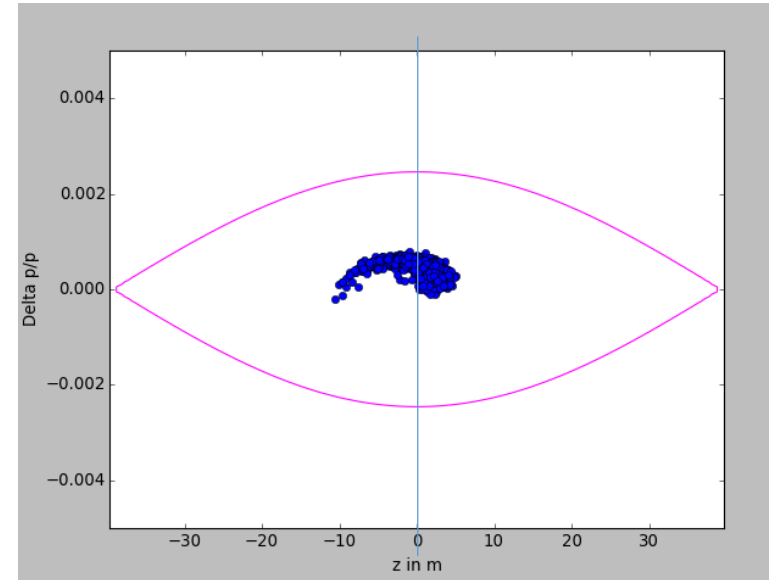
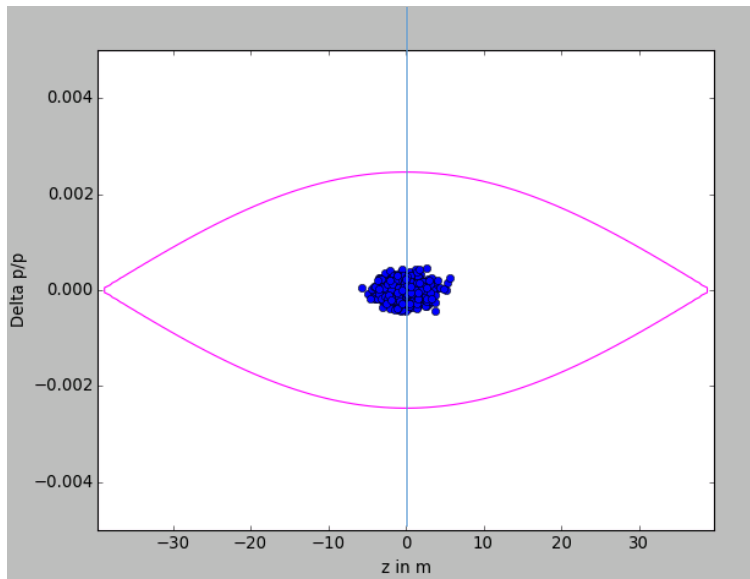
- Use a smaller bunch length (e.g. divide by 10)
- Generate ~ 1000 particles
- Add e.g. 5m to all particles: `bunch.z = bunch.z+5`
- Track particles for ~ 5000 turns
- Plot the distribution every 100 turns: `if i%100 == 0:`



→ What will happen?

Impact of phase mismatch

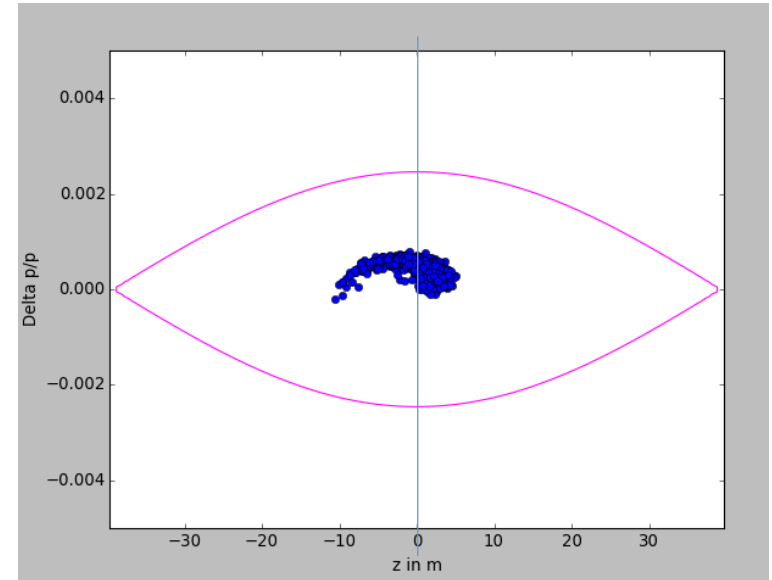
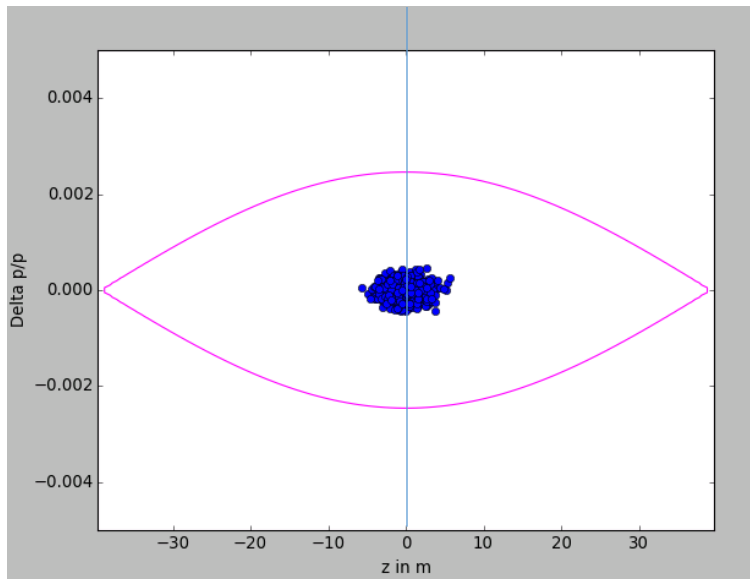
- Use a smaller bunch length (e.g. divide by 10)
- Use ~1000 particles
- Add e.g. 5m to all particles: `bunch.z = bunch.z+5`
- Track particles for ~5000 turns
- Plot the distribution every 100 turns: `if i%100 == 0:`



→ Filamentation and eventually?

Impact of phase mismatch

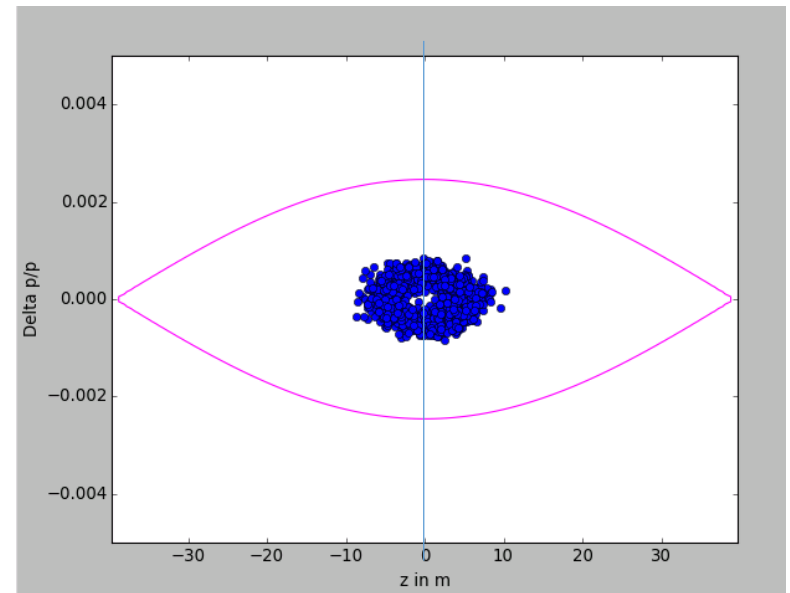
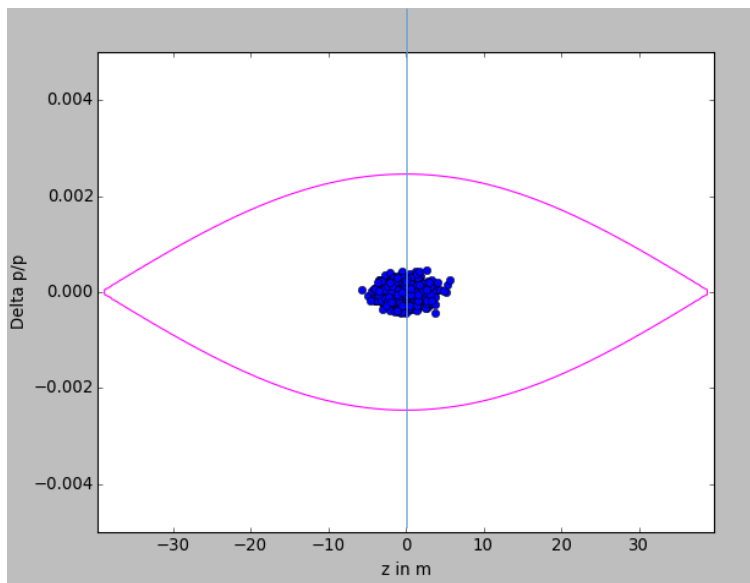
- Use a smaller bunch length (e.g. divide by 10)
- Use ~ 1000 particles
- Add e.g. 5m to all particles:
- Track particles for ~ 500000 turns
- Plot the distribution every 1000 turns



→ Filamentation and eventually?

Impact of phase mismatch

- Use a smaller bunch length (e.g. divide by 10)
- Use ~ 1000 particles
- Add e.g. 5m to all particles:
- Track particles for ~ 500000 turns
- Plot the distribution every 100 turns:



- Eventually emittance growth
- Try also with larger mismatch (20 m)

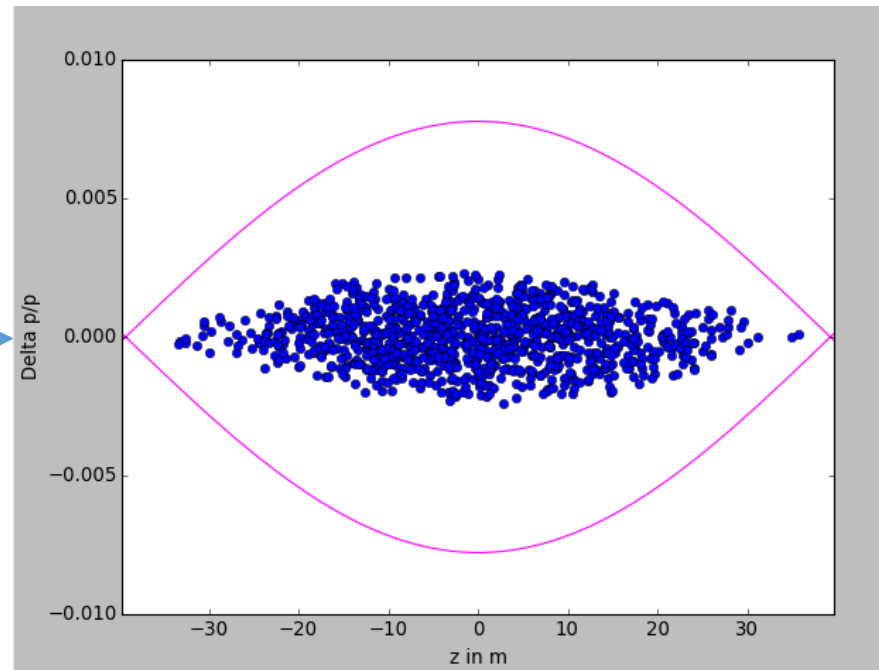
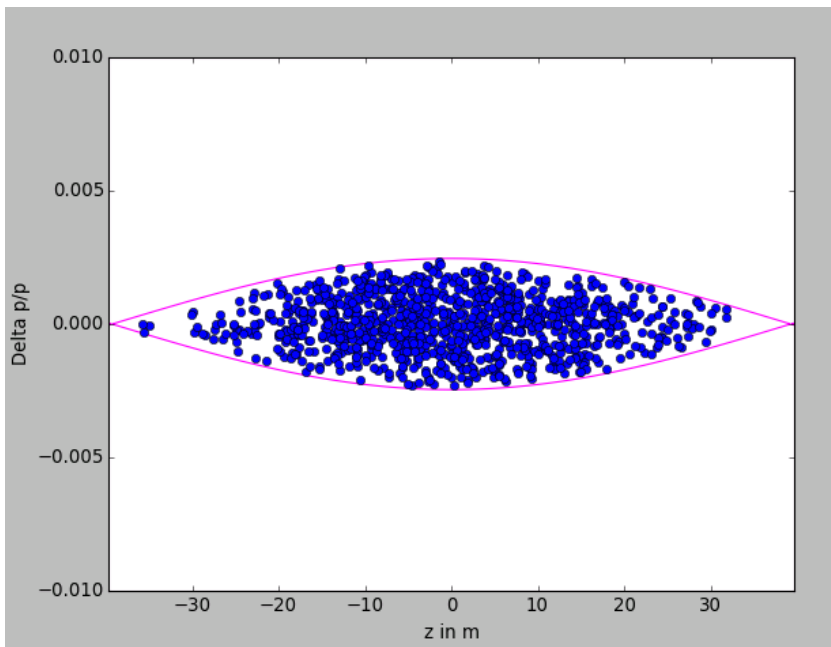
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Impact of voltage mismatch (too high)

- Use the **nominal** bunch length
- track 1000 particles for **500 turns (plot every 10 turns)**
- Change V_{RF} to **$V_{RF} * 10$** after the bunch is matched to V_{RF}

```
# Define RF systems  
rfsystems = RFSystems(circumference, [harmonic], [V_rf*10], [phi_offset],  
                      alpha_c_array, gamma, p_increment=p_increment_0)
```

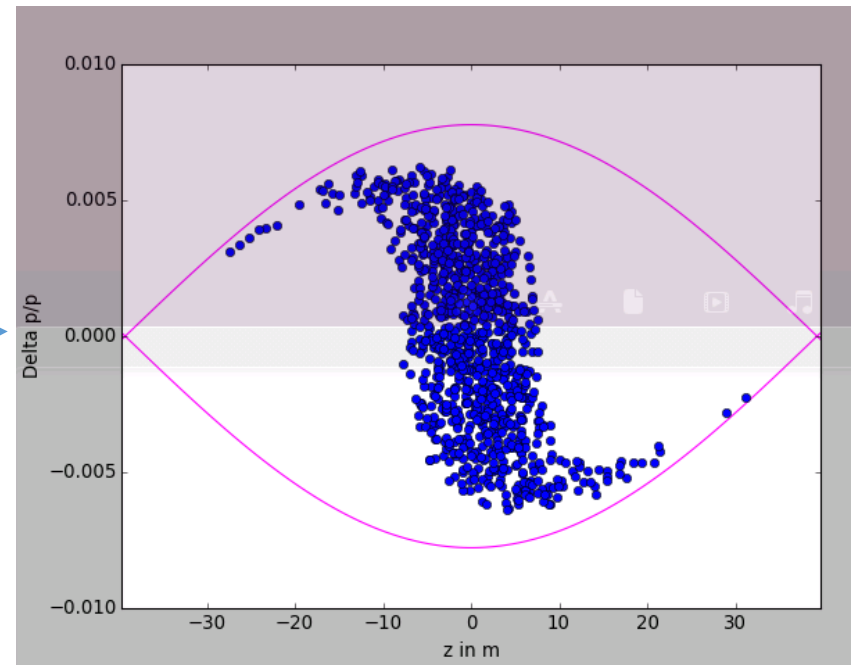
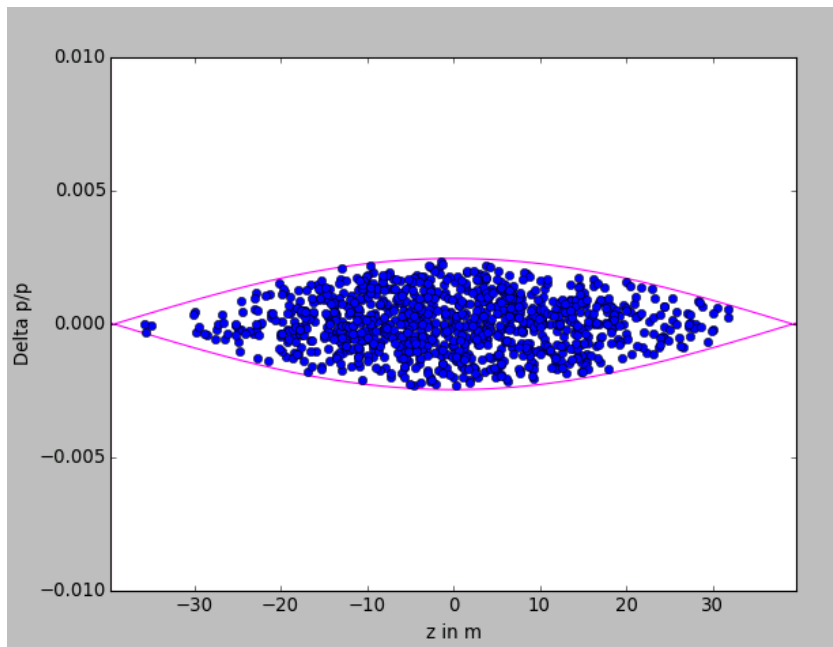


→ What will happen?

Impact of voltage mismatch (too high)

- Use the **nominal** bunch length
- track 1000 particles for **500 turns (plot every 10 turns)**
- Change V_{RF} to **$V_{RF} * 10$** after the bunch is matched to V_{RF}

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rfsystems = RFSystems(circumference, [harmonic], [V_rf*10], [phi_offset],  
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```



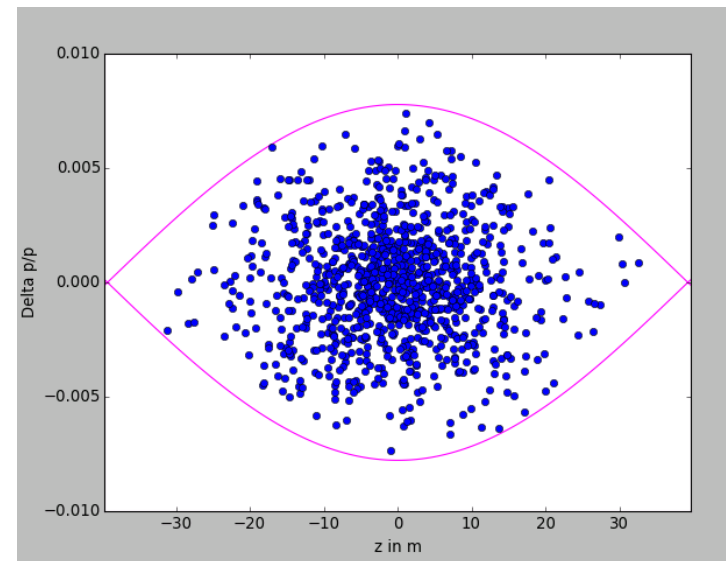
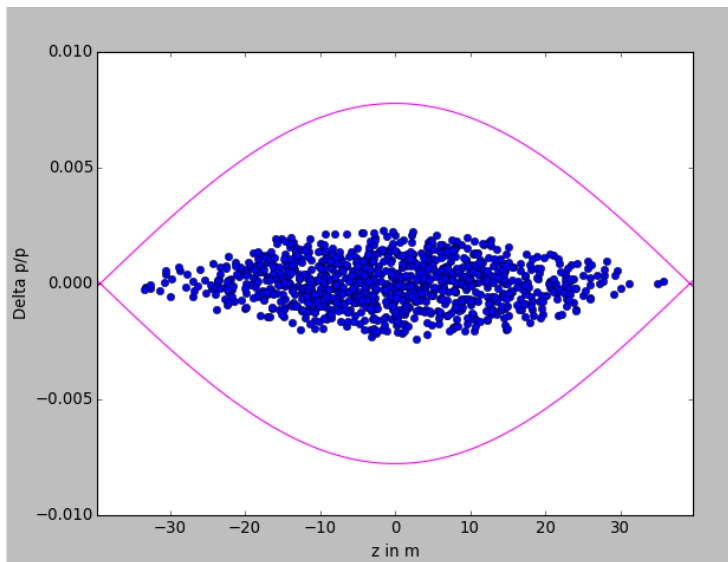
→ Bunch rotation

→ Bunch shortening, if phased correctly, but...

Impact of voltage mismatch (too high)

- Use the nominal bunch length
- track 1000 particles for 5000 turns (plot every 100 turns)
- Change V_{RF} to $V_{RF} * 10$ after the bunch is matched to V_{RF}

```
# Define RF systems  
rfsystems = RFSystems(circumference, [harmonic], [V_rf*10], [phi_offset],  
                      alpha_c_array, gamma, p_increment=p_increment_0)
```

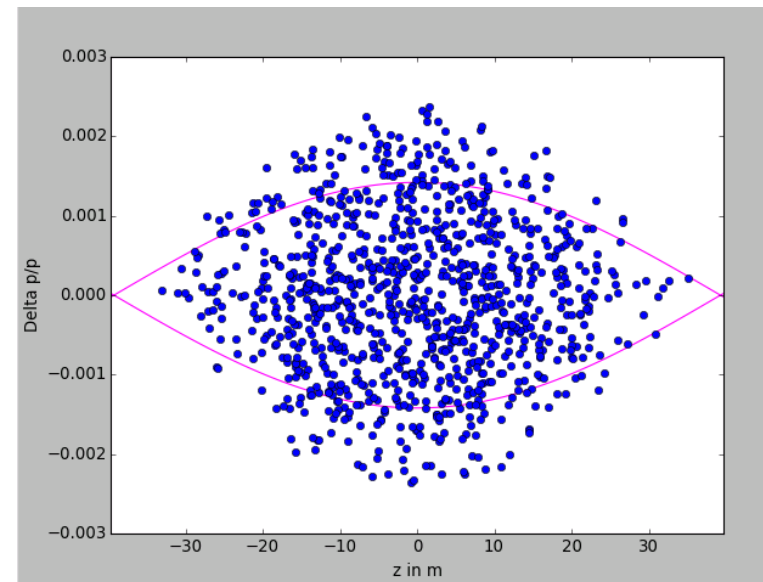
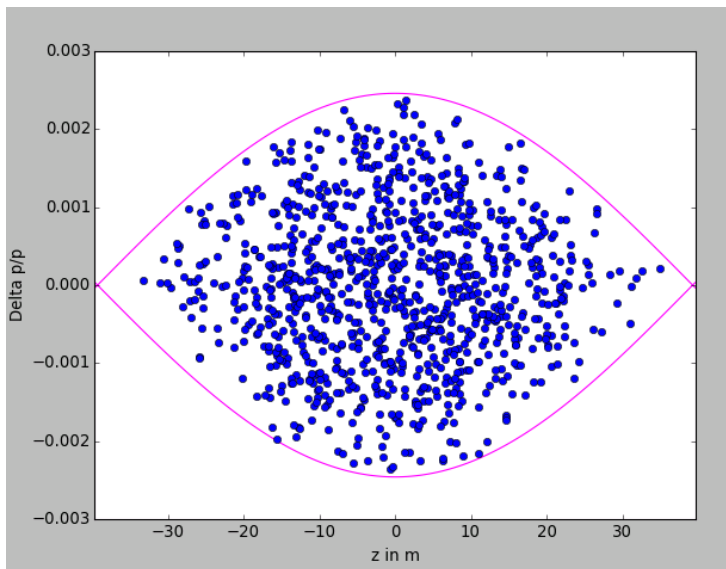


→ But should not wait too long!

→ Emittance growth!

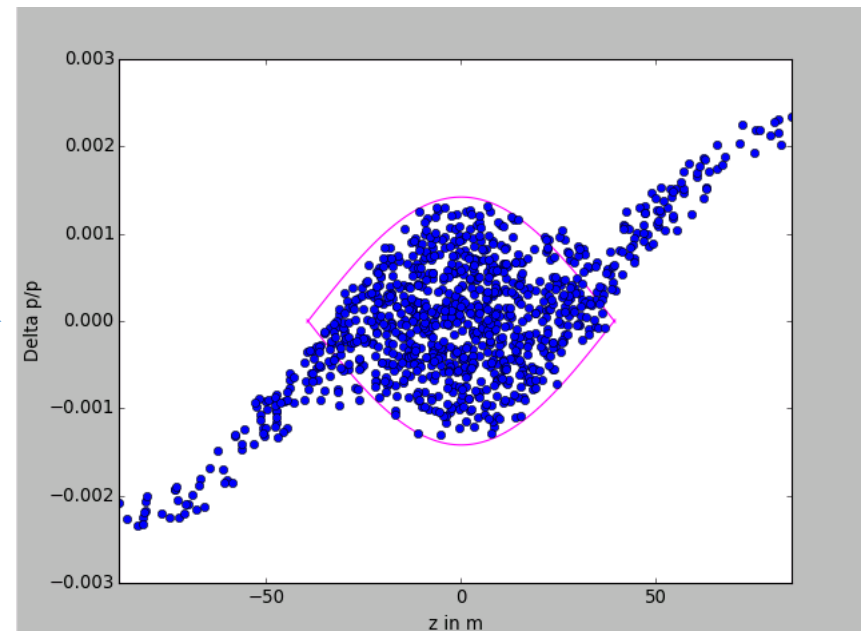
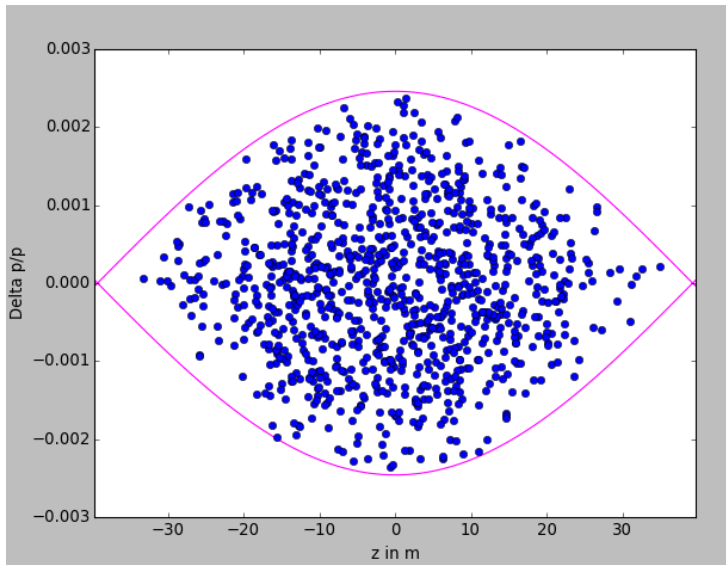
Impact of voltage mismatch (too low)

- Use the nominal bunch length
- track 1000 particles for 500 turns (plot every 10 turns)
- Change V_{RF} to $V_{RF}/10$ after the bunch is matched to V_{RF}
- Adapt the deltap plotting limits



Impact of voltage mismatch (too low)

- Use the nominal bunch length
- track 1000 particles for 500 turns (plot every 10 turns)
- Change V_{RF} to $V_{RF}/10$ after the bunch is matched to V_{RF}
- Adapt the deltap plotting limits



Agenda

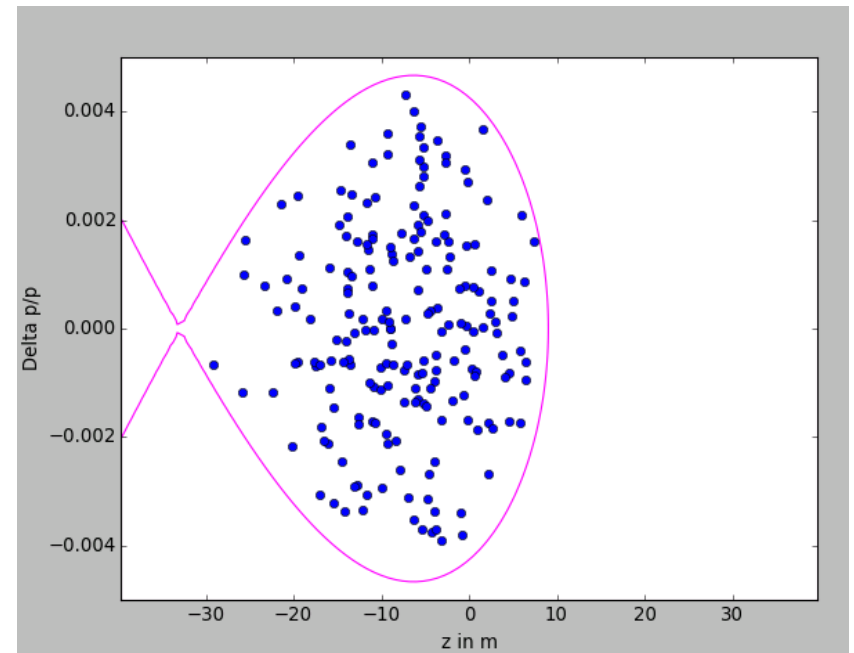
- Goals
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- Introduction to PyHEADTAIL
- Structure of ipython files
- **Tracking examples**
 - Impact of phase mismatch
 - Impact of voltage mismatch
 - **Impact of acceleration**
- Setting up the environment
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 - Ubuntu
 - Anaconda
 - PyHEADTAIL

Impact of acceleration

- Use the nominal bunch length
- Use 200 particles
- Use $\dot{B}=2.2 \text{ T/s}$
- Does it work? Why?

Impact of acceleration

- Use the nominal bunch length
- Use 1000 particles
- Use $B\dot{=}2.2$ T/s
- Change VRF to 200 kV
- Compute the synchronous phase
- Plot the phase space in ϕ [degrees], ΔE [MeV]



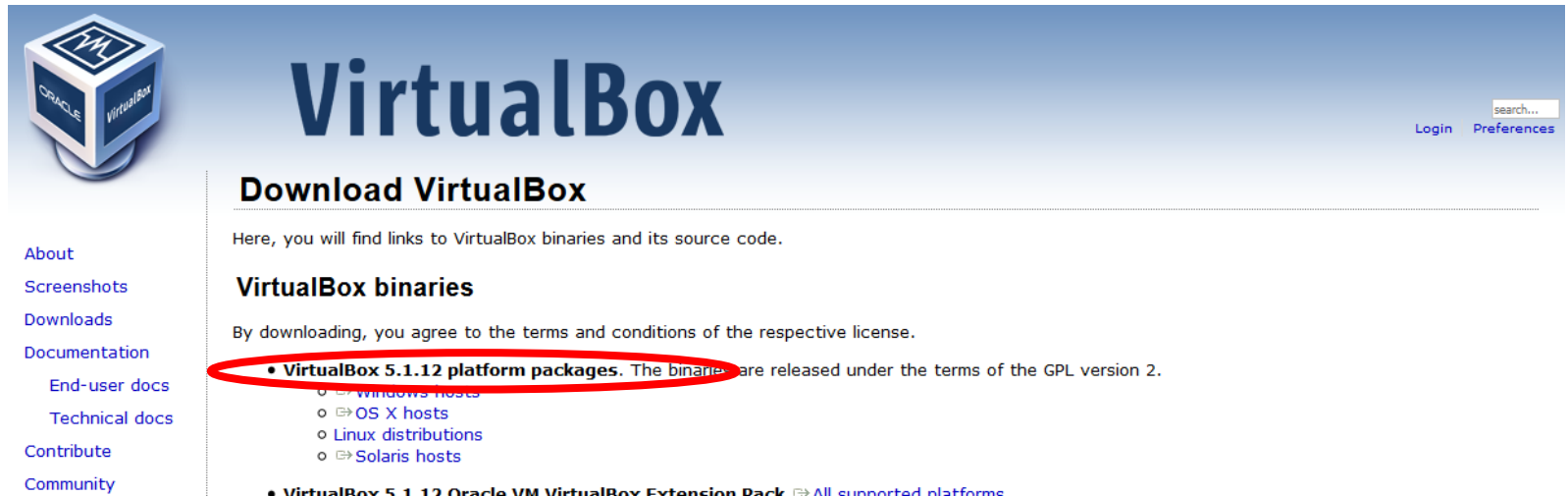
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Setting up the operating system

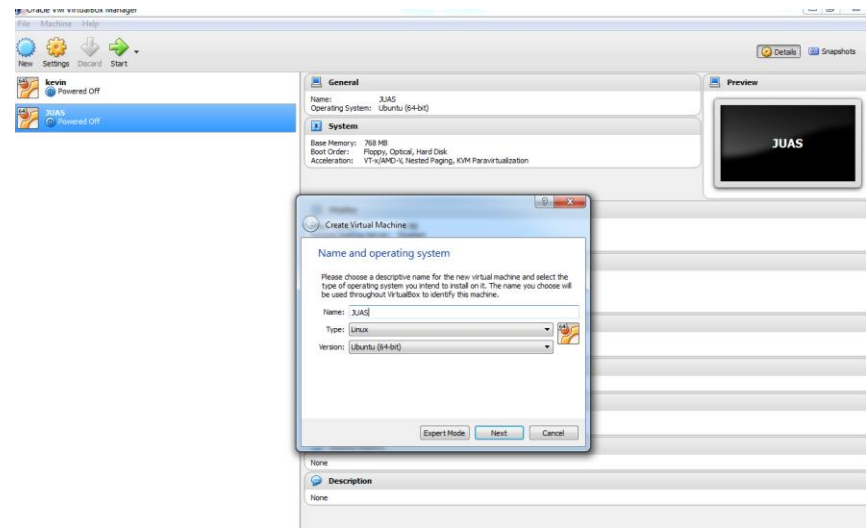
- Install a virtual box to get an Ubuntu environment

Ex: <https://www.virtualbox.org/wiki/Downloads>



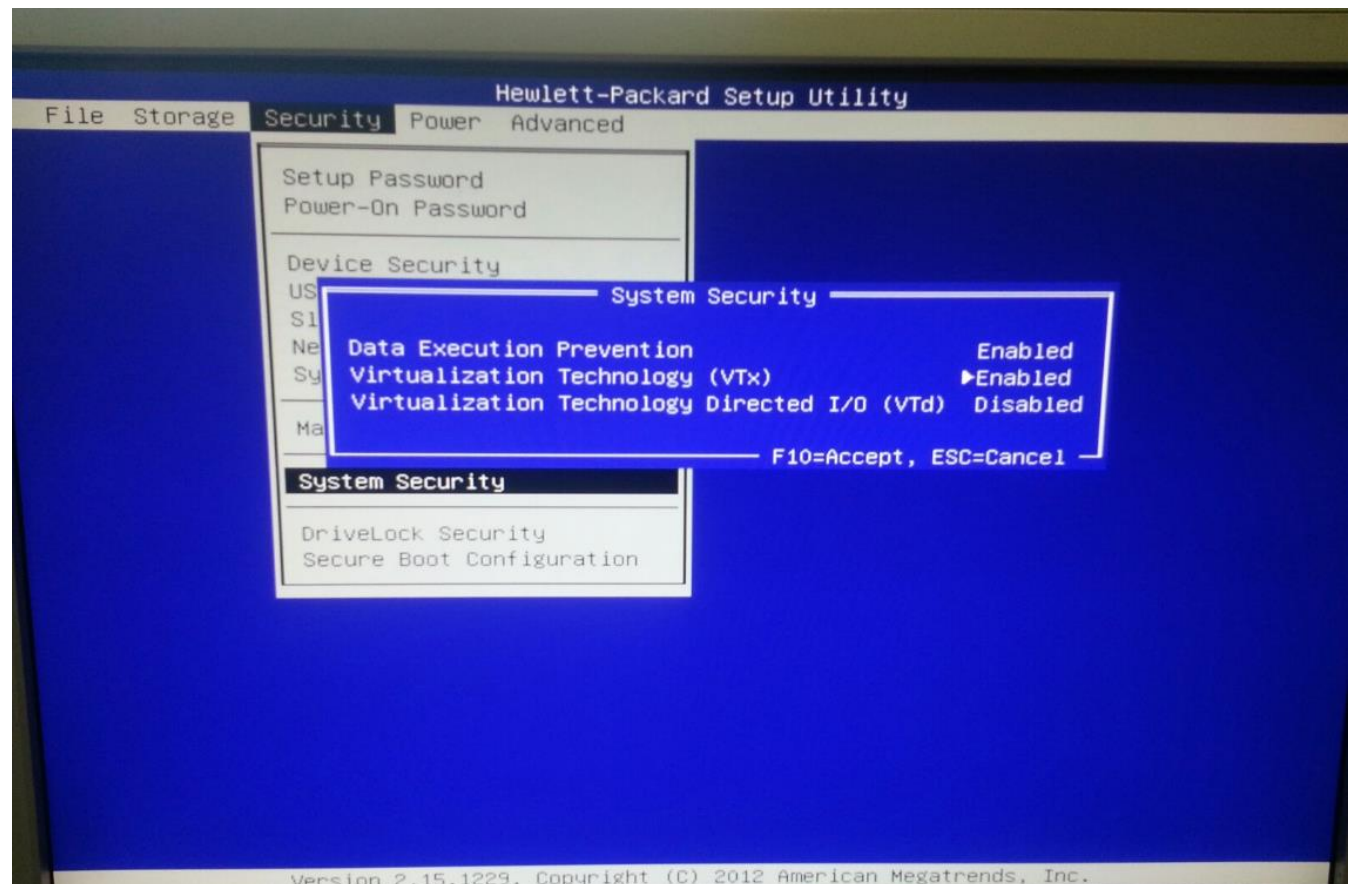
The screenshot shows the VirtualBox website homepage. On the left is a navigation menu with links for About, Screenshots, Downloads, Documentation, End-user docs, Technical docs, Contribute, and Community. The main content area features the VirtualBox logo and a heading 'Download VirtualBox'. Below this, it states 'Here, you will find links to VirtualBox binaries and its source code.' and 'VirtualBox binaries'. A red circle highlights the text 'VirtualBox 5.1.12 platform packages. The binaries are released under the terms of the GPL version 2.' followed by a list of host types: windows hosts, OS X hosts, Linux distributions, and Solaris hosts. At the bottom, there is a link for 'VirtualBox 5.1.12 Oracle VM VirtualBox Extension Pack' and 'All supported platforms'.

- Create a new virtual machine with Ubuntu 64-bit

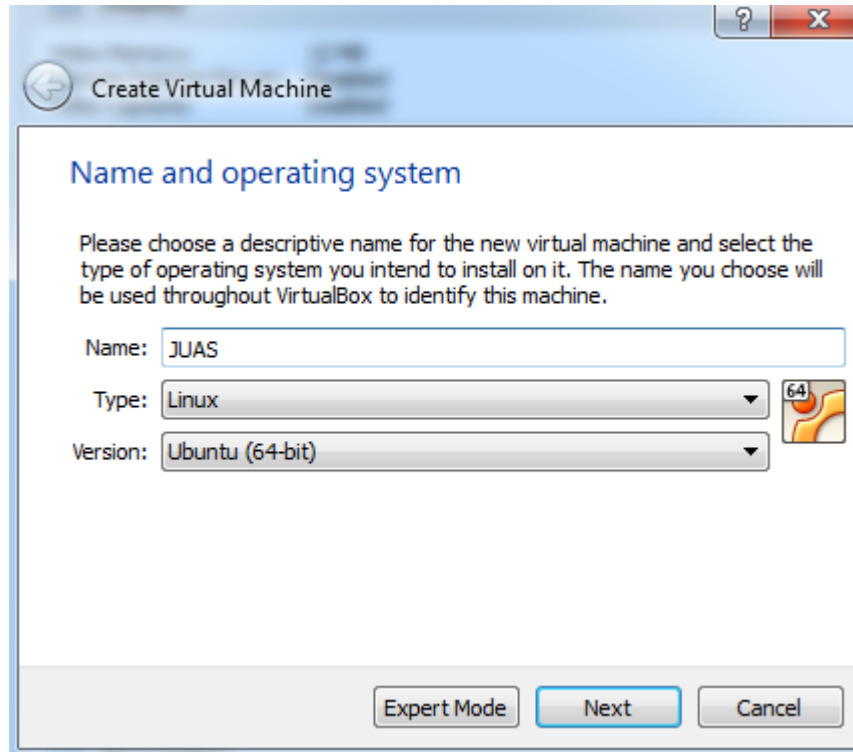


The screenshot shows the Oracle VM VirtualBox Manager interface. The main window displays a list of virtual machines: 'Kevin' (Powered Off) and 'JUAS' (Powered Off). The 'JUAS' machine is selected, and its settings are visible in the 'General' tab. The 'Name' is 'JUAS' and the 'Operating System' is 'Ubuntu (64-bit)'. A 'Create Virtual Machine' dialog box is open in the foreground, showing the 'Name and operating system' section. The 'Name' is 'JUAS', the 'Type' is 'Linux', and the 'Version' is 'Ubuntu (64-bit)'. The dialog box has 'Expert Mode', 'Next', and 'Cancel' buttons.

In case of a “VTx” error or an impossibility to select 64 bit operating system when launching the virtual machine, one needs to turn Virtualization Technology (VTx) on in the BIOS (see chapter 10.3 in <https://www.virtualbox.org/manual/ch10.html#hwvirt>)



Create virtual machine




Create Virtual Machine

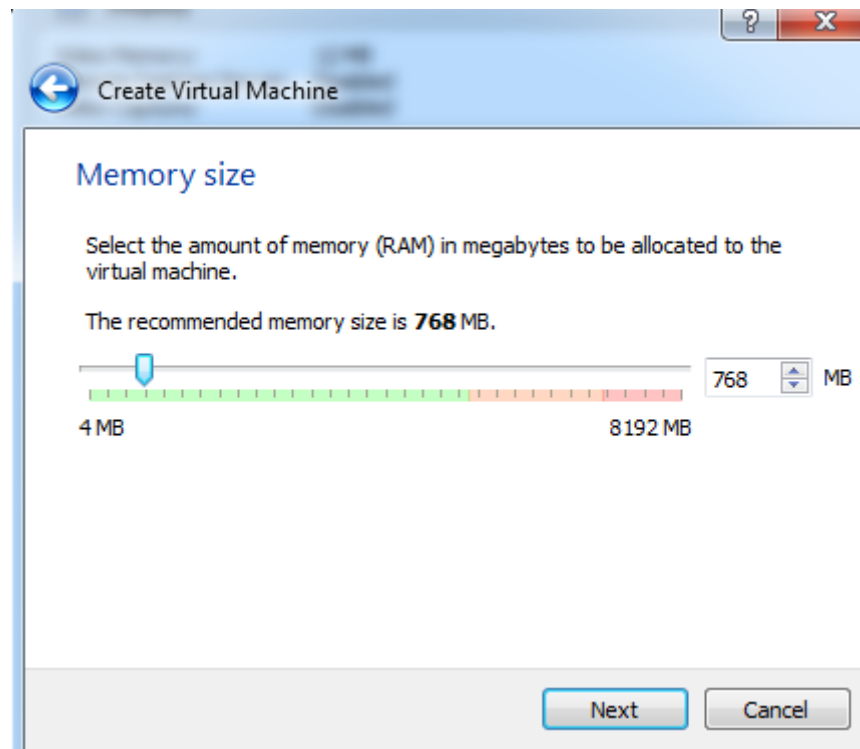
Name and operating system

Please choose a descriptive name for the new virtual machine and select the type of operating system you intend to install on it. The name you choose will be used throughout VirtualBox to identify this machine.

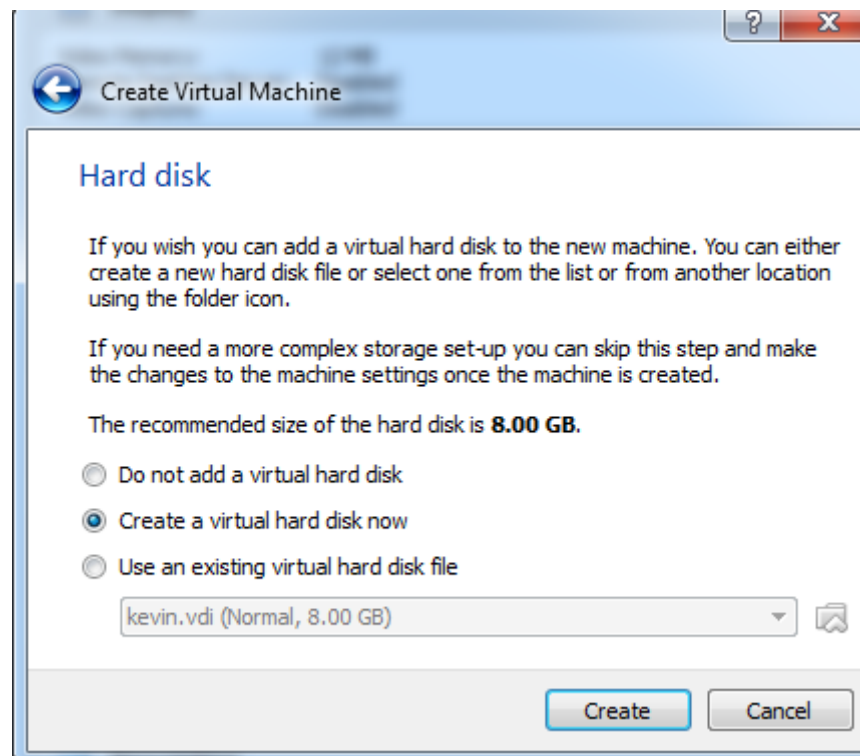
Name:

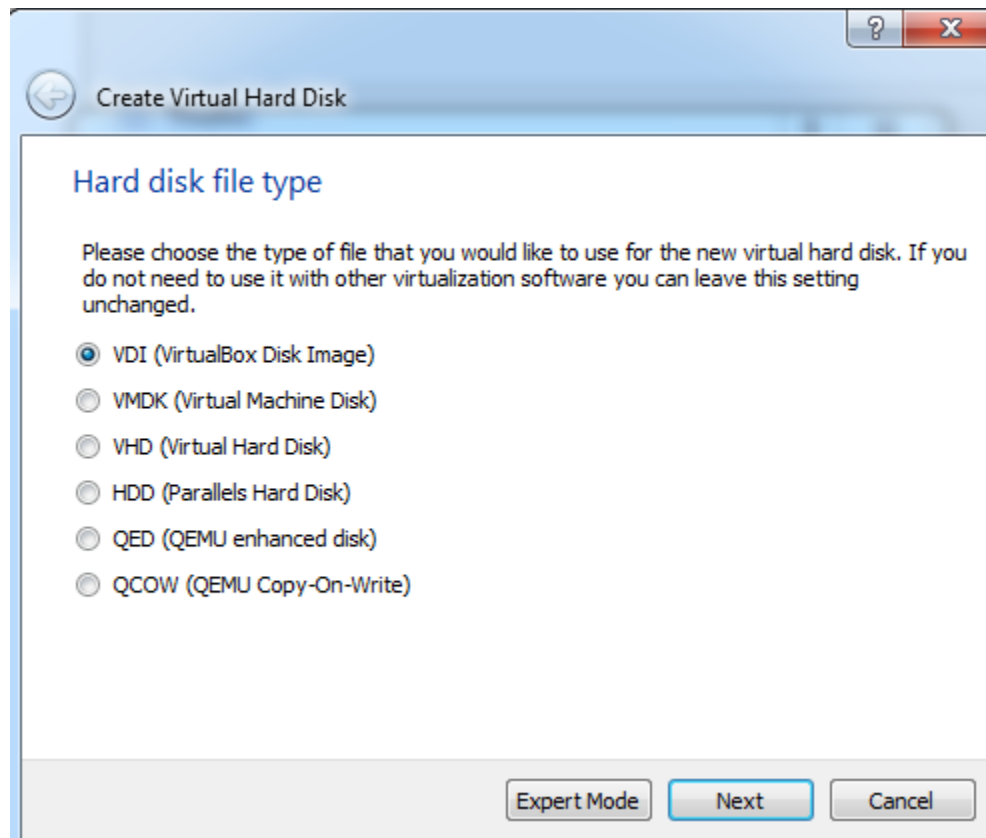
Type: 

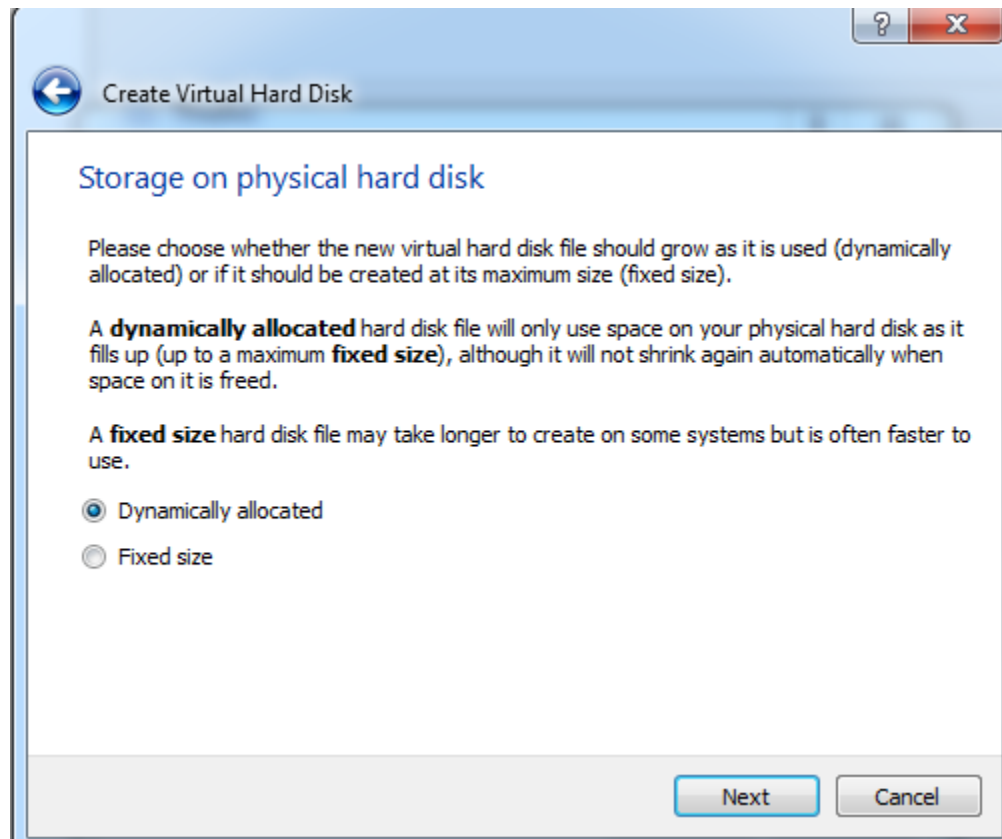
Version:

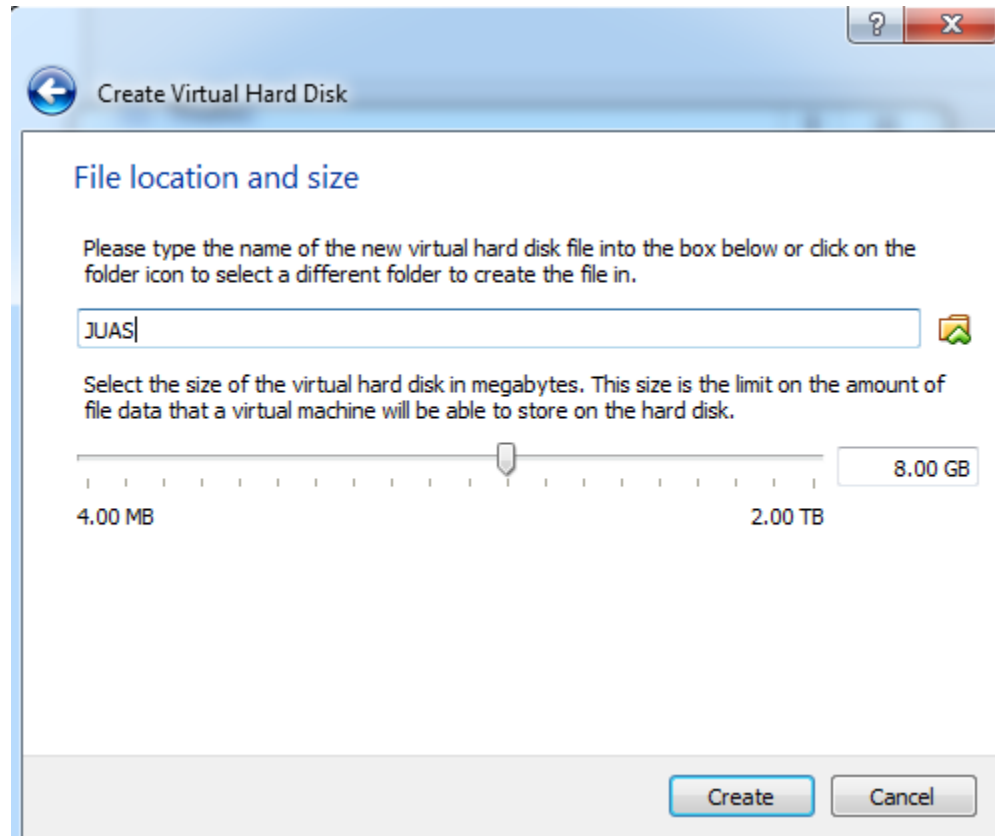


Actually 2 GB is recommended by Ubuntu. This setting can be modified later.



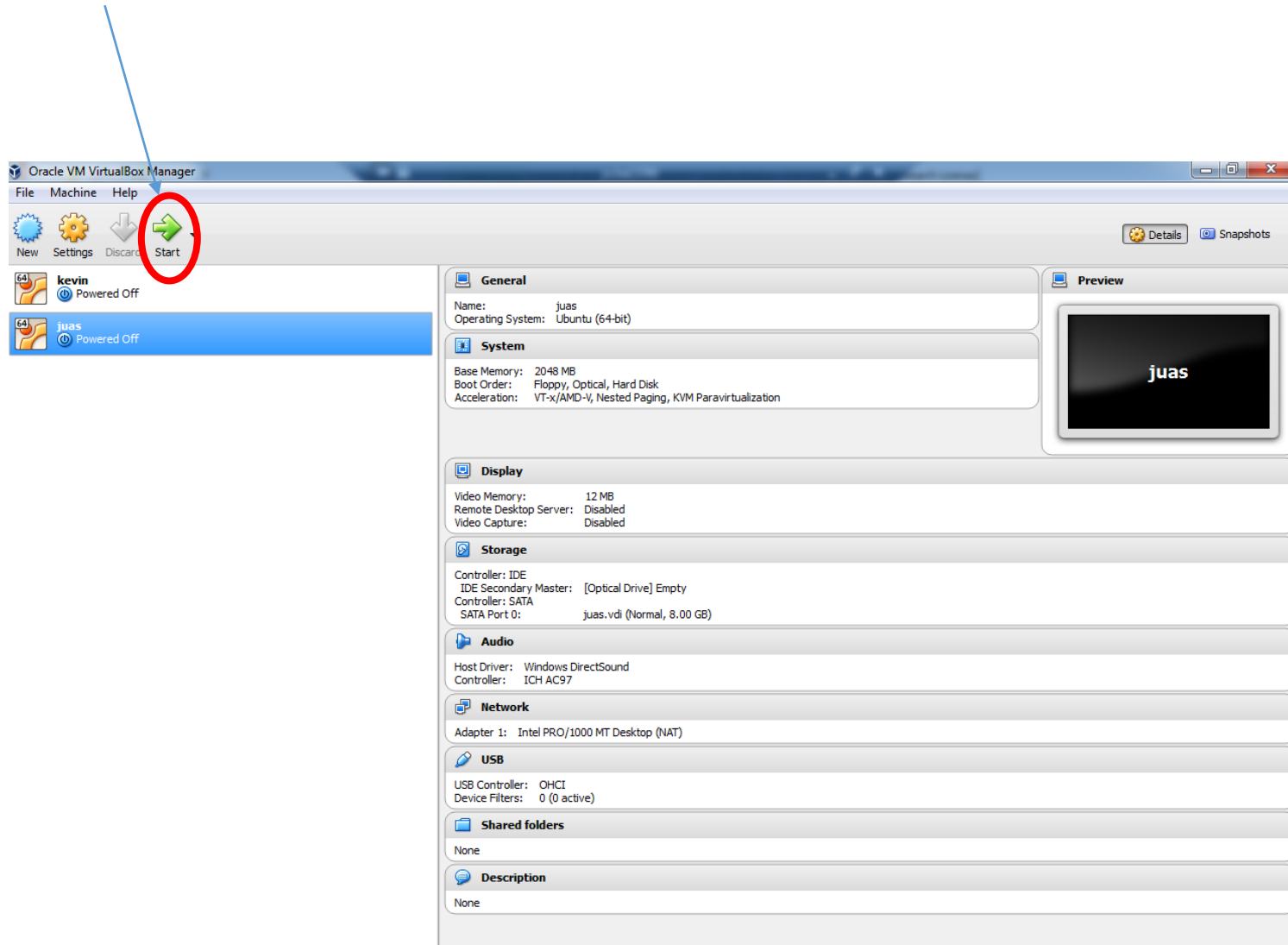






→ Assign 20 GB if possible.

Start the VM



The screenshot shows the Oracle VM VirtualBox Manager interface. The 'Start' button in the top toolbar is circled in red, with a blue arrow pointing to it from the title 'Start the VM'. The interface displays two virtual machines: 'kevin' and 'juas', both in a 'Powered Off' state. The 'juas' VM is selected, and its configuration details are shown on the right side of the window.

General
Name: juas
Operating System: Ubuntu (64-bit)

System
Base Memory: 2048 MB
Boot Order: Floppy, Optical, Hard Disk
Acceleration: VT-x/AMD-V, Nested Paging, KVM Paravirtualization

Display
Video Memory: 12 MB
Remote Desktop Server: Disabled
Video Capture: Disabled

Storage
Controller: IDE
IDE Secondary Master: [Optical Drive] Empty
Controller: SATA
SATA Port 0: juas.vdi (Normal, 8.00 GB)

Audio
Host Driver: Windows DirectSound
Controller: ICH AC97

Network
Adapter 1: Intel PRO/1000 MT Desktop (NAT)

USB
USB Controller: OHCI
Device Filters: 0 (0 active)

Shared folders
None

Description
None

Preview
juas

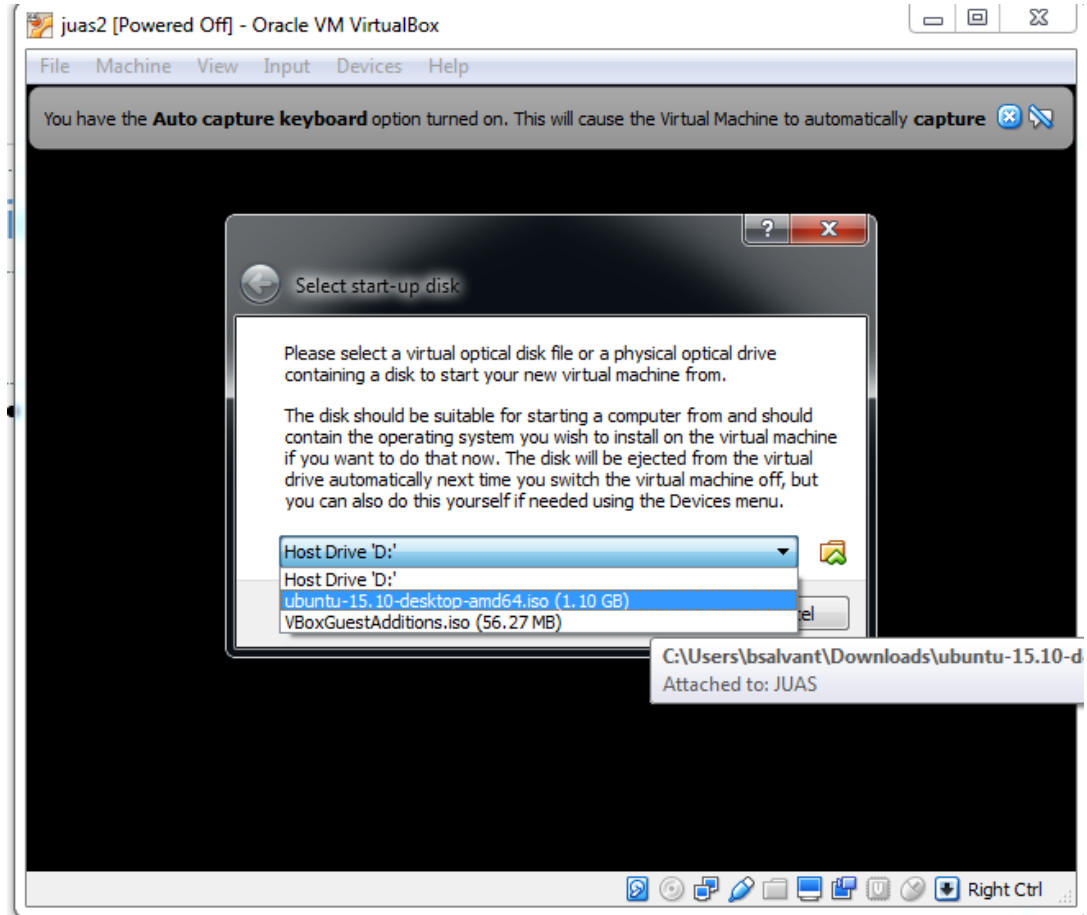
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Download ubuntu

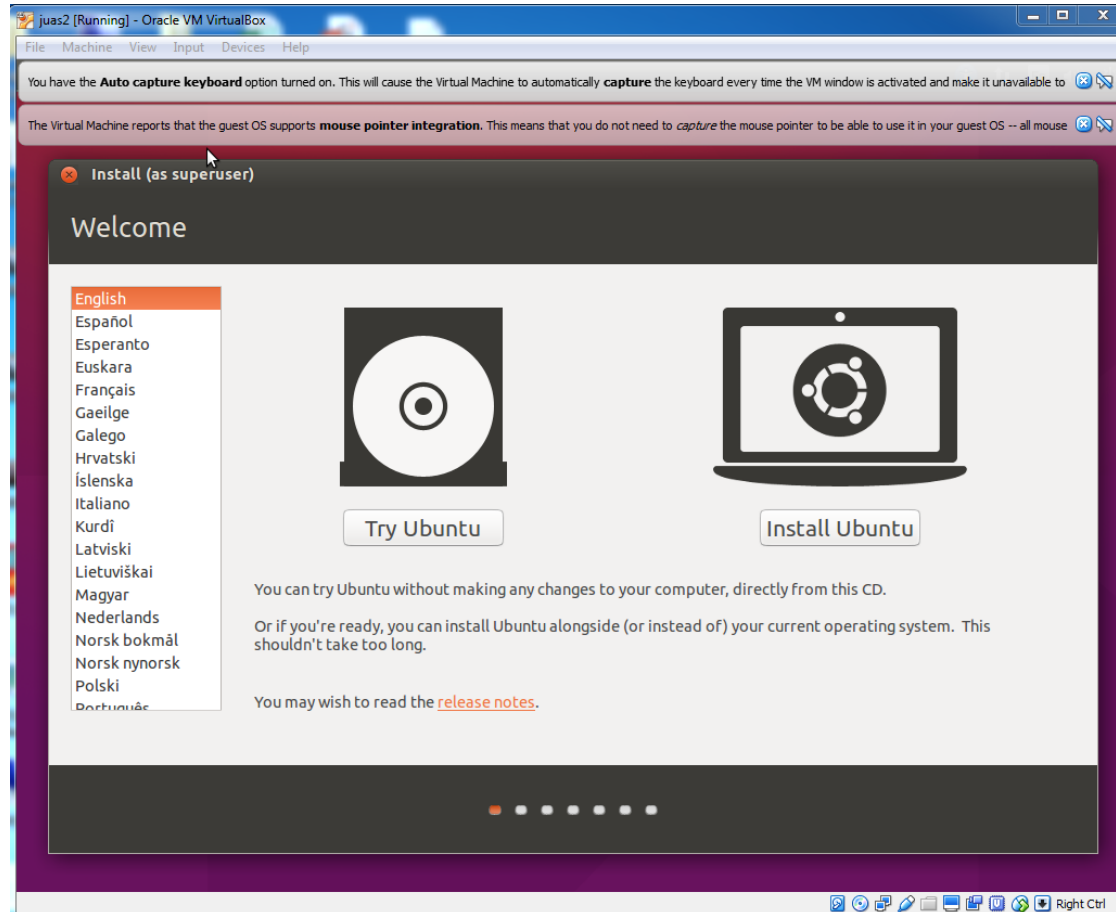
- <https://www.ubuntu.com/download/desktop>

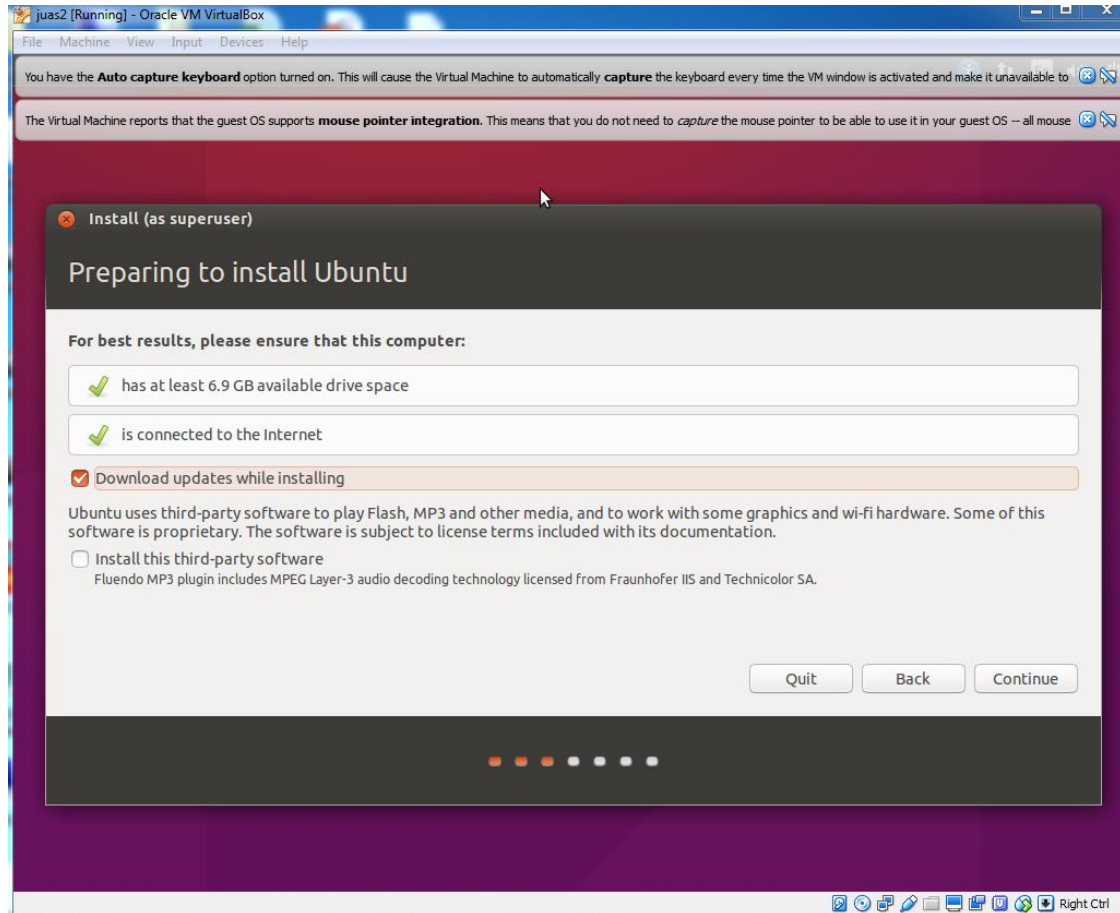
Select Ubuntu start up disk



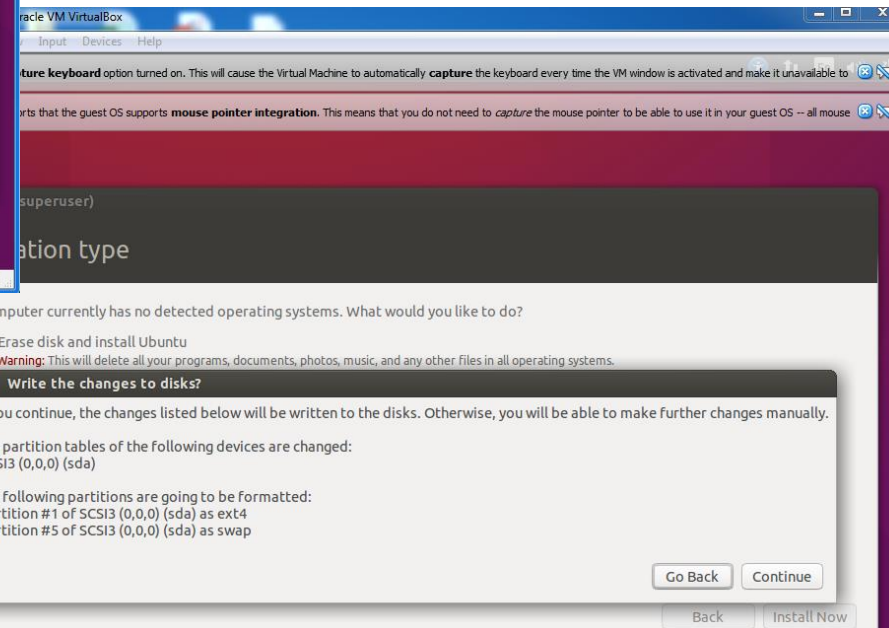
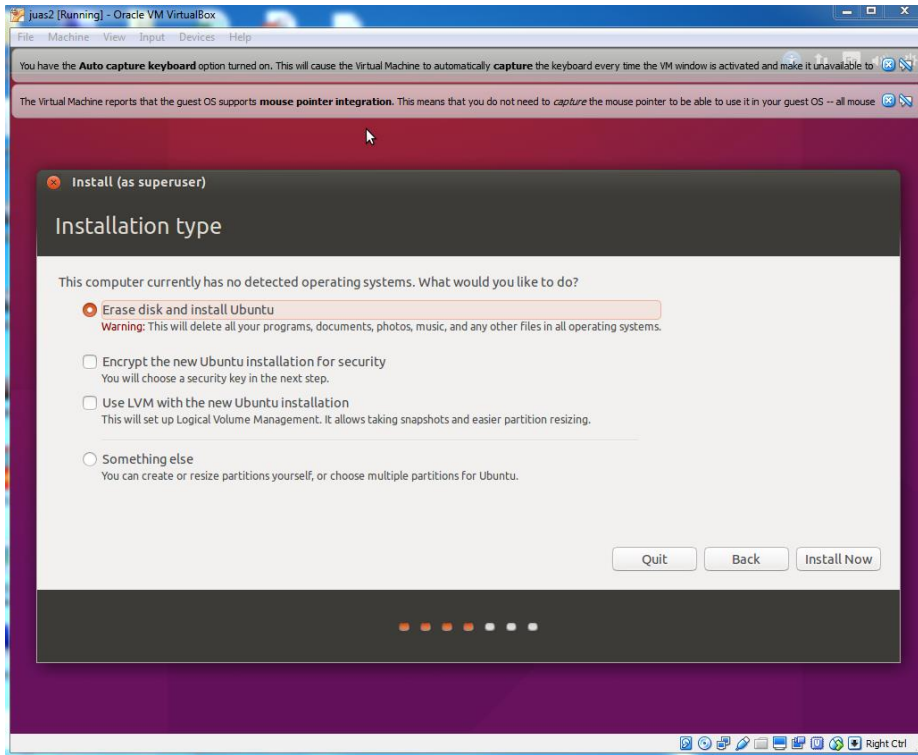
You may have to click on the small green arrow to browse for your Ubuntu installation image

Install ubuntu





Keep default installation and “install now” and confirm with “continue” on the pop up



Then choose time zone and login details

Then wait for file copy and Ubuntu installation restart the machine after it is complete

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Install anaconda

Go to <https://www.continuum.io/downloads> and follow instructions

The screenshot shows a Firefox browser window with the URL <https://www.continuum.io/downloads>. The page content includes:

- Navigation tabs: [Download for Windows](#), [Download for OSX](#), and [Download for Linux](#) (selected).
- Section: **Anaconda 4.2.0 For Linux**
- Text: "Anaconda is BSD licensed which gives you permission to use Anaconda commercially and for redistribution."
- Link: [Changelog](#)
- Instructions:
 1. Download the installer
 2. Optional: Verify data integrity with [MD5](#) or [SHA-256](#) [More info](#)
 3. In your terminal window type one of the below and follow the instructions:
- Python 3.5 version:
 - 64-BIT INSTALLER (455M) (green button)
 - 32-BIT INSTALLER (373M) (green button)
- Python 2.7 version (circled in red):
 - 64-BIT INSTALLER (446M) (blue button)
 - 32-BIT INSTALLER (365M) (blue button)
- Terminal commands:

```
bash Anaconda3-4.2.0-Linux-x86_64.sh
```

```
bash Anaconda2-4.2.0-Linux-x86_64.sh
```
- NOTE: Include the "bash" command even if you are not using the bash shell.

Do not forget to add Anaconda to path
(answer yes or add the path yourself in the
“.bashrc” configuration file)

```
Python 2.7.12 :: Continuum Analytics, Inc.  
creating default environment...  
installation finished.  
Do you wish the installer to prepend the Anaconda2 install location  
to PATH in your /home/juas/.bashrc ?  [yes]  [no]  
[no] >>>  
You may wish to edit your .bashrc or prepend the Anaconda2 install location:  
  
$ export PATH=/home/juas/anaconda2/bin:$PATH  
  
Thank you for installing Anaconda2!
```

- Restart the shell (“>>source .bashrc” from the home directory)
- Check that “>>ipython notebook” works

- Anaconda contains Python and all the necessary libraries to run PyHEADTAIL.

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Installing PyHEADTAIL

- website for installation and information:
<https://github.com/PyCOMPLETE/PyHEADTAIL>

```
juas@juas-VirtualBox:~/Downloads$ sudo apt install git
```

```
juas@juas-VirtualBox:~/Downloads$ git clone https://github.com/PyCOMPLETE/PyHEADTAIL
```

```
juas@juas-VirtualBox:~$ cd PyHEADTAIL/
```

```
juas@juas-VirtualBox:~/PyHEADTAIL$ make
```

Add path to the end of the .bashrc to be able to import PyHeadtail from everywhere:

```
juas@juas-VirtualBox:~$ gedit ~/.bashrc
```

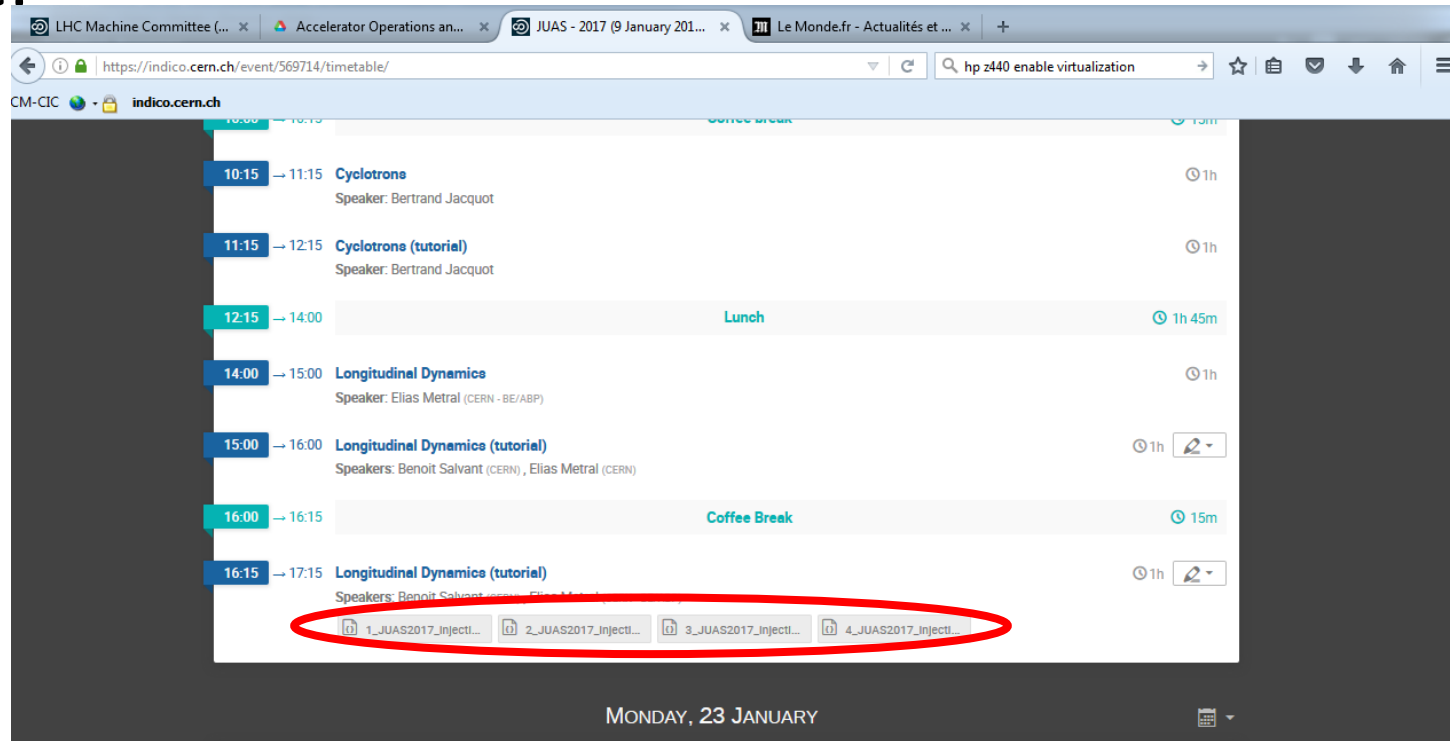
```
# Anaconda
export PATH=/home/juas/anaconda2/bin:$PATH

# PyHEADTAIL
export PYTHONPATH=/home/juas/:$PYTHONPATH
```

This only updates once a new terminal is opened (or do `source ~/.bashrc`)

Finding the examples

- From JUAS Indico site in the last tutorial of the course:

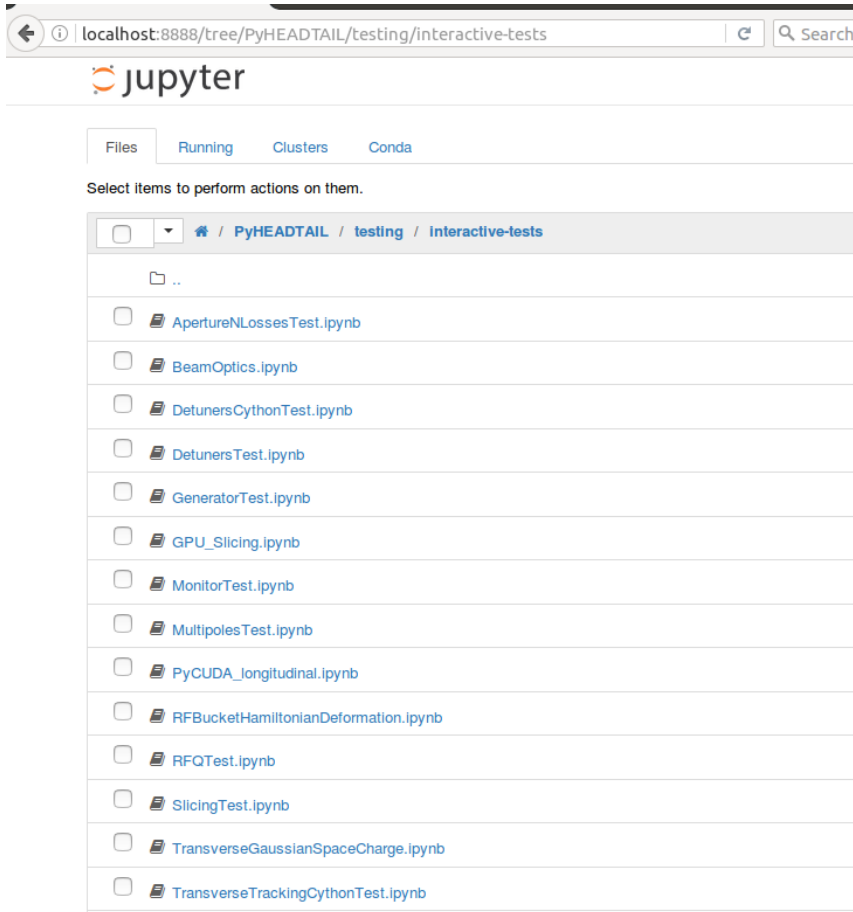


The screenshot shows a web browser displaying the Indico timetable for the JUAS 2017 event. The page lists several sessions with their titles, speakers, and durations. At the bottom of the page, there are four download links for files named '1_JUAS2017_Inject...', '2_JUAS2017_Inject...', '3_JUAS2017_Inject...', and '4_JUAS2017_Inject...'. These links are circled in red, indicating they are the files to be downloaded.

Time	Session Title	Speaker(s)	Duration
10:15	Cyclotrons	Bertrand Jacquot	1h
11:15	Cyclotrons (tutorial)	Bertrand Jacquot	1h
12:15	Lunch		1h 45m
14:00	Longitudinal Dynamics	Elias Metral (CERN - BE/ABP)	1h
15:00	Longitudinal Dynamics (tutorial)	Benoit Salvant (CERN), Elias Metral (CERN)	1h
16:00	Coffee Break		15m
16:15	Longitudinal Dynamics (tutorial)	Benoit Salvant (CERN), Elias Metral (CERN)	1h

→ Download the files and open them browsing with “ipython notebook”

Other examples available with the PyHEADTAIL install



The screenshot shows a web browser window at the URL `localhost:8888/tree/PyHEADTAIL/testing/interactive-tests`. The Jupyter logo is visible at the top. Below the logo are tabs for `Files`, `Running`, `Clusters`, and `Conda`. A message says "Select items to perform actions on them." Below this is a file browser interface with a breadcrumb path: `PyHEADTAIL / testing / Interactive-tests`. The file list includes:

- `ApertureNLossesTest.ipynb`
- `BeamOptics.ipynb`
- `DetunersCythonTest.ipynb`
- `DetunersTest.ipynb`
- `GeneratorTest.ipynb`
- `GPU_Slicing.ipynb`
- `MonitorTest.ipynb`
- `MultipolesTest.ipynb`
- `PyCUDA_longitudinal.ipynb`
- `RFBucketHamiltonianDeformation.ipynb`
- `RFQTest.ipynb`
- `SlicingTest.ipynb`
- `TransverseGaussianSpaceCharge.ipynb`
- `TransverseTrackingCythonTest.ipynb`

Other interesting examples to run

- Doublet beam creation
- PS to SPS transfer
- Double and triple splitting

Changing variables

JUAS

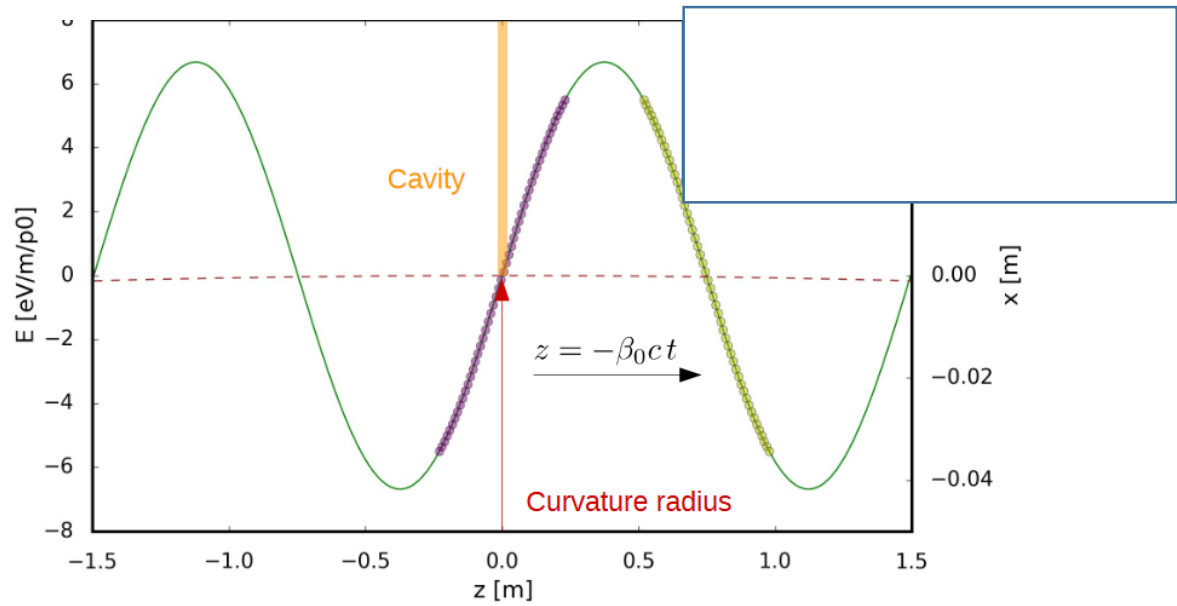
Johns Hopkins Accelerator School

- ◆ Change of variables if one wants to use $(\Phi, \Delta E)$ or $(\Delta t, \Delta E)$ instead of $(\Phi, d\Phi/dt)$

$$\begin{aligned}\Delta\phi &= \phi - \phi_s \\ &= \omega_{RF} \Delta t \\ &= h \omega_s \Delta t\end{aligned}$$

$$\Delta p = \frac{\Delta E}{\beta_s c}$$

$$\dot{\phi} = -\frac{\eta h c}{\beta_s E_s R_s} \Delta E$$



$$\Delta\phi[\text{rad}] = -\frac{h}{R} z[\text{m}]$$

$$\Delta E = \frac{\Delta p}{p} m_0 \gamma \beta^2 c^2$$

For protons

$$\Delta E[\text{GeV}] = \frac{\Delta p}{p} 0.938 \gamma \beta^2$$