

PyHEADTAIL examples

All our thanks to
D. Amorim, Kevin Li and Michael Schenk
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

- 1h00 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**
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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!

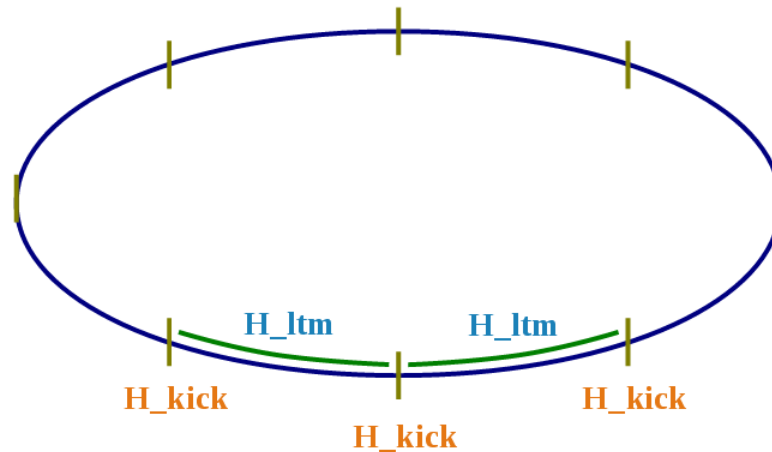
introduction to PyHEADTAIL

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
```

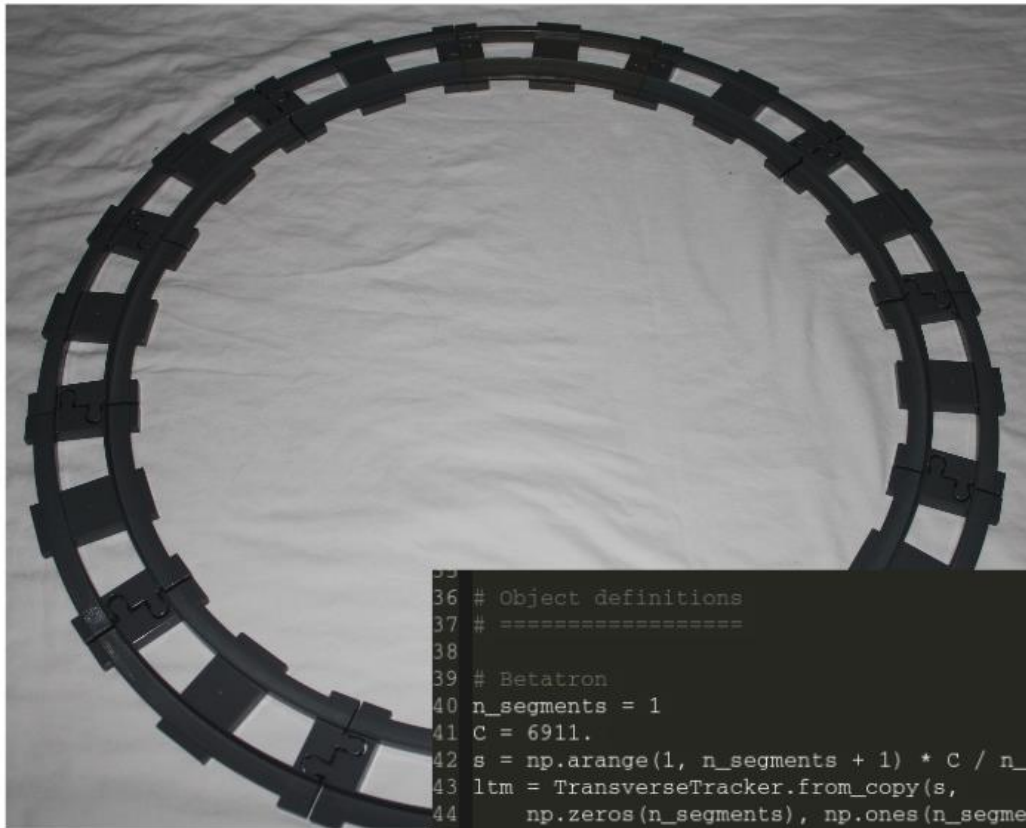


introduction to PyHEADTAIL



A real world example

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- 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
```

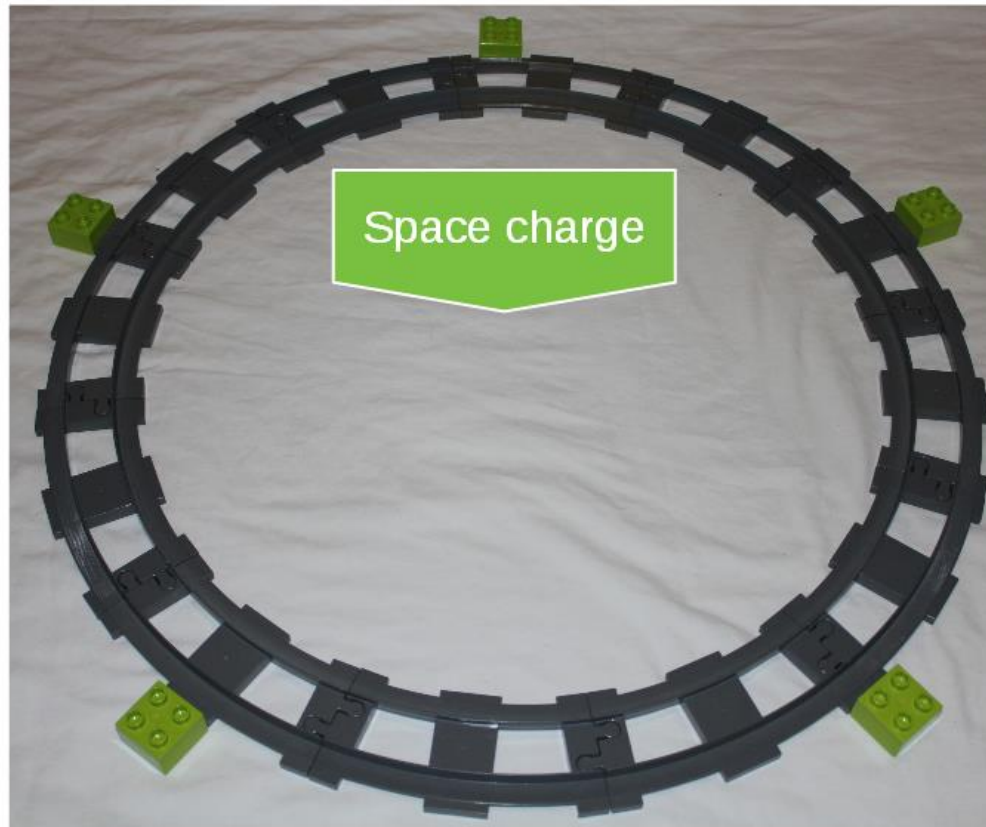


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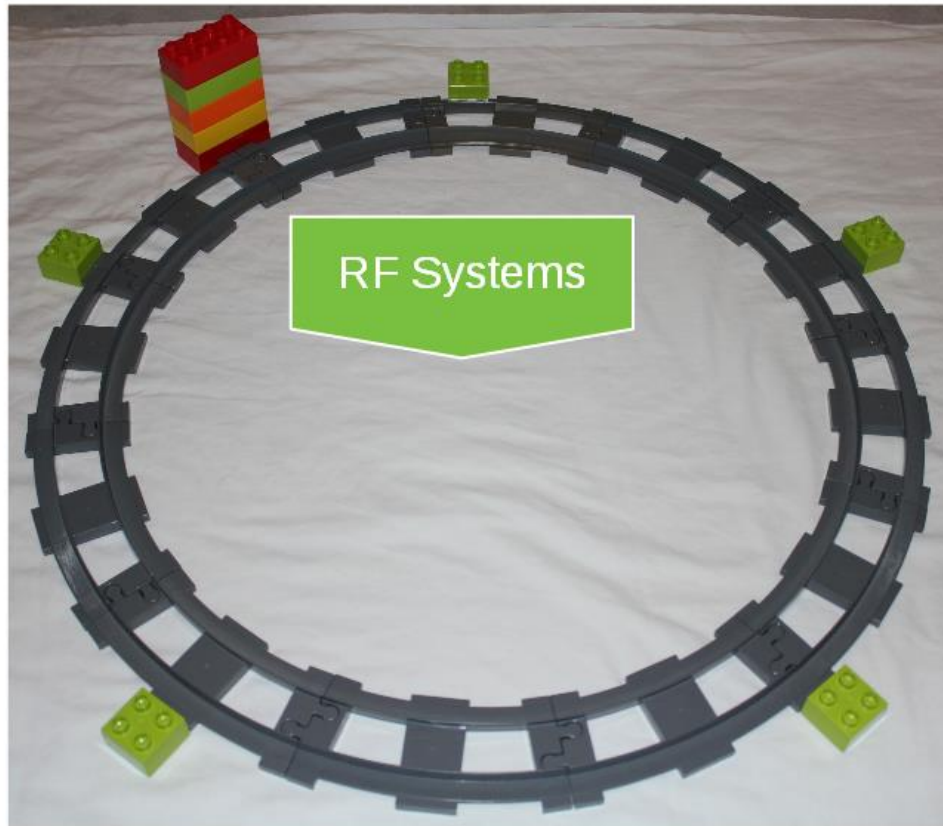
- Load Python packages and modules
- Build linear periodic transfer maps
- Add (collective) kick elements

introduction to PyHEADTAIL



A real world example

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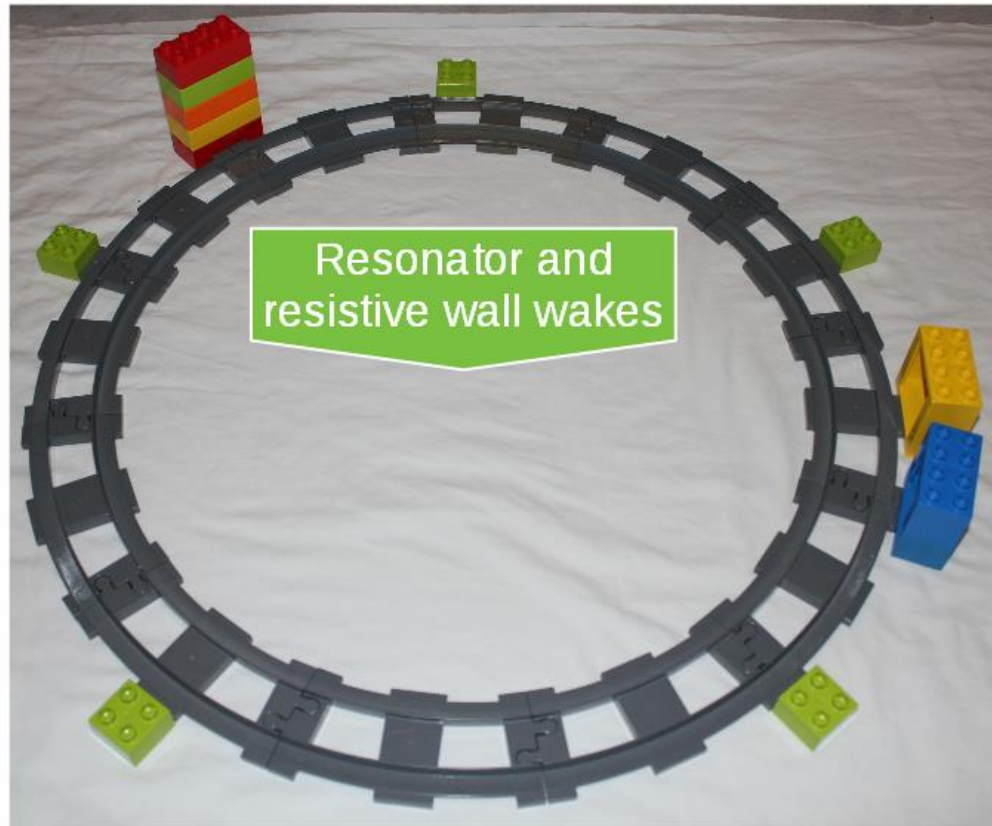
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introduction to PyHEADTAIL



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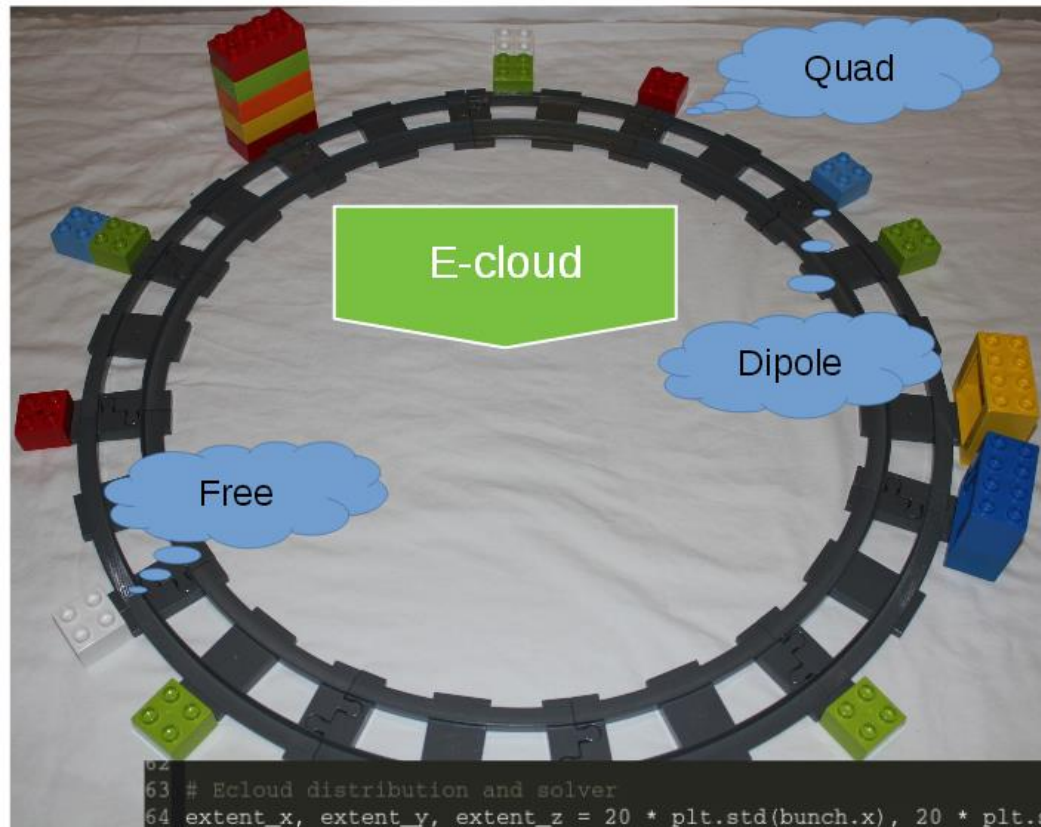
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introduction to PyHEADTAIL



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
```

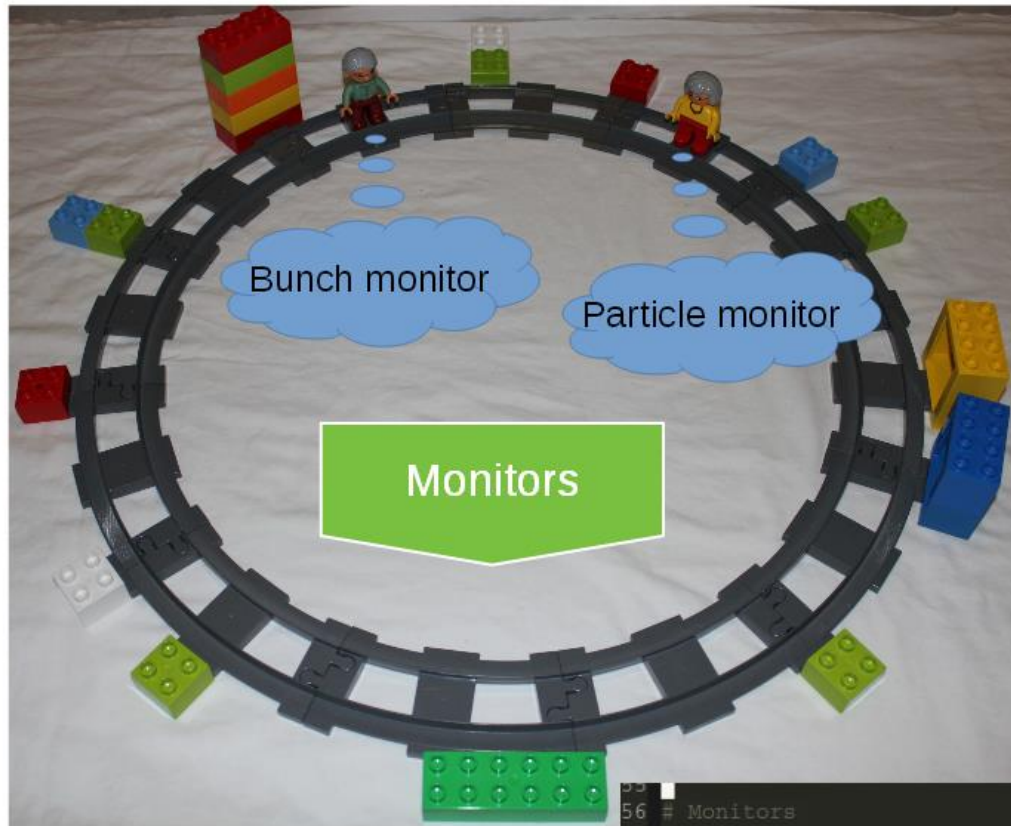


introduction to PyHEADTAIL



A real world example

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USPAS 2015



- Load Python packages and modules
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```
56 # Monitors
57 bunchmonitor = BunchMonitor('bunch', n_turns, slices)
58 particlemonitor = ParticleMonitor('particles', n_turns, slices)
```

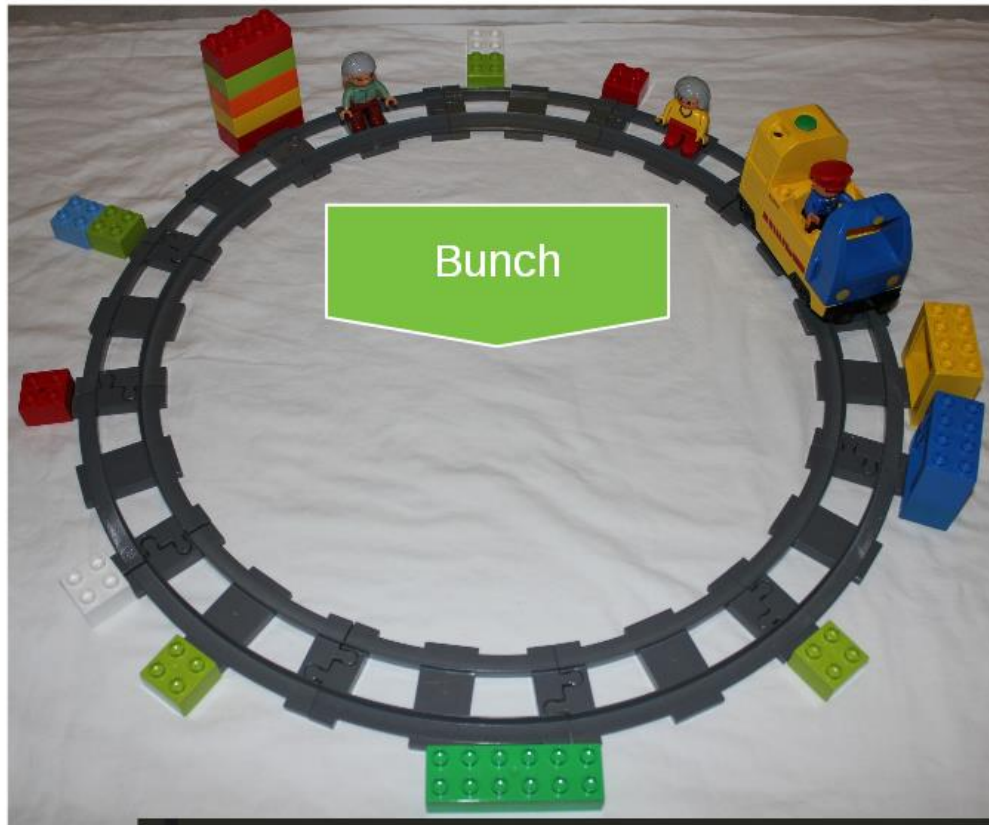


introduction to PyHEADTAIL



A real world example

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Kevin Li
USPAS 2015



- 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
```

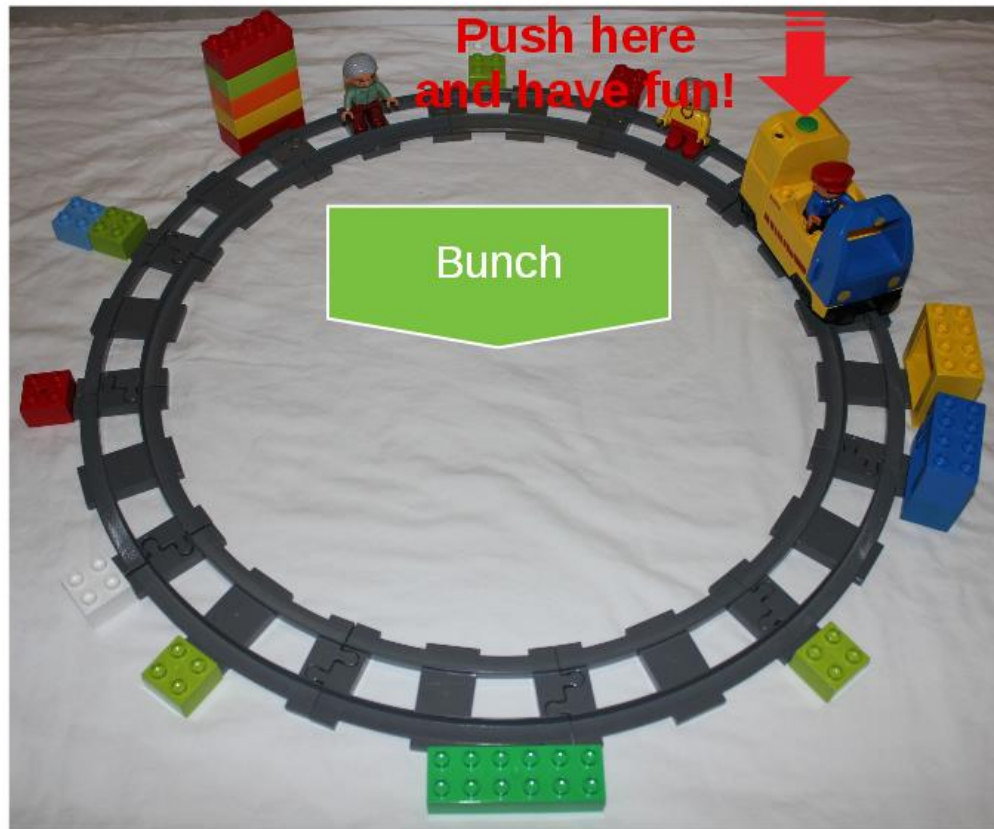


introduction to PyHEADTAIL



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

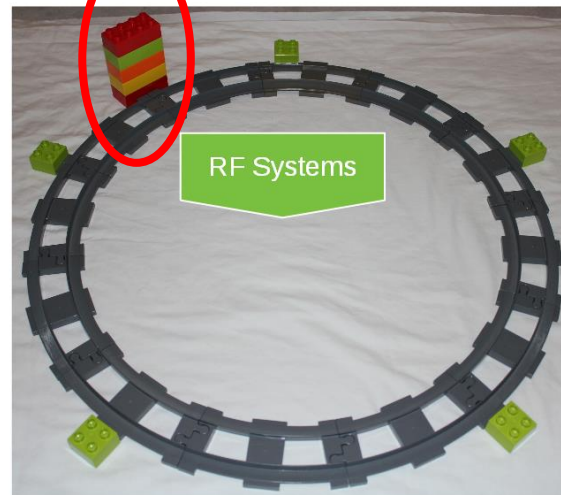
- ◆ The motion of the particles can be tracked turn by turn using the recurrence relation (between turn n and turn $n+1$)

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

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

JUAS - Jan 2016 - E.Métral

A real world example



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

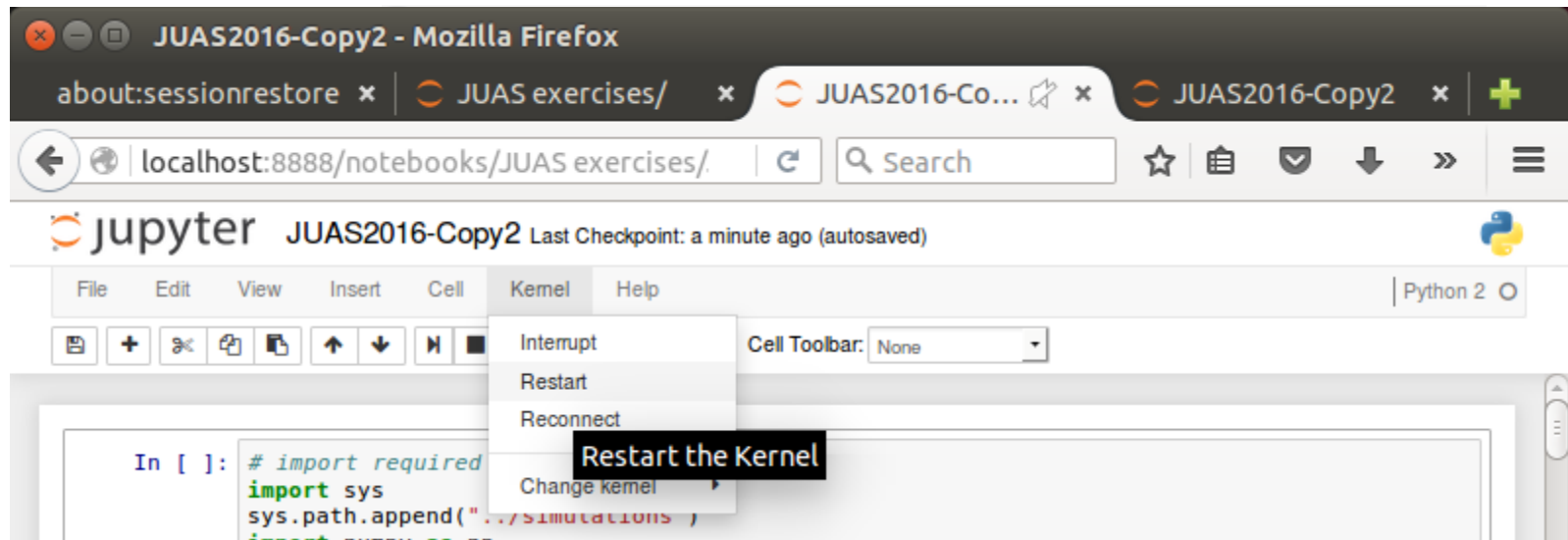
```
In [ ]: # plot phase space
plt.close()
plt.ion()
fig = plt.figure(1)
for i in np.arange(0, 50, 1):
    # track the particles
    rfsystems.track(bunch)
    if i%10 == 0:
        # monitor the particles
        bucket = rfsystems.get_bucket(gamma=bunch.gamma)
        # plot the RF bucket envelope
        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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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

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

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

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

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

20 kV at injection

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

$\Rightarrow \gamma_t \approx 6.1$

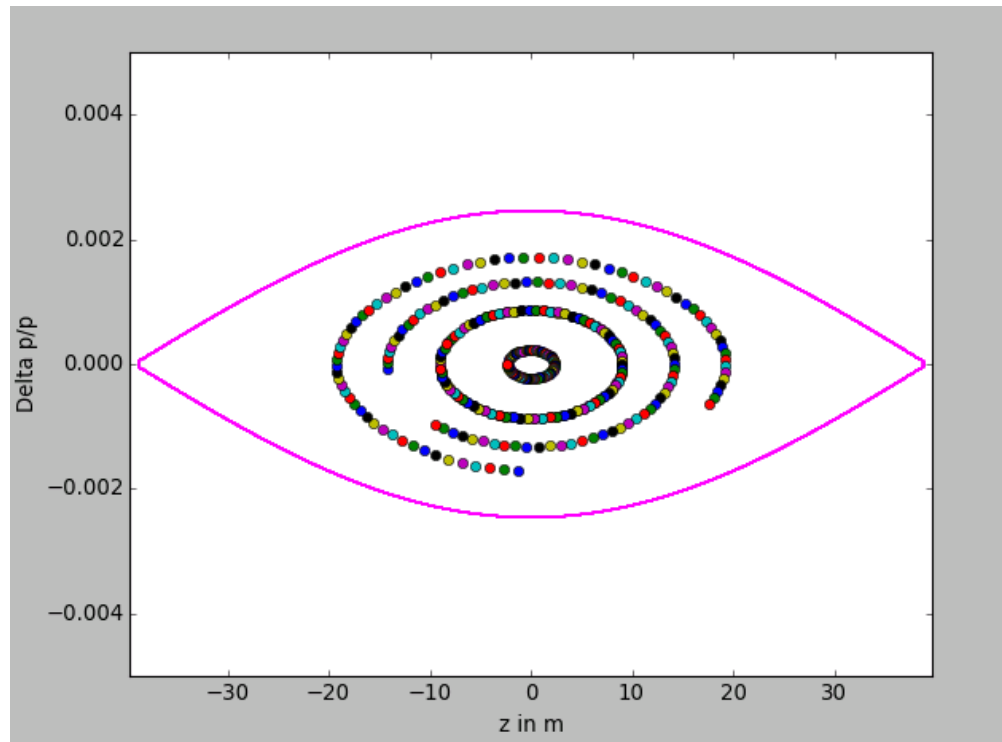
$\text{harmonic} = 8$

$\text{gamma_tr} = 6.1$

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

Let's run that example

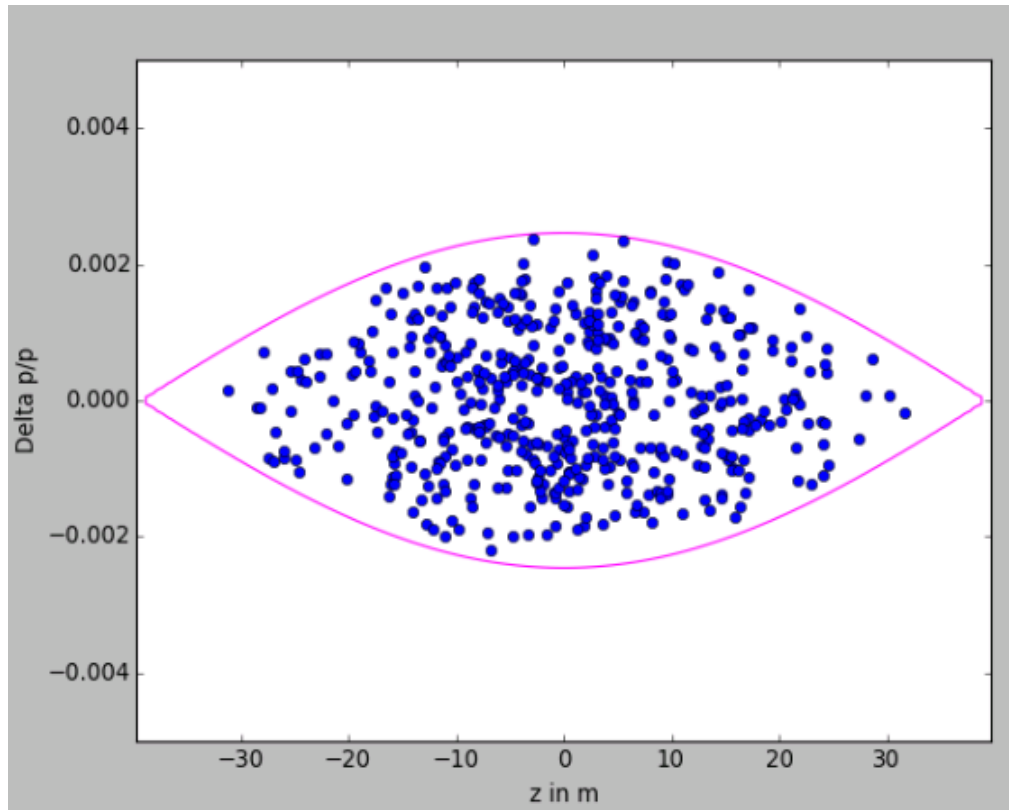
- Use a small number of particles ($n_{\text{particles}} \sim 10$) to observe the synchrotron motion at large and small amplitudes



→ 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

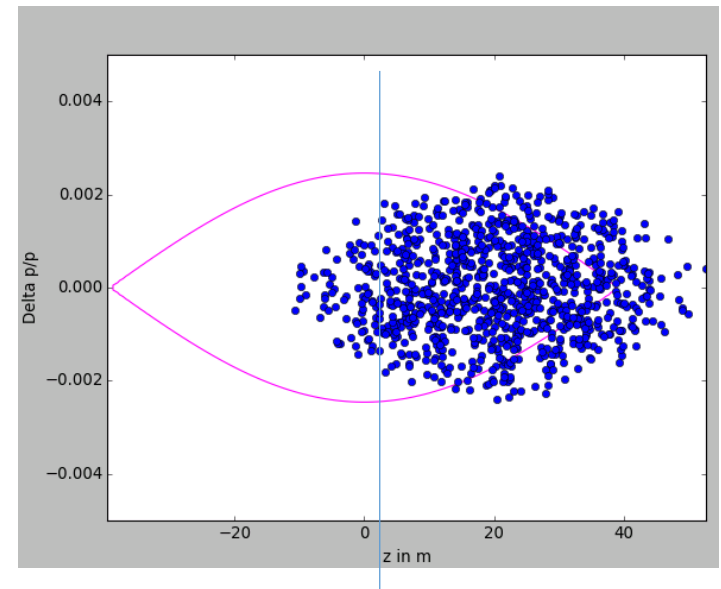
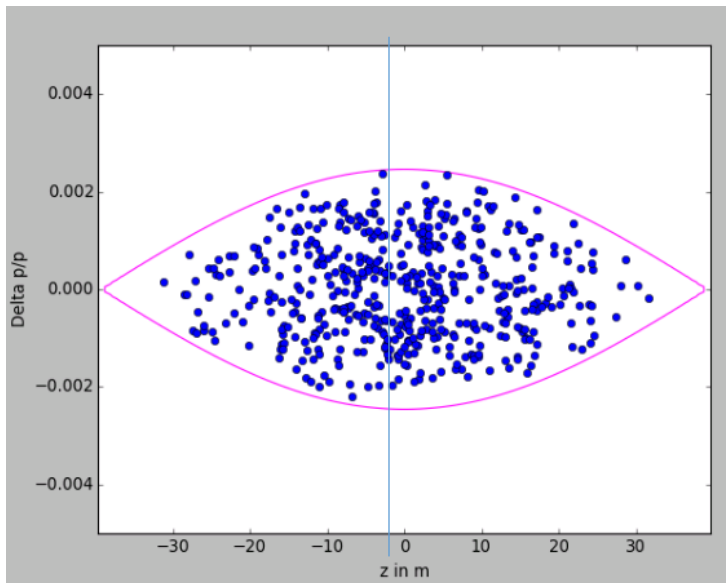
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 - Impact of voltage mismatch
 - Impact of acceleration
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Impact of phase mismatch

- 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
```

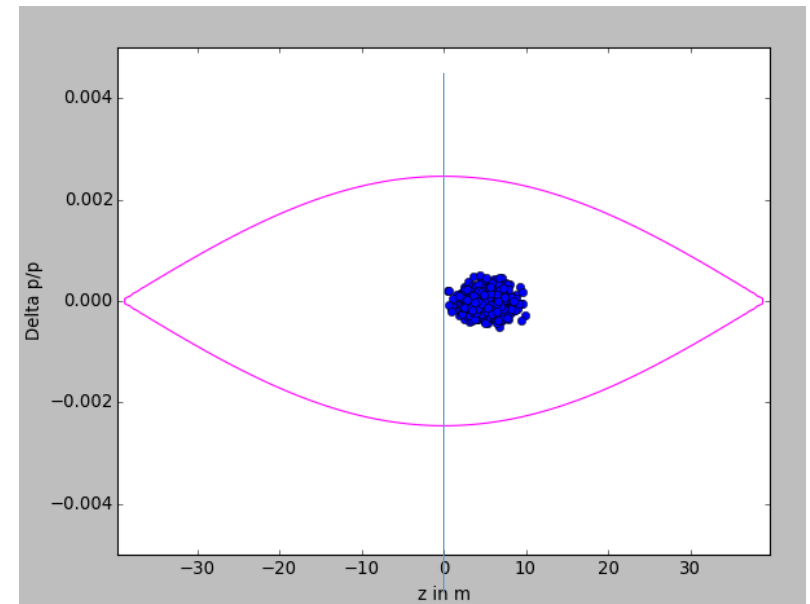
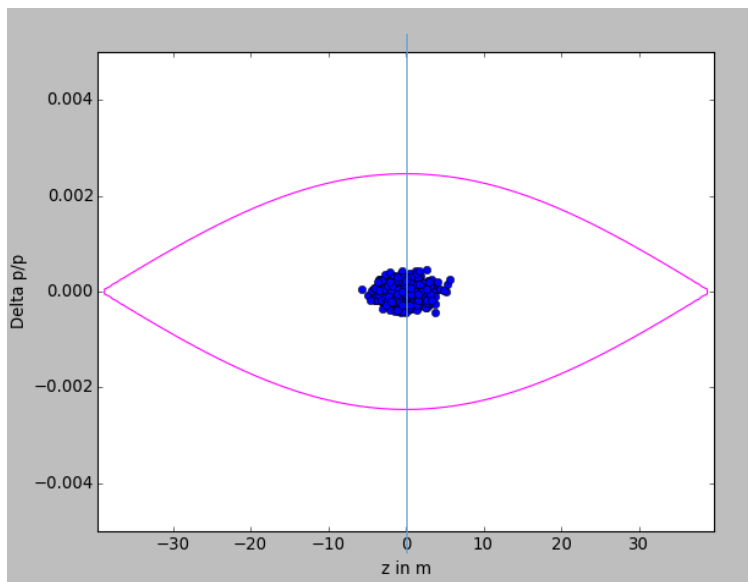
```
if i%100 == 0:
```



→ What will happen?

Impact of phase mismatch

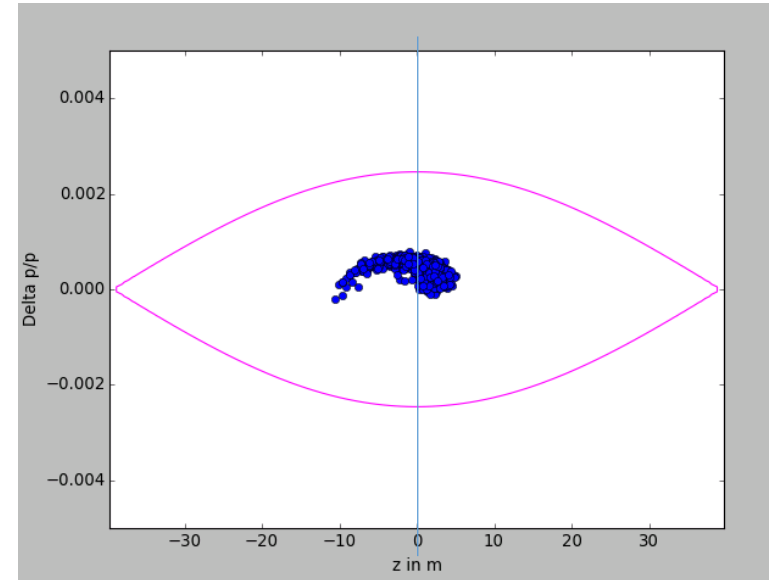
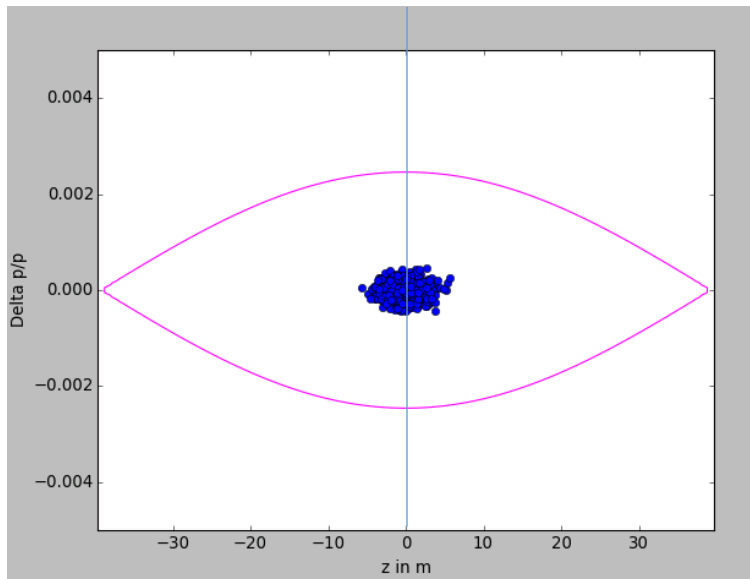
- Use a smaller bunch length (e.g. divide by 10)
- Generate ~ 1000 particles
- Add e.g. 5m to all particles: `bunch.z = bunch.z+5`
- Track particles for ~ 5000 turns
- Plot the distribution every 100 turns: `if i%100 == 0:`



→ What will happen?

Impact of phase mismatch

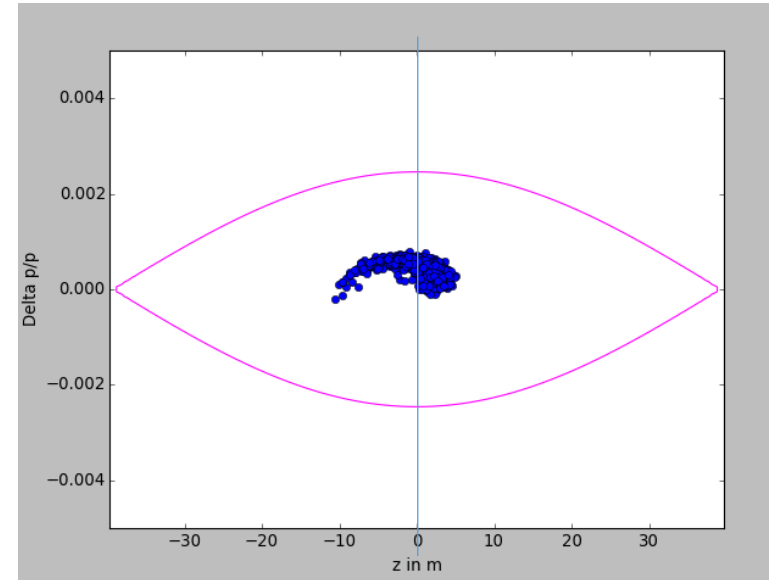
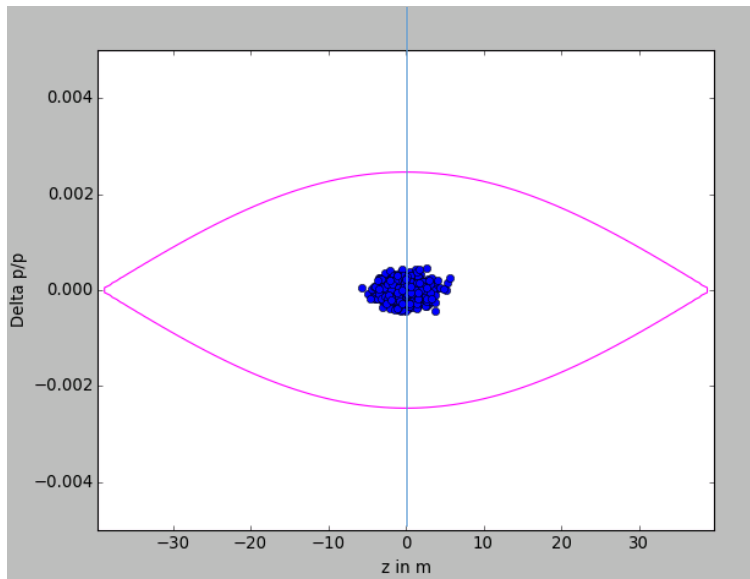
- Use a smaller bunch length (e.g. divide by 10)
- Use ~1000 particles
- Add e.g. 5m to all particles: `bunch.z = bunch.z+5`
- Track particles for ~5000 turns
- Plot the distribution every 100 turns: `if i%100 == 0:`



→ Filamentation and eventually?

Impact of phase mismatch

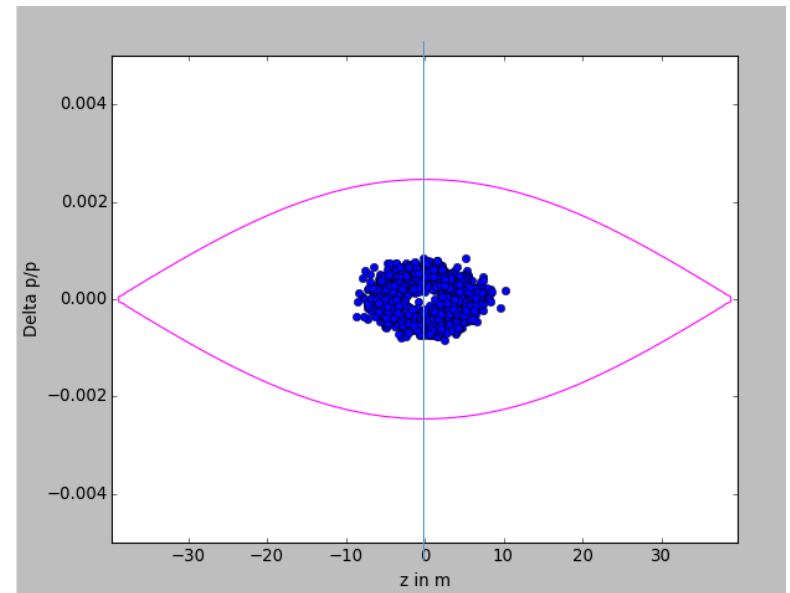
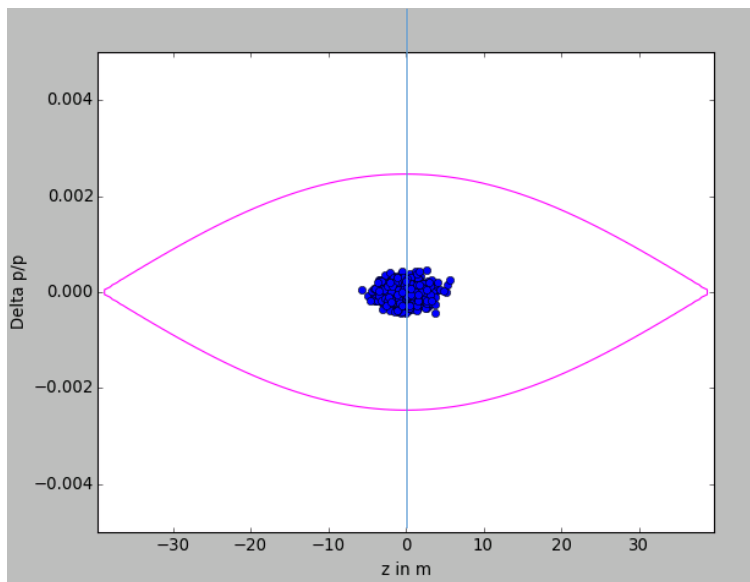
- Use a smaller bunch length (e.g. divide by 10)
- Use ~ 1000 particles
- Add e.g. 5m to all particles:
- Track particles for ~ 500000 turns
- Plot the distribution every 1000 turns



→ 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)

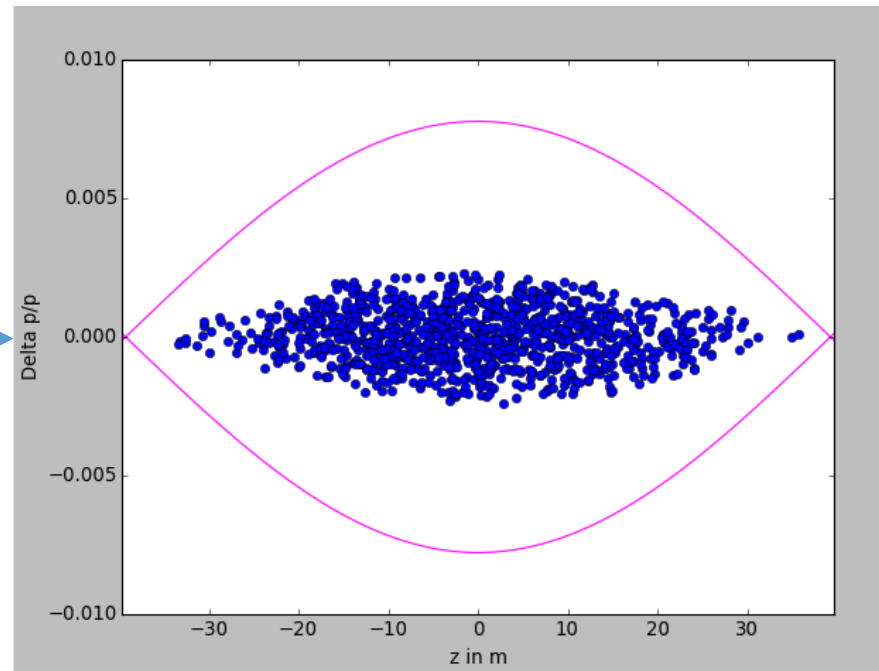
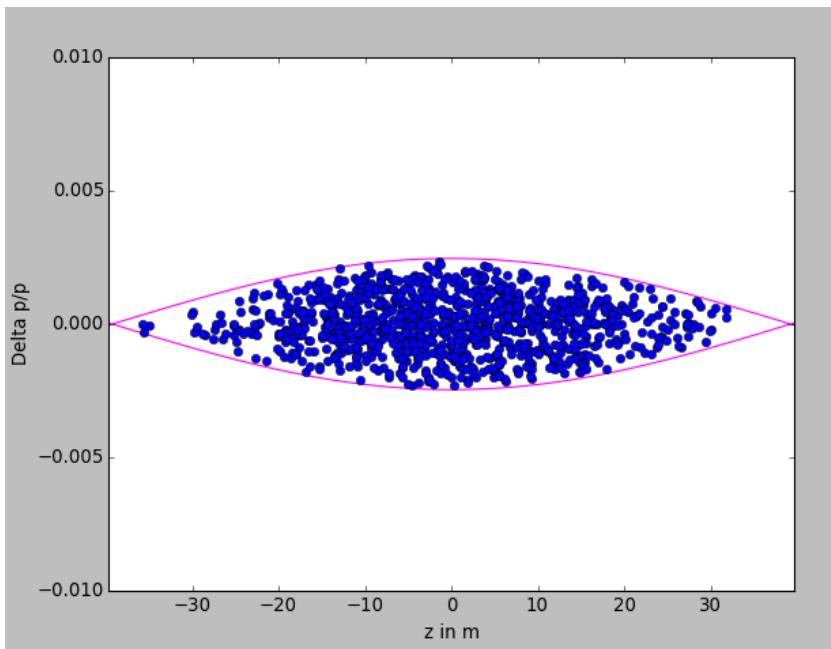
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Impact of voltage mismatch (too high)

- 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)
```

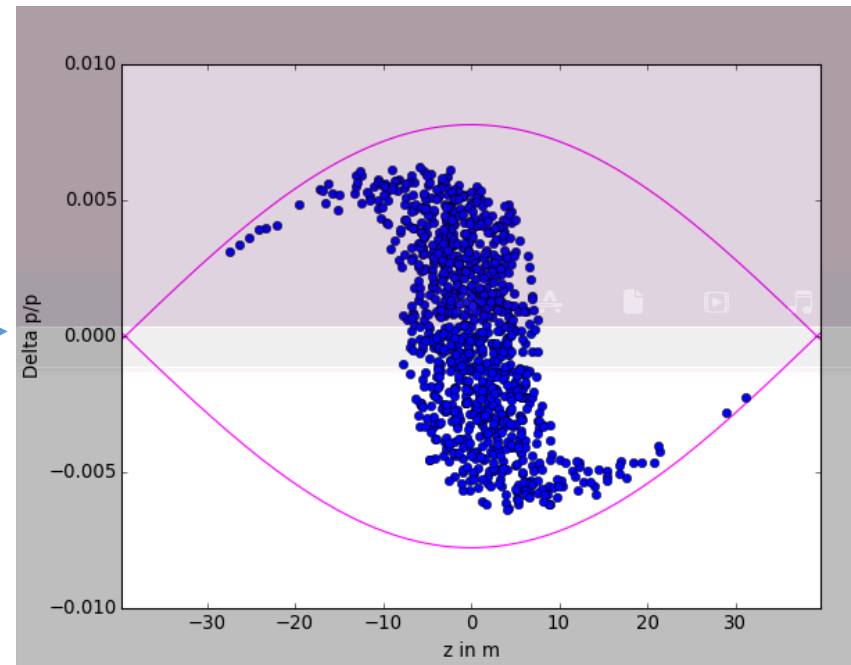
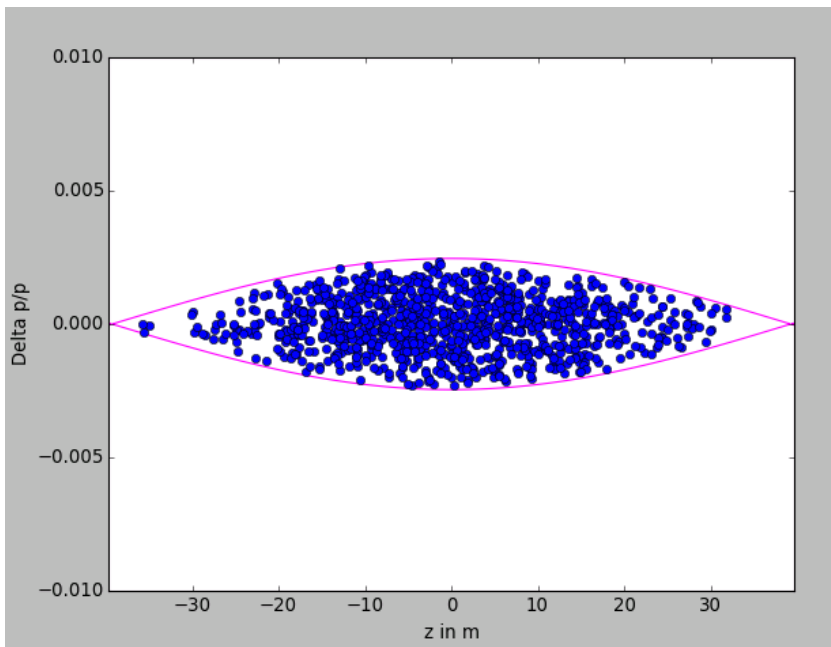


→ What will happen?

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],  
alpha_c_array, gamma, p_increment=p_increment_0)
```



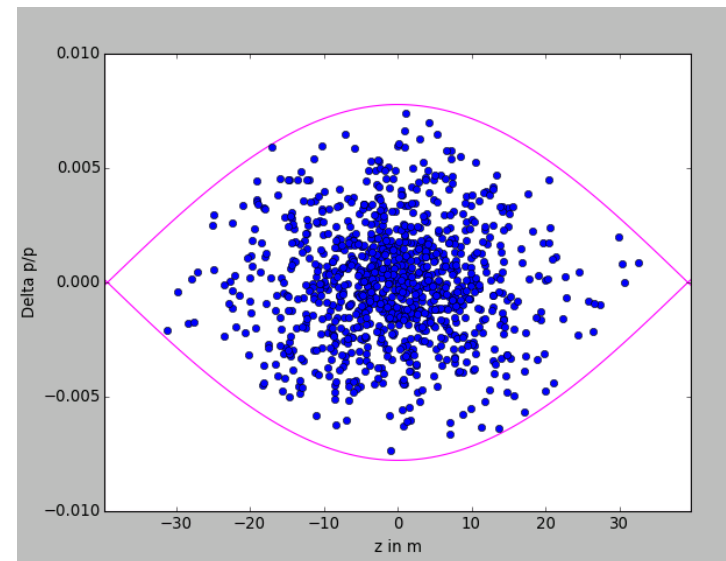
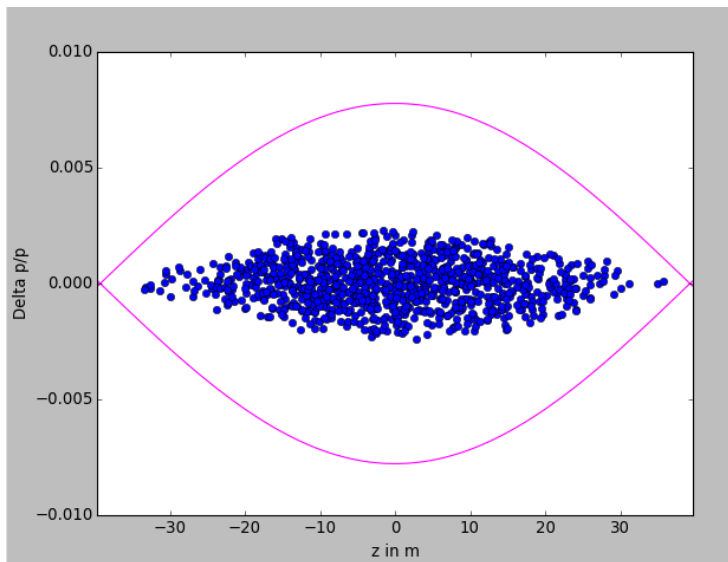
→ Bunch rotation

→ 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)
```

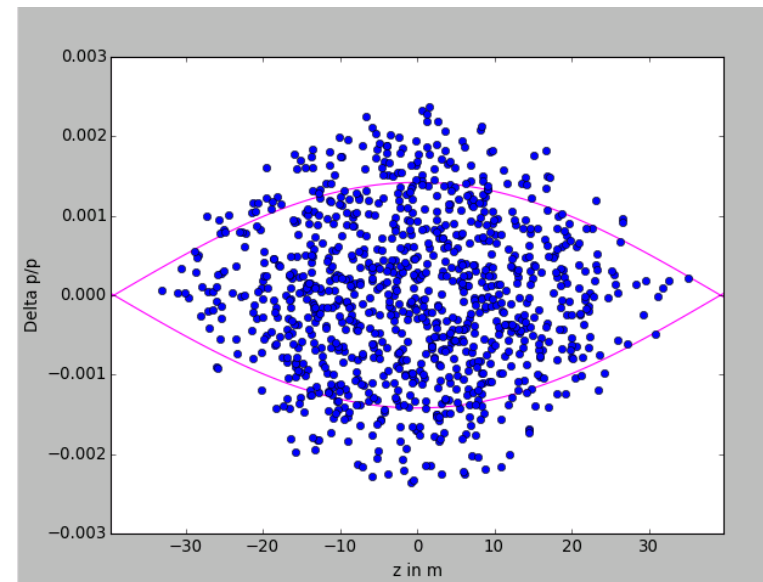
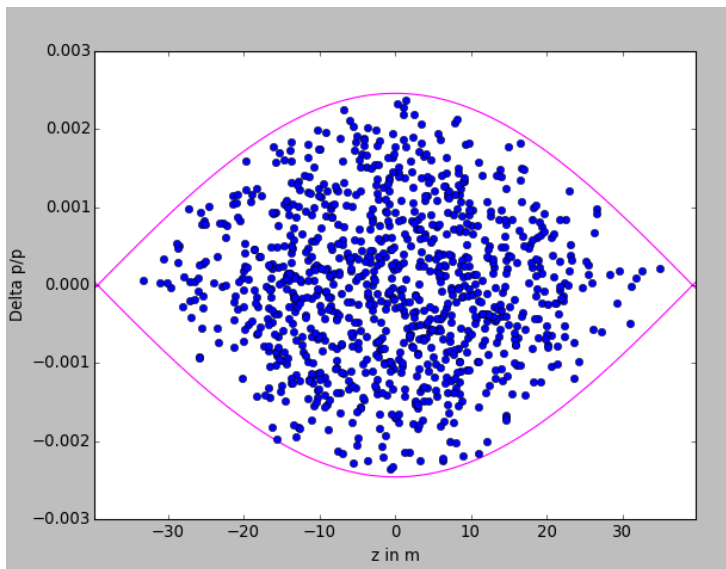


→ But should not wait too long!

→ Emittance growth!

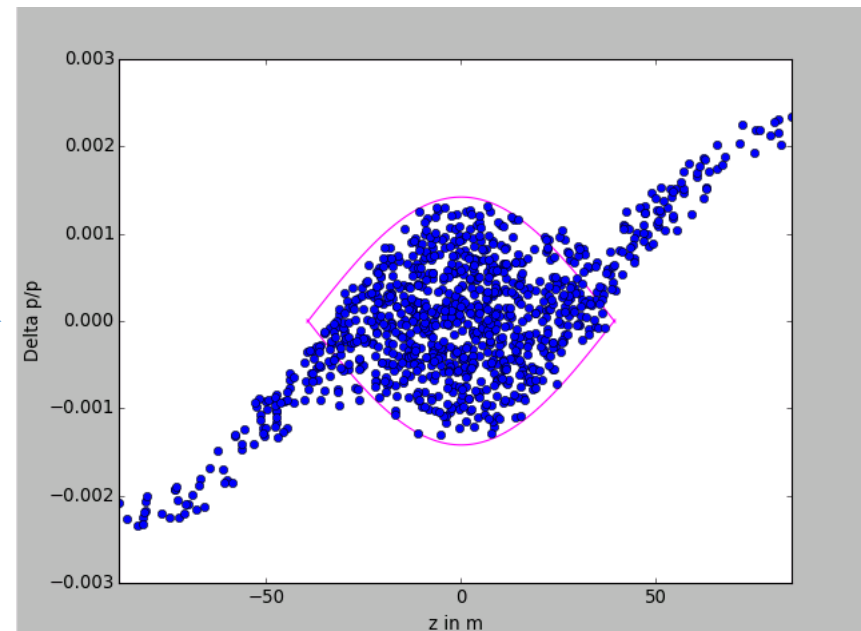
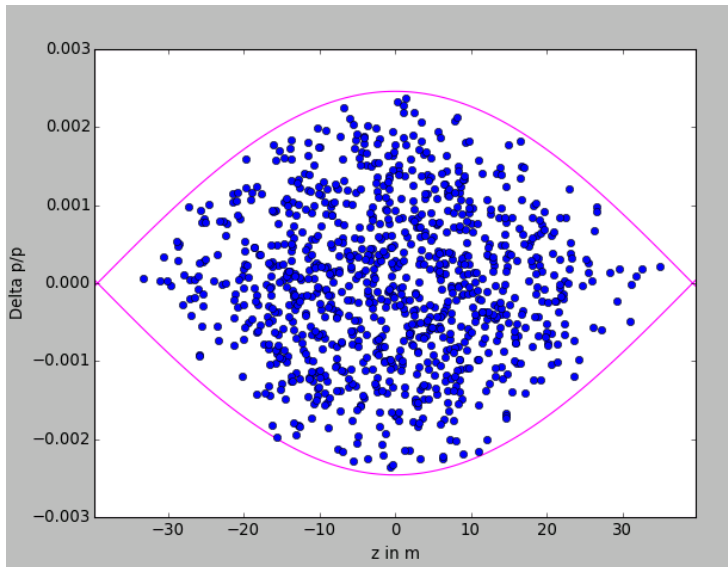
Impact of voltage mismatch (too low)

- 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

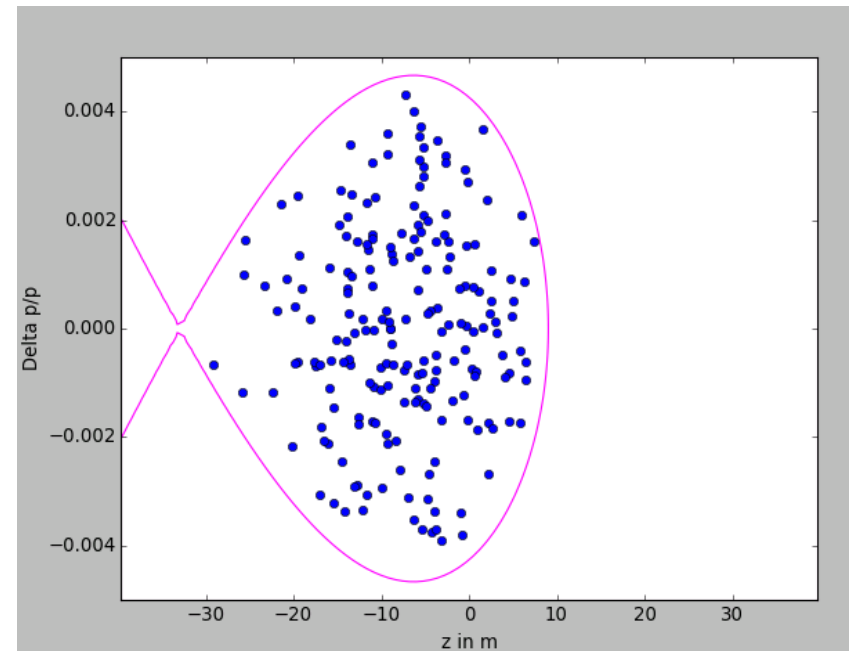
- Goals
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- Structure of ipython files
- **Tracking examples**
 - Impact of phase mismatch
 - Impact of voltage mismatch
 - **Impact of acceleration**
- Setting up the environment
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 - Ubuntu
 - Anaconda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



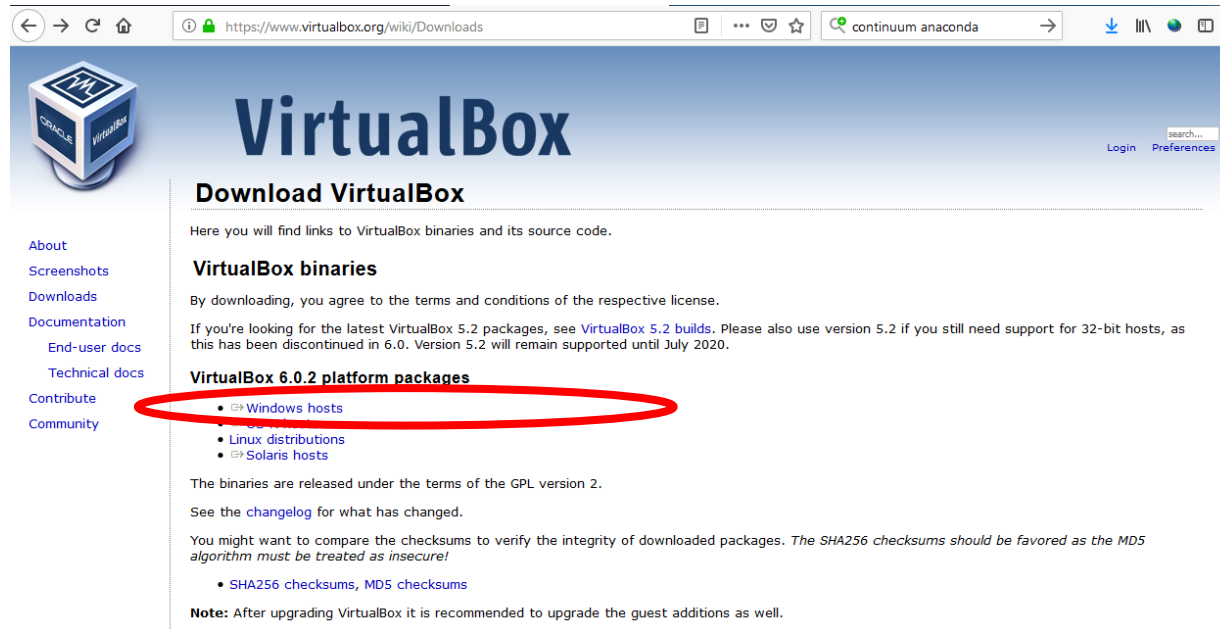
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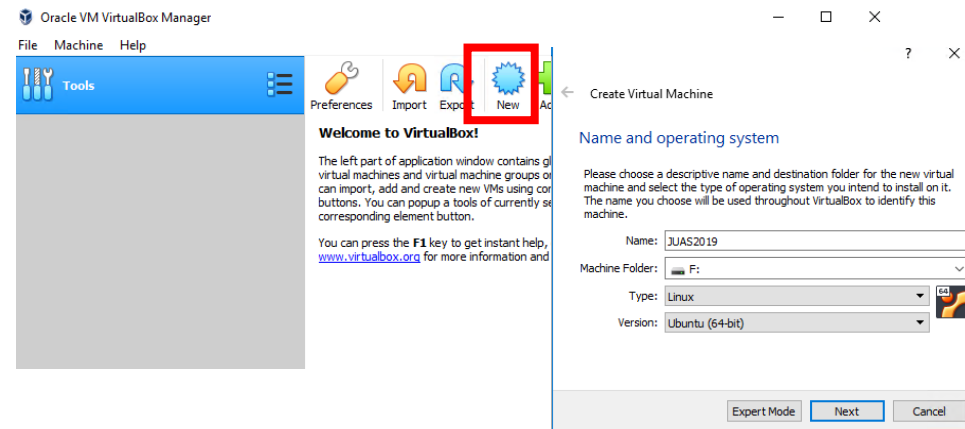
Setting up the operating system

- Install a virtual box to get an Ubuntu environment

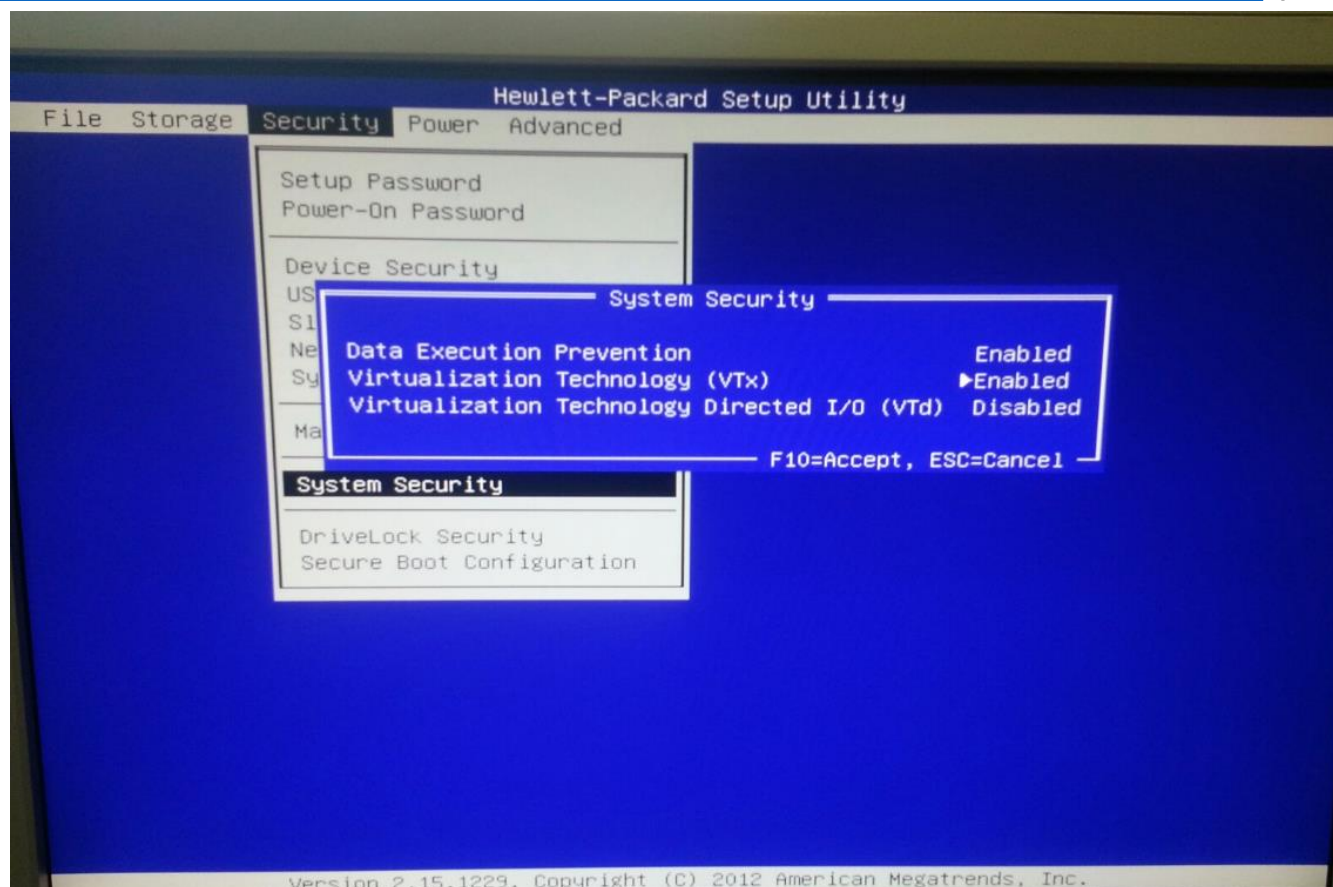
Ex: <https://www.virtualbox.org/wiki/Downloads>



- Create a new virtual machine with Ubuntu 64-bit



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

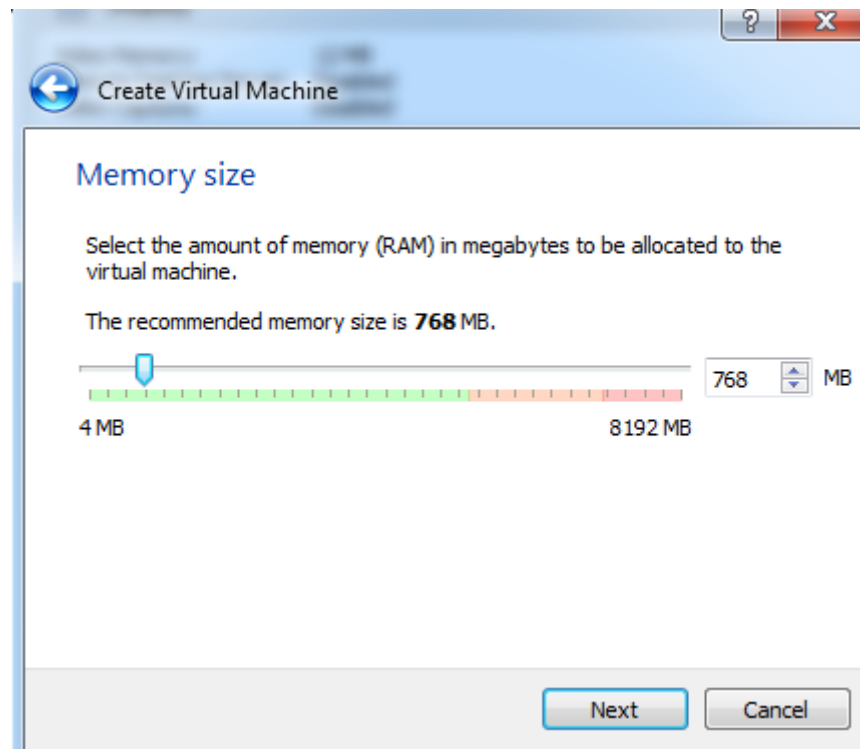
Please choose a descriptive name and destination folder for the new virtual machine and select the type of operating system you intend to install on it. The name you choose will be used throughout VirtualBox to identify this machine.

Name:

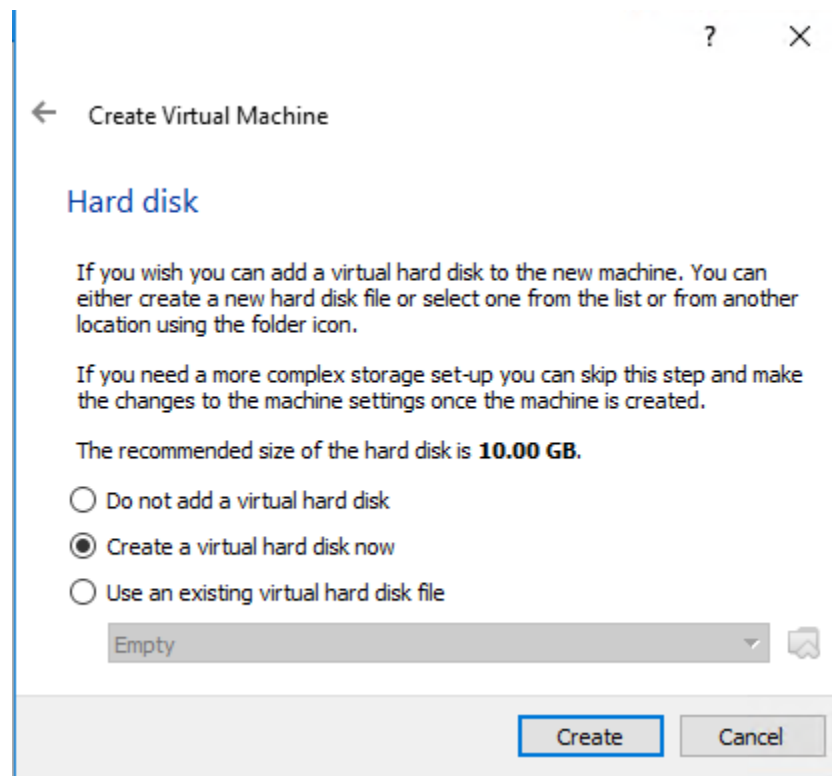
Machine Folder:

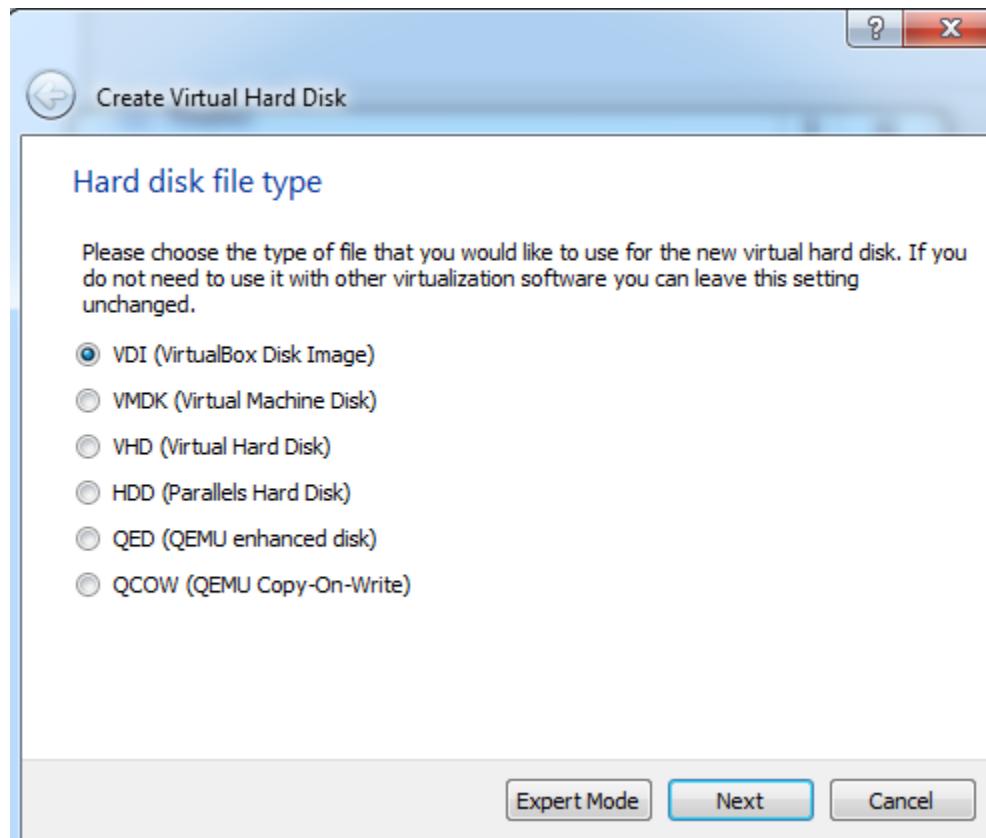
Type: 

Version:



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







← Create Virtual Hard Disk

Storage on physical hard disk

Please choose whether the new virtual hard disk file should grow as it is used (dynamically allocated) or if it should be created at its maximum size (fixed size).

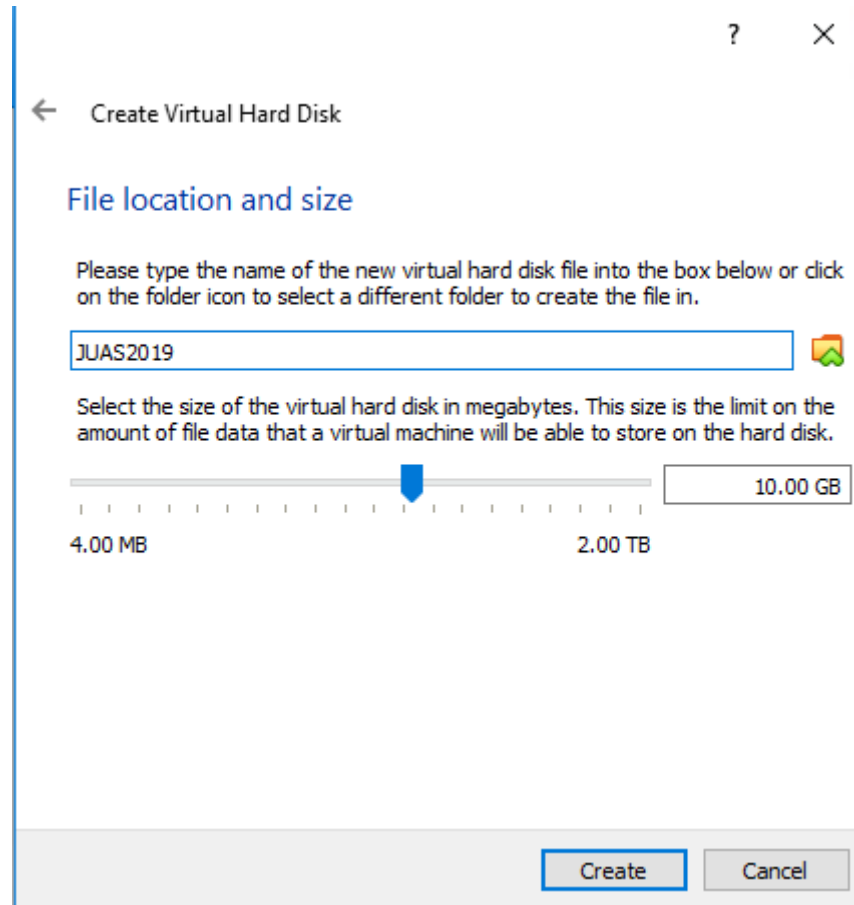
A **dynamically allocated** hard disk file will only use space on your physical hard disk as it fills up (up to a maximum **fixed size**), although it will not shrink again automatically when space on it is freed.

A **fixed size** hard disk file may take longer to create on some systems but is often faster to use.

- Dynamically allocated
- Fixed size

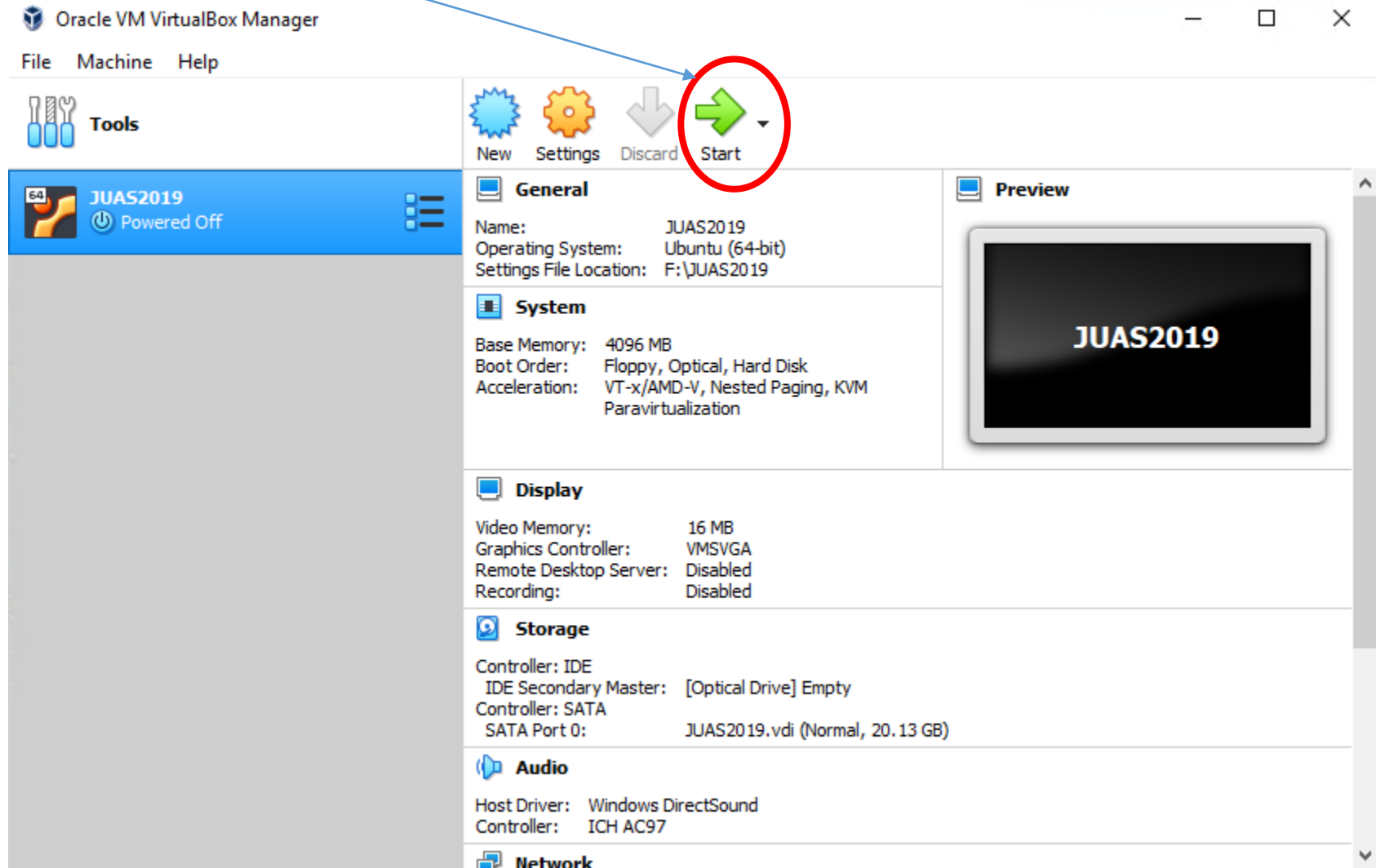
Next

Cancel



→ Assign 20 GB if possible.

Start the VM



The screenshot shows the Oracle VM VirtualBox Manager interface. The window title is "Oracle VM VirtualBox Manager". The menu bar includes "File", "Machine", and "Help". The "Tools" section contains icons for "New", "Settings", "Discard", and "Start". The "Start" button, represented by a green arrow, is circled in red. A blue arrow points from the "Start the VM" text to this button. The main area displays the configuration for a VM named "JUAS2019", which is currently "Powered Off". The configuration is organized into several sections: General, System, Display, Storage, Audio, and Network. The "Preview" window shows a black screen with the text "JUAS2019" in white.

Oracle VM VirtualBox Manager

File Machine Help

Tools

New Settings Discard Start

JUAS2019
Powered Off

General
Name: JUAS2019
Operating System: Ubuntu (64-bit)
Settings File Location: F:\JUAS2019

System
Base Memory: 4096 MB
Boot Order: Floppy, Optical, Hard Disk
Acceleration: VT-x/AMD-V, Nested Paging, KVM Paravirtualization

Display
Video Memory: 16 MB
Graphics Controller: VMSVGA
Remote Desktop Server: Disabled
Recording: Disabled

Storage
Controller: IDE
IDE Secondary Master: [Optical Drive] Empty
Controller: SATA
SATA Port 0: JUAS2019.vdi (Normal, 20.13 GB)

Audio
Host Driver: Windows DirectSound
Controller: ICH AC97

Network

Preview
JUAS2019

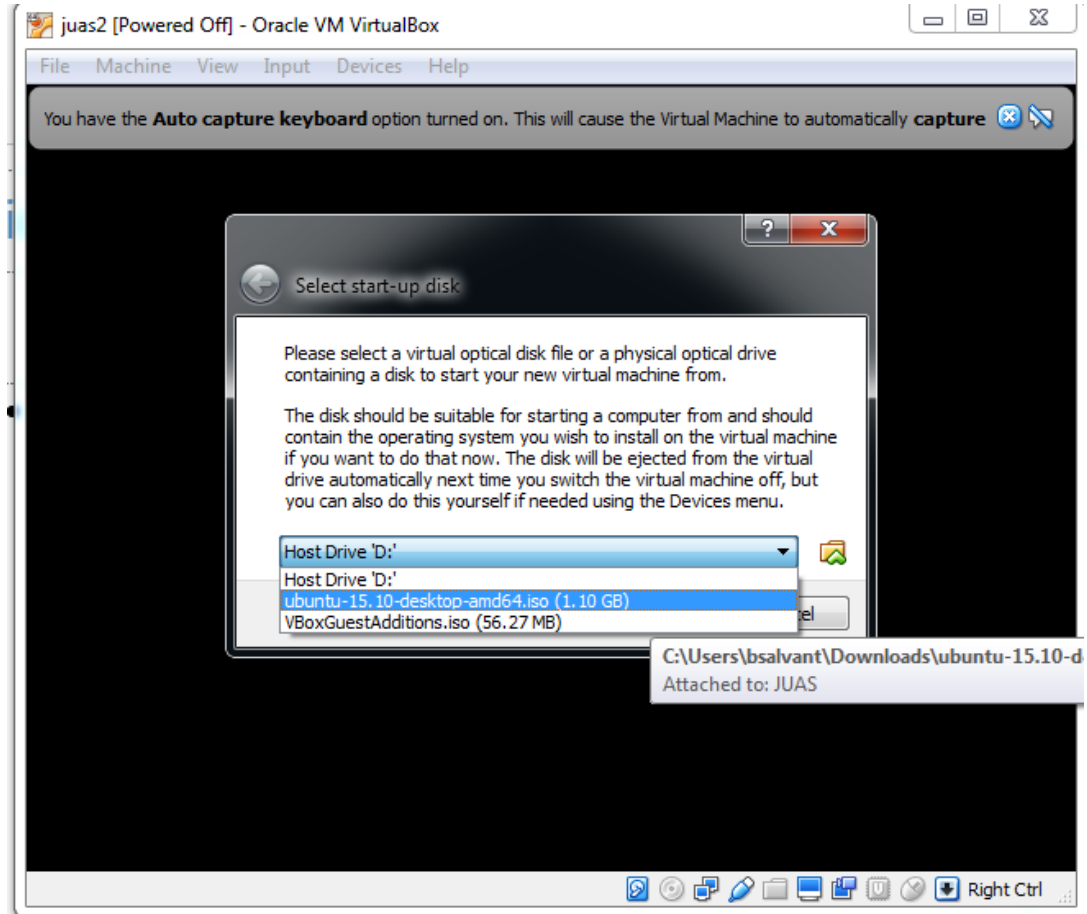
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Download ubuntu

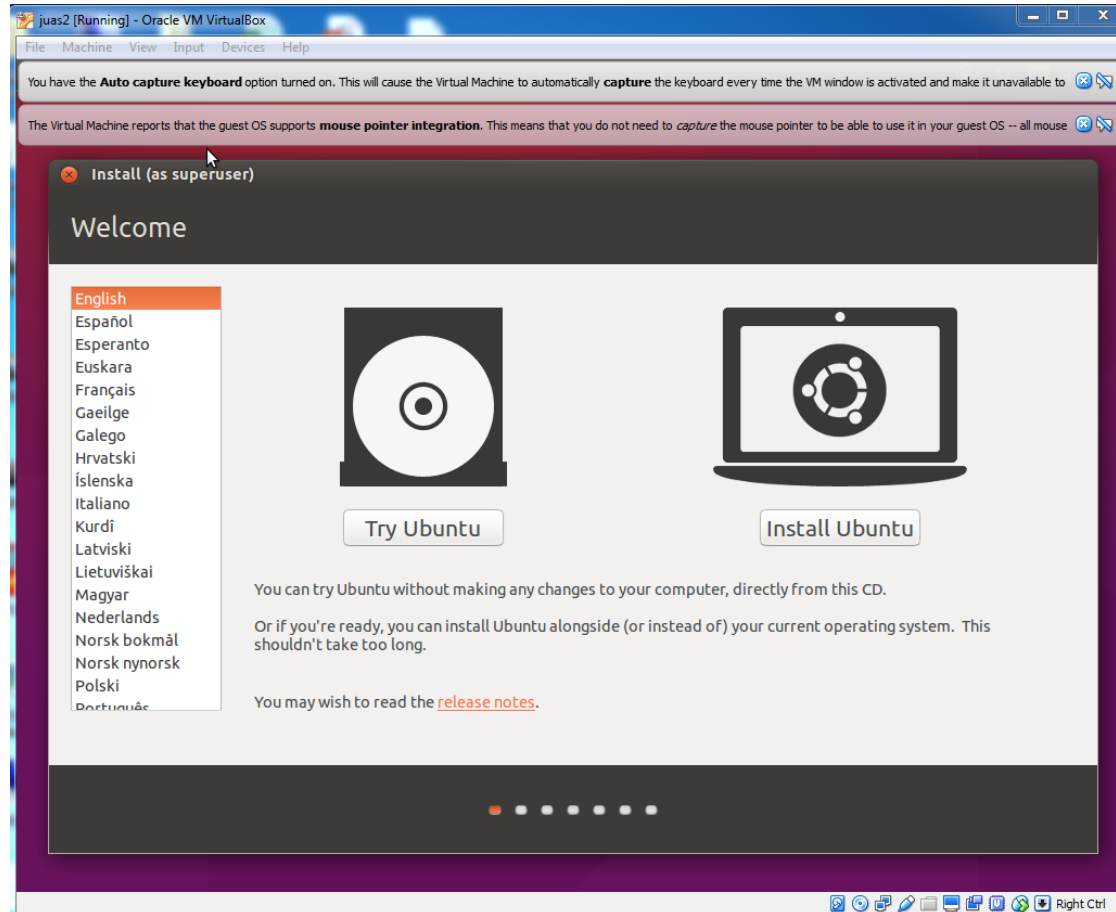
- <https://www.ubuntu.com/download/desktop>

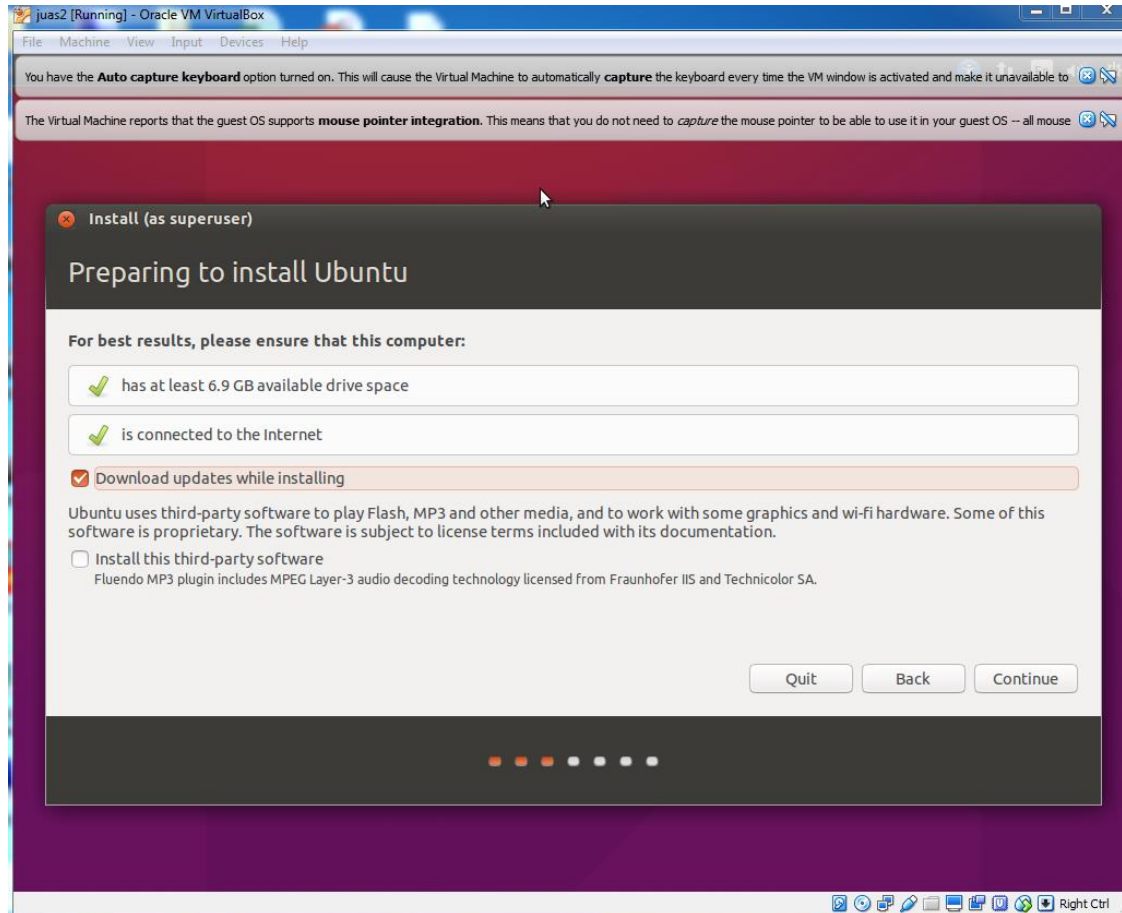
Select Ubuntu start up disk



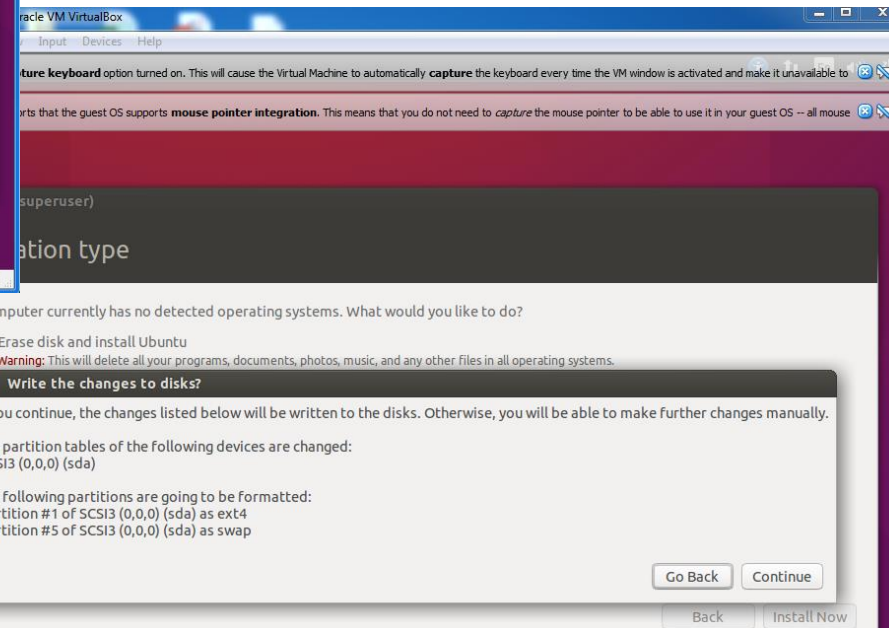
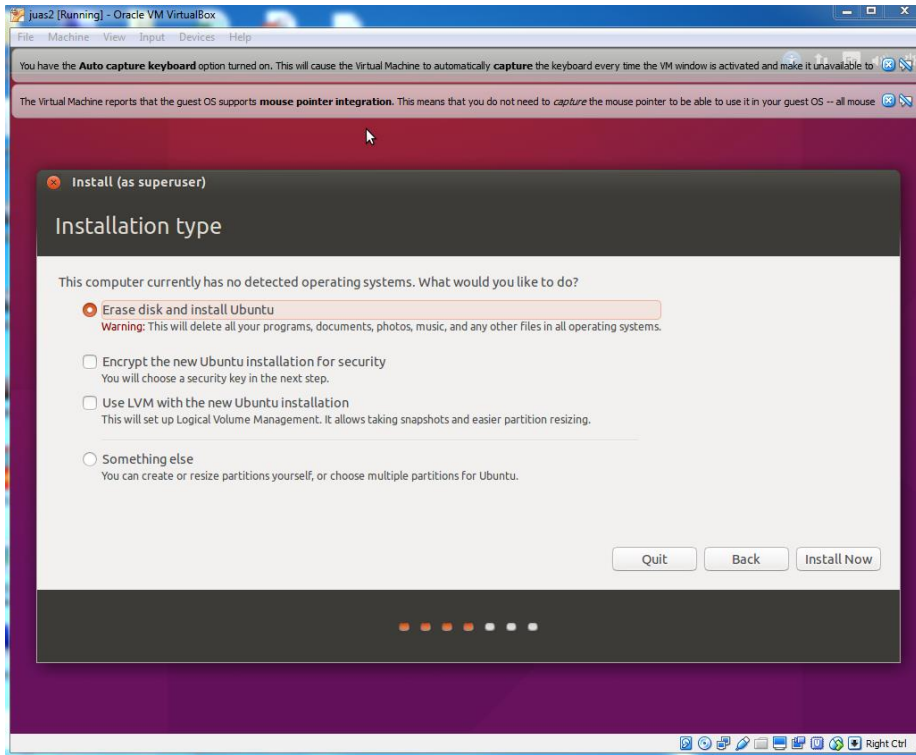
You may have to click on the small green arrow to browse for your Ubuntu installation image

Install ubuntu





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



Then choose time zone and login details

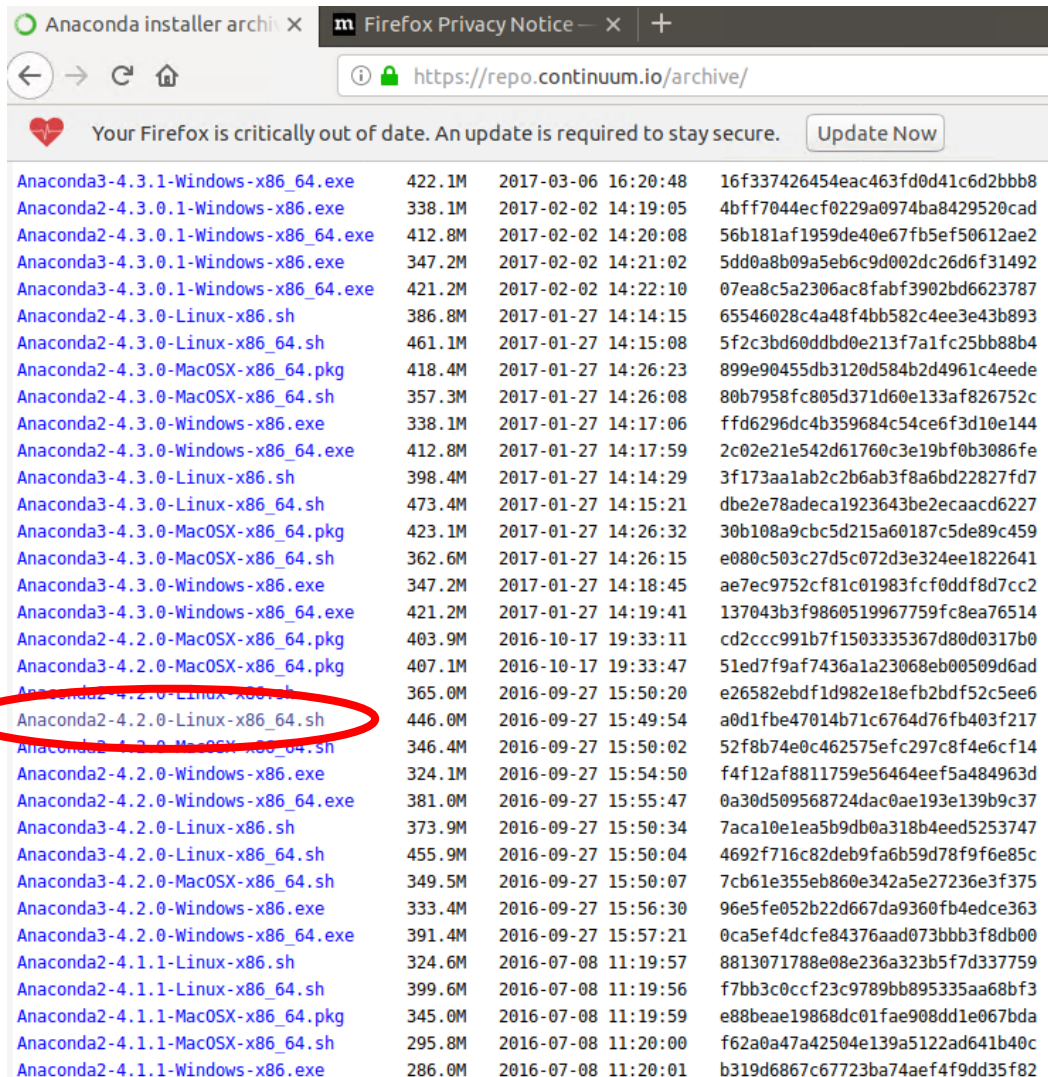
Then wait for file copy and Ubuntu installation restart the machine after it is complete

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Install anaconda

Go to <https://repo.continuum.io/archive/> and download the link “Anaconda2-4.2.0-Linux-x86_64.sh”



Anaconda3-4.3.1-Windows-x86_64.exe	422.1M	2017-03-06 16:20:48	16f337426454eac463fd0d41c6d2bbb8
Anaconda2-4.3.0.1-Windows-x86.exe	338.1M	2017-02-02 14:19:05	4bff7044ecf0229a0974ba8429520cad
Anaconda2-4.3.0.1-Windows-x86_64.exe	412.8M	2017-02-02 14:20:08	56b181af1959de40e67fb5ef50612ae2
Anaconda3-4.3.0.1-Windows-x86.exe	347.2M	2017-02-02 14:21:02	5dd0a8b09a5eb6c9d002dc26d6f31492
Anaconda3-4.3.0.1-Windows-x86_64.exe	421.2M	2017-02-02 14:22:10	07ea8c5a2306ac8fabf3902bd6623787
Anaconda2-4.3.0-Linux-x86.sh	386.8M	2017-01-27 14:14:15	65546028c4a48f4bb582c4ee3e43b893
Anaconda2-4.3.0-Linux-x86_64.sh	461.1M	2017-01-27 14:15:08	5f2c3bd60ddb0e213f7a1fc25bb88b4
Anaconda2-4.3.0-MacOSX-x86_64.pkg	418.4M	2017-01-27 14:26:23	899e90455db3120d584b2d4961c4eede
Anaconda2-4.3.0-MacOSX-x86_64.sh	357.3M	2017-01-27 14:26:08	80b7958fc805d371d60e133af826752c
Anaconda2-4.3.0-Windows-x86.exe	338.1M	2017-01-27 14:17:06	ffd6296dc4b359684c54ce6f3d10e144
Anaconda2-4.3.0-Windows-x86_64.exe	412.8M	2017-01-27 14:17:59	2c02e21e542d61760c3e19bf0b3086fe
Anaconda3-4.3.0-Linux-x86.sh	398.4M	2017-01-27 14:14:29	3f173aa1ab2c2b6ab3f8a6bd22827fd7
Anaconda3-4.3.0-Linux-x86_64.sh	473.4M	2017-01-27 14:15:21	dbe2e78adeca1923643be2ecaacd6227
Anaconda3-4.3.0-MacOSX-x86_64.pkg	423.1M	2017-01-27 14:26:32	30b108a9cbc5d215a60187c5de89c459
Anaconda3-4.3.0-MacOSX-x86_64.sh	362.6M	2017-01-27 14:26:15	e080c503c27d5c072d3e324ee1822641
Anaconda3-4.3.0-Windows-x86.exe	347.2M	2017-01-27 14:18:45	ae7ec9752cf81c01983fcf0ddf8d7cc2
Anaconda3-4.3.0-Windows-x86_64.exe	421.2M	2017-01-27 14:19:41	137043b3f9860519967759f9c8ea76514
Anaconda2-4.2.0-MacOSX-x86_64.pkg	403.9M	2016-10-17 19:33:11	cd2ccc991b7f1503335367d80d0317b0
Anaconda3-4.2.0-MacOSX-x86_64.pkg	407.1M	2016-10-17 19:33:47	51ed7f9af7436a1a23068eb00509d6ad
Anaconda2-4.2.0-Linux-x86.sh	365.0M	2016-09-27 15:50:20	e26582ebddf1d982e18efb2bdf52c5ee6
Anaconda2-4.2.0-Linux-x86_64.sh	446.0M	2016-09-27 15:49:54	a0d1fbc47014b71c6764d76fb403f217
Anaconda2-4.2.0-MacOSX-x86_64.sh	346.4M	2016-09-27 15:50:02	52f8b74e0c462575efc297c8f4e6cf14
Anaconda2-4.2.0-Windows-x86.exe	324.1M	2016-09-27 15:54:50	f4f12af8811759e56464ee5a484963d
Anaconda2-4.2.0-Windows-x86_64.exe	381.0M	2016-09-27 15:55:47	0a30d509568724dac0ae193e139b9c37
Anaconda3-4.2.0-Linux-x86.sh	373.9M	2016-09-27 15:50:34	7aca10e1ea5b9db0a318b4eed5253747
Anaconda3-4.2.0-Linux-x86_64.sh	455.9M	2016-09-27 15:50:04	4692f716c82deb9fa6b59d78f9f6e85c
Anaconda3-4.2.0-MacOSX-x86_64.sh	349.5M	2016-09-27 15:50:07	7cb61e355eb860e342a5e27236e3f375
Anaconda3-4.2.0-Windows-x86.exe	333.4M	2016-09-27 15:56:30	96e5fe052b22d667da9360fb4edce363
Anaconda3-4.2.0-Windows-x86_64.exe	391.4M	2016-09-27 15:57:21	0ca5ef4dcfe84376aad073bbb3f8db00
Anaconda2-4.1.1-Linux-x86.sh	324.6M	2016-07-08 11:19:57	8813071788e08e236a323b5f7d337759
Anaconda2-4.1.1-Linux-x86_64.sh	399.6M	2016-07-08 11:19:56	f7bb3c0ccf23c9789bb895335aa68bf3
Anaconda2-4.1.1-MacOSX-x86_64.pkg	345.0M	2016-07-08 11:19:59	e88beae19868dc01fae908dd1e067bda
Anaconda2-4.1.1-MacOSX-x86_64.sh	295.8M	2016-07-08 11:20:00	f62a0a47a42504e139a5122ad641b40c
Anaconda2-4.1.1-Windows-x86.exe	286.0M	2016-07-08 11:20:01	b319d6867c67723ba74aef4f9dd35f82

→ Then install it by typing

```
bash Anaconda2-4.2.0-Linux-x86_64.sh
```

In the download directory

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 2.7 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`)

In case of issues with PyHEADTAIL installation, most likely it is that fortran and C need to be installed:

- `sudo apt-get install gfortran`
- `Sudo apt-get install gcc`

Finding the examples

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

The screenshot shows a web browser window displaying the Indico timetable for the event 'JUAS - 2019 (7 January 2019 - 10 X)'. The URL is <https://indico.cern.ch/event/779575/timetable/>. The date is Tuesday, 22 January. The timetable is as follows:

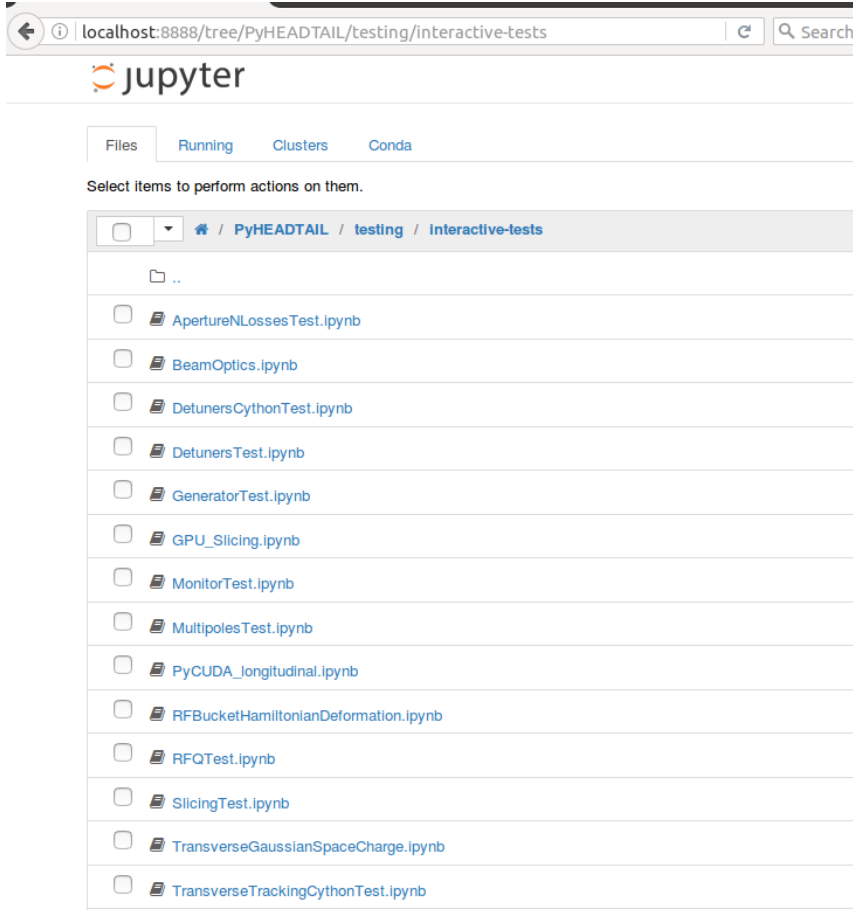
Time	Activity	Speakers	Duration
09:00 → 10:00	Longitudinal Dynamics	Benoit Salvant (CERN), Elias Metral (CERN - BE/ABP)	1h
10:00 → 10:15	Coffee break		15m
10:15 → 11:15	Longitudinal Dynamics	Benoit Salvant (CERN), Elias Metral (CERN)	1h
11:15 → 12:15	Longitudinal Dynamics	Benoit Salvant (CERN), Elias Metral (CERN - BE/ABP)	1h
12:15 → 14:00	Lunch		1h 45m
14:00 → 15:00	Longitudinal Dynamics	Benoit Salvant (CERN), Elias Metral (CERN - BE/ABP)	1h

The files for the 14:00-15:00 session are circled in red:

- 1_JUAS2017_injecti...
- 2_JUAS2017_injecti...
- 3_JUAS2017_injecti...
- 4_JUAS2017_injecti...
- 4_JUAS2017_injecti...

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

Other examples available with the PyHEADTAIL install



The screenshot shows a web browser window with the address bar displaying `localhost:8888/tree/PyHEADTAIL/testing/interactive-tests`. The Jupyter logo is visible at the top. Below the logo, there are tabs for `Files`, `Running`, `Clusters`, and `Conda`. The `Files` tab is active, and the page displays a file browser interface. The breadcrumb path is `PyHEADTAIL / testing / Interactive-tests`. Below the breadcrumb, there is a list of files, each with a checkbox and a file icon:

- ..
- 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

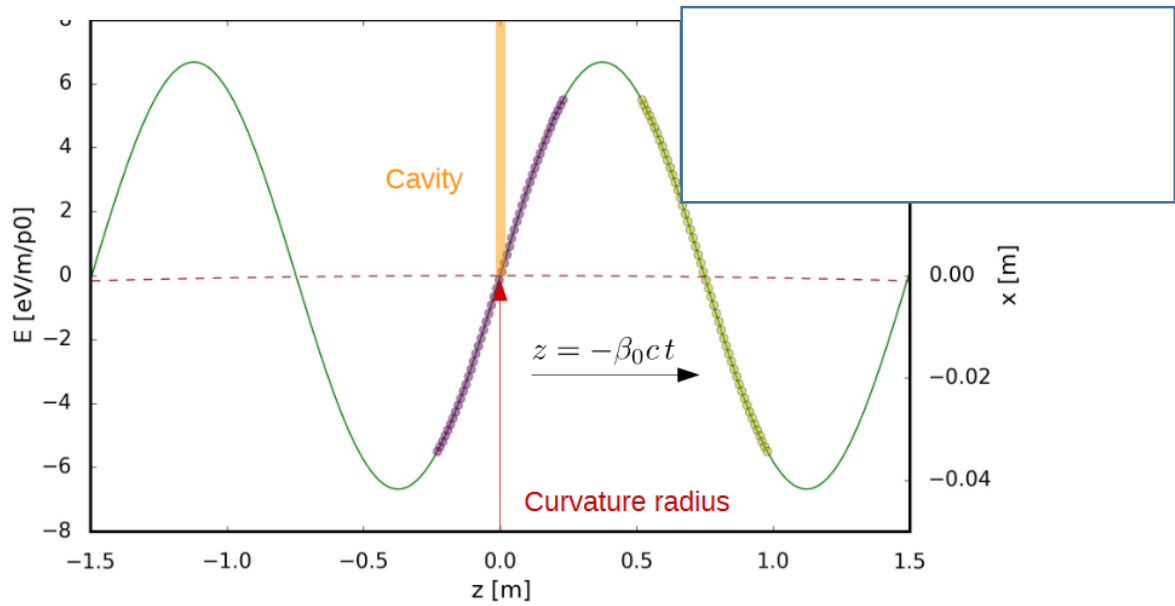
Johns Hopkins Accelerator School

- ◆ 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

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