

# Update (29/05/2012)

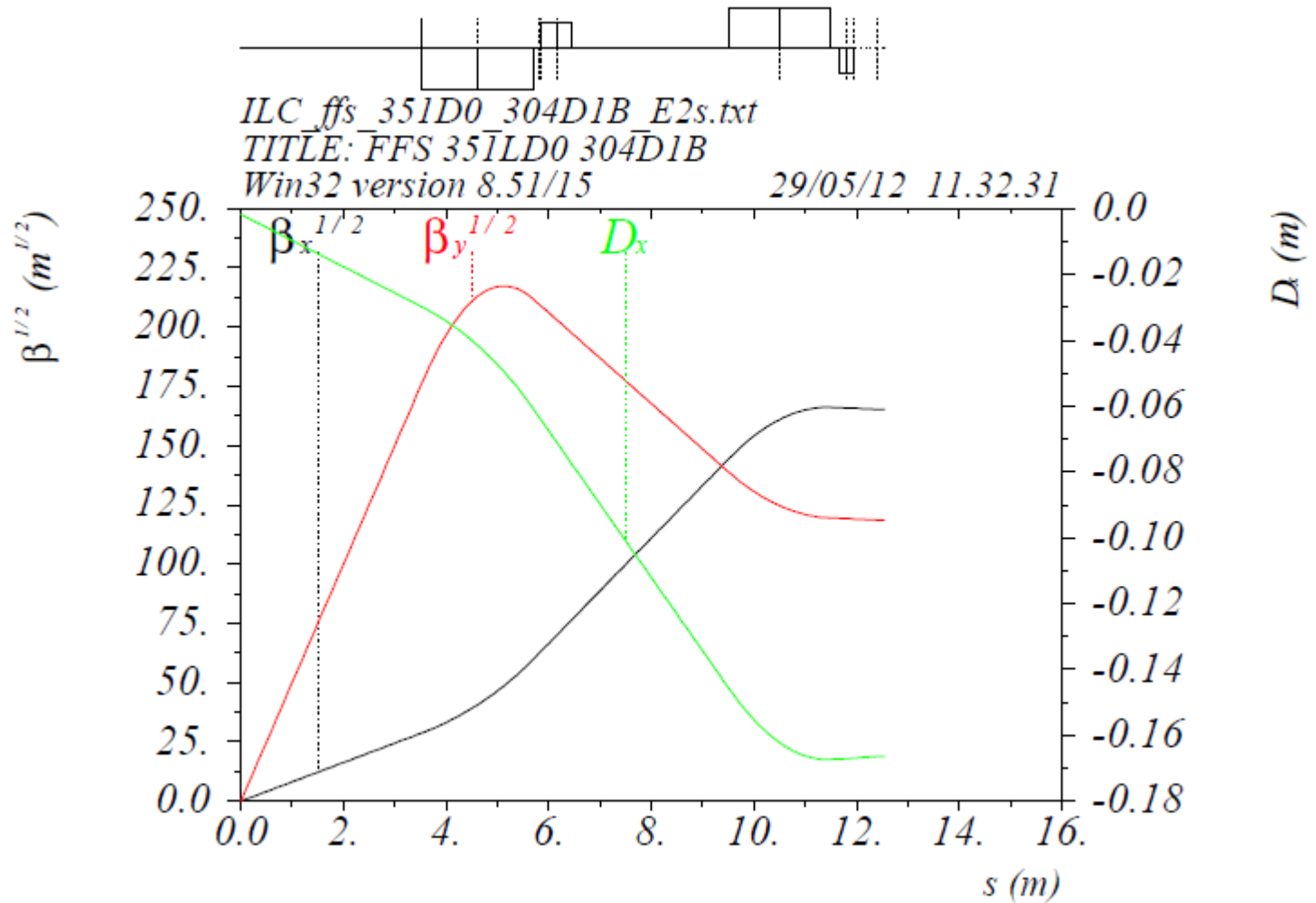
- MAD-X for Windows XP in Oxford – can run it now ( some editing of files was required)
- MAD-X via SSH in DESY/CERN is running but..
- I can use converted BDS lattice to first step matching
- Take 250 GeV CoM settings the task is to:
- match to the new  $\beta_x = 13\text{mm}$  and  $\beta_y = 0.41\text{mm}$
- and remove dispersion which is non=zero in the present lattice –should vary quadrupoles downstream
- $KQF3 := 0.011524797231;$        $kqf3 = 0.01147786643;$
- $KQD0 := -0.15229179766;$        $kqd0 = -0.1522525301;$
- $KQF1 := 0.07234909233;$        $kqf1 = 0.07254916217;$
- $KQF5 := 0.011062408721;$        $kqf5 = 0.01097346172;$
- $KQD4 := -0.858256166518E-2;$        $kqd4 = -0.008605842666;$

$$\beta_x = 0.01299999482; \quad \beta_y = 0.0004105255407; \quad \alpha_x = 3.852111358e-010;$$
$$\alpha_y = 3.623848278e-010; \quad Dx = -1.285273381e-010$$

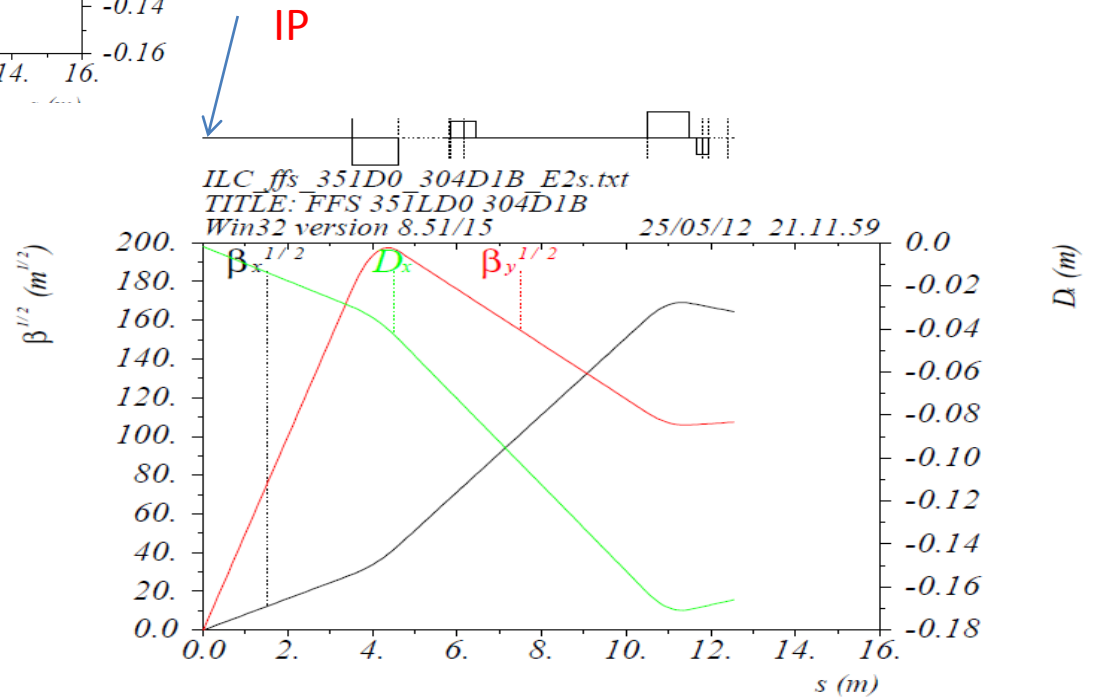
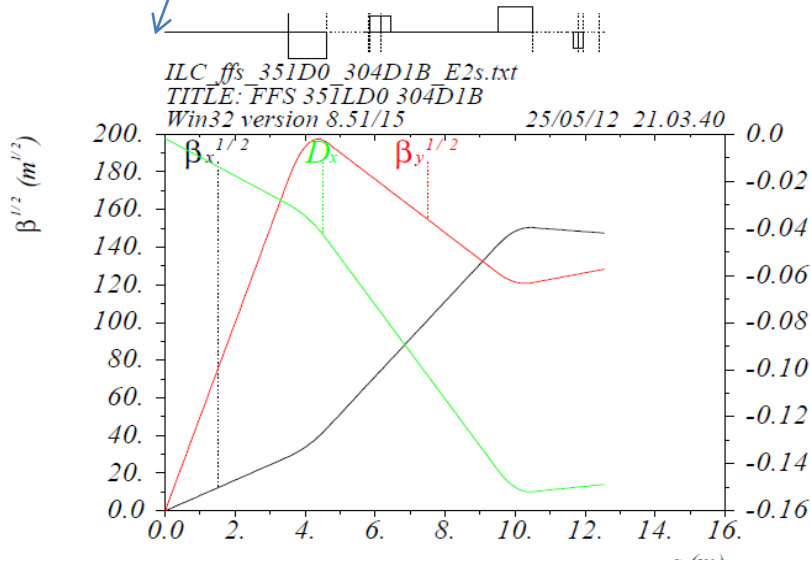
# Update(cont.)

- When I use the full matching program ( with call to Python)from Rogelio, some error message appears:
- *file "getSigmasMixed25.py", line 10, in ?  
from mapclassGaussianDelta25 import \*  
File  
"/afs/desy.de/user/m/malyshev/private/ILC\_BDS/RepILCcopy/mapclassGaussianDelta25.py", line 4, in ? from numpy  
import \*  
ImportError: No module named numpy*
- *Nevertheless*
- `./getSigmasMixed25.py` is running and I got a list of sigma values. Thus some mistake in redirection to CERN afs happens somewhere.

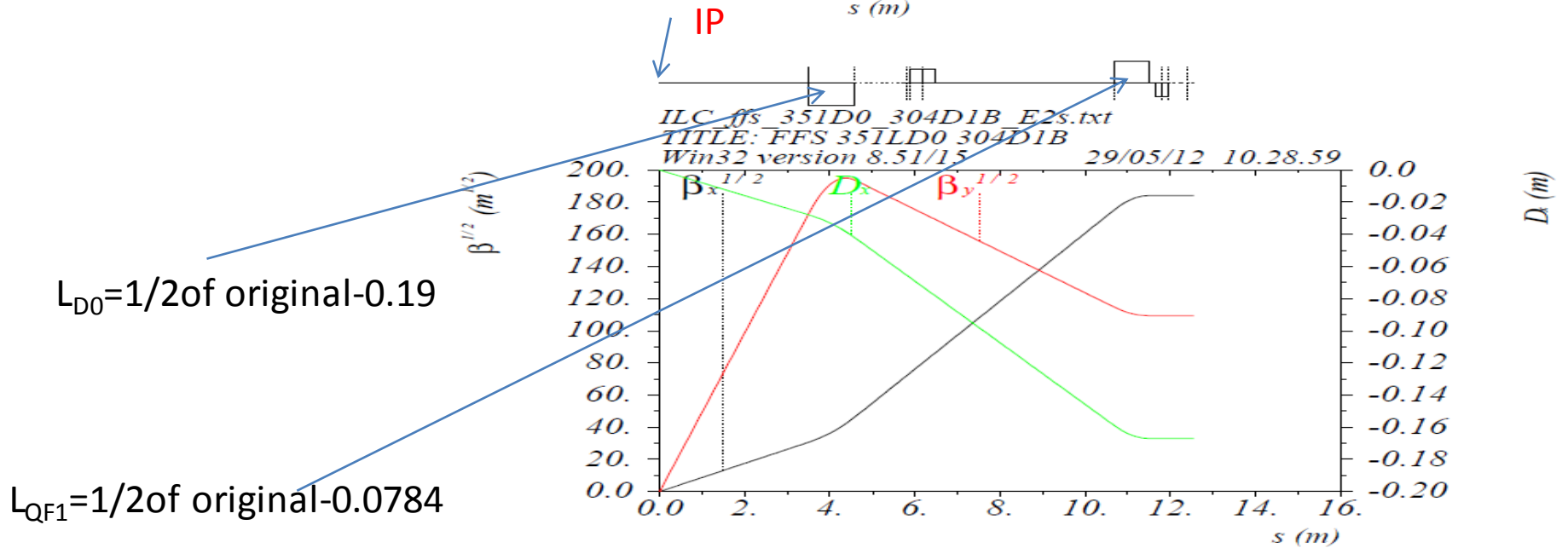
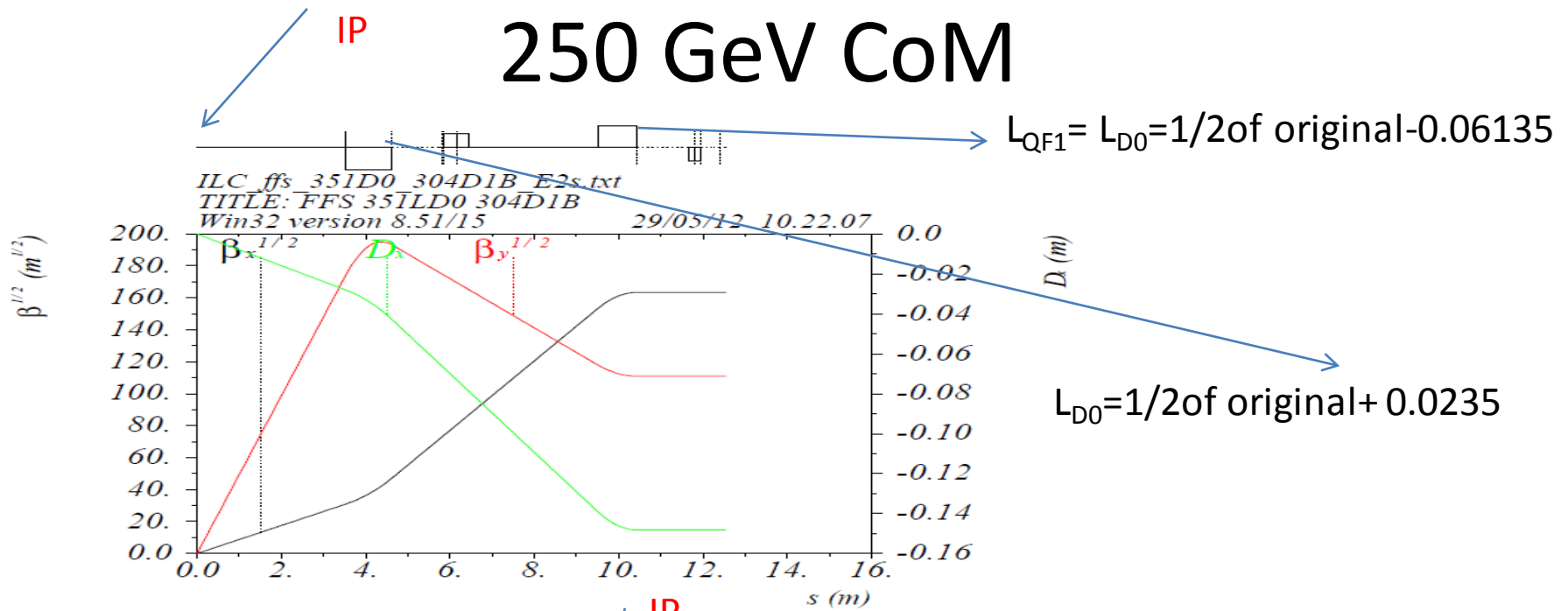
# 500 GeV CoM final doublet



# 250 GeV CoM (unmatched)



# 250 GeV CoM



# Plans

- Get MAD-X matching program running (update Python in DESY or/and short visit to CERN?)
- 250 GeV CoM lattice optimisation over range of parameters