



Matrix-Element Method Tutorial: MadWeight

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- Instruction
 - ➔ Theoretical reminder
 - ➔ description of the exercise
- Installation
- Tutorial
- Conclusion

- Associate to each experimental event characterised by \mathbf{p}^{vis} , the **probability** $\mathcal{P}(\mathbf{p}^{vis}|\alpha)$ to be produced and observed following a theoretical assumption α

$$\mathcal{P}(\mathbf{p}^{vis}|\alpha) = \frac{1}{\sigma_{\alpha}^{vis}} \int d\Phi dx_1 dx_2 |M_{\alpha}(\mathbf{p})|^2 W(\mathbf{p}, \mathbf{p}^{vis})$$

- $|M_{\alpha}(\mathbf{p})|^2$ is the squared matrix element
- $W(\mathbf{p}, \mathbf{p}^{vis})$ is the transfer function
- $\int d\Phi dx_1 dx_2$ is the phase-space integral
- σ_{α}^{vis} is the cross-section (after cuts)

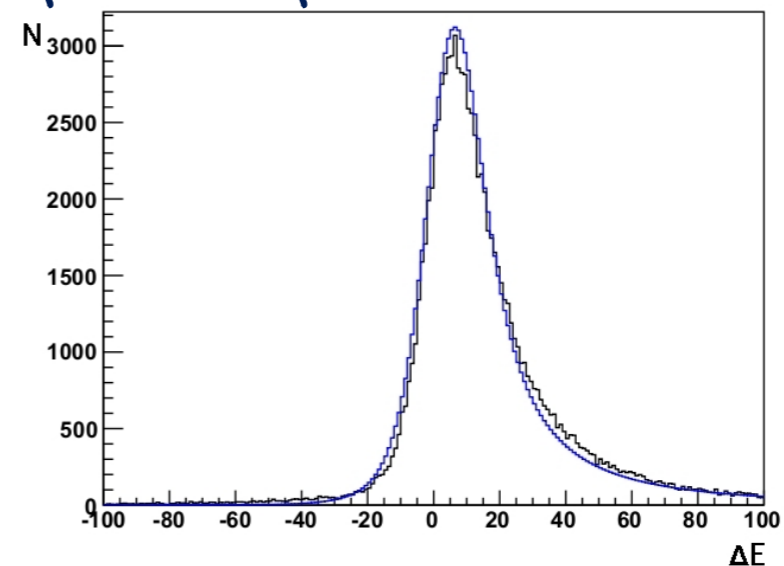
$$\mathcal{P}(\mathbf{p}^{vis}|\alpha) = \frac{1}{\sigma_\alpha} \int d\Phi dx_1 dx_2 |M_\alpha(\mathbf{p})|^2 W(\mathbf{p}, \mathbf{p}^{vis})$$

Four Elements:

- cross-section
- matrix-element
- transfer function
- integration

Computed via

- MadGraph5
- MadGraph5
- fitted from MC



- Madweight


MadWeight returns:

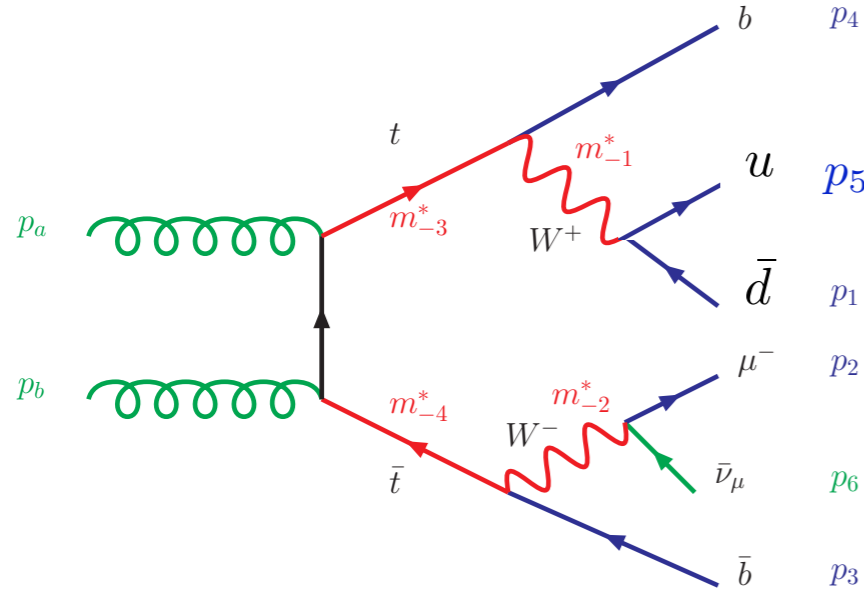
$$\mathcal{P}(\mathbf{p}^{vis} | \alpha) = \frac{1}{\sigma_\alpha} \int d\Phi dx_1 dx_2 |M_\alpha(\mathbf{p})|^2 W(\mathbf{p}, \mathbf{p}^{vis})$$

- Likelihood need to be correctly normalized

$$L(\alpha) = \prod_{i=1}^N \mathcal{P}(\mathbf{p}_i^{vis} | \alpha)$$

$$-\ln(L(\alpha)) = -\sum_{i=1}^N \ln(P_i^{MW}) + N * \ln(\sigma_\alpha)$$

- 
- 2009: MadGraph4 Implementation
 - 2011: Private Implementation in MadGraph5
 - Initial State Radiation Support
 - SubProcess grouping (speed)
 - NWA (speed)
 - 2013: MadWeight5
 - Improve cluster support (speed)
 - MC over jet/parton assignment (speed)
 - pre-training (speed)
 - better multi-channel (speed)
 - 2014: MadWeight5
 - Support for multi-transfert function estimated on the same phase-space point (speed)
 - Module of preselection of the jet/parton assignment (speed)



| | | | | | | | | | | |
|---|---|--------|-------|-------|------|------|-----|------|-----|-----|
| 0 | | | 12 | 3587 | | | | | | |
| 1 | 1 | 0.935 | 5.230 | 83.80 | 0.00 | 1.0 | 0.0 | 0.00 | 0.0 | 0.0 |
| 2 | 4 | -0.161 | 1.878 | 85.60 | 9.66 | 7.0 | 0.0 | 1.10 | 0.0 | 0.0 |
| 3 | 4 | -0.223 | 5.295 | 45.64 | 5.43 | 3.0 | 0.0 | 0.30 | 0.0 | 0.0 |
| 4 | 4 | 0.695 | 2.208 | 37.99 | 7.68 | 8.0 | 0.0 | 3.63 | 0.0 | 0.0 |
| 5 | 4 | 1.164 | 3.357 | 49.01 | 6.95 | 13.0 | 0.0 | 2.66 | 0.0 | 0.0 |
| 6 | 6 | 0.000 | 6.035 | 39.48 | 0.00 | 0.0 | 0.0 | 0.00 | 0.0 | 0.0 |

Which jet correspond to which parton?

MW4: one integral per permutation

MW5: Monte-Carlo over the permutation

| process | tf | permutation | Sum/MonteCarlo |
|------------------|-------|-------------|----------------|
| tt semi leptonic | delta | 24 | 7.5 |
| tt semi leptonic | gauss | 24 | 2 |
| tt di leptonic | gauss | 2 | 0.6 |
| w+ j j | delta | 2 | 1.5 |
| tth (semi lept) | gauss | 720 | 20 |

| process | perm | MW4 | MW5 |
|------------------|------|----------|-------|
| tt semi lept | 24 | 1h16 | 41s |
| tt fully lept | 2 | 46s | 10s |
| tth semi lept | 720 | > 2 days | 10min |
| tth semi lept | 48 | > 3h | 6min |
| tth fully lept | 24 | > 1h | 1min |
| h> w+ w- > 1lept | 2 | 59s | < 5s |
| h> w+ w- > 2lept | 1 | 8s | < 5s |
| z b b | 24 | 39m | 18s |
| zh | 24 | 43m | < 5s |

running on 1 core of a Intel core i7 2.3Ghz

Installation:

- Download the tar-ball from indico
 - everything is inside
 - you are ready to go.

Exercise 1:

- Extract the top-quark mass from a sample of $t\bar{t}$ events (fully leptonic decay) at parton level
- Use the likelihood method (do not forget to normalize it)

Exercise 2:

- Same but at detector level
- Show the sensitivity to the transfer function

Exercise 3:

- Add background

- Download the package from indico
- `$> tar -xzip Tutorial.tgz`
- You have
 - ➔ `Tutorial/samples/`: samples of events that we will use
 - ➔ `Tutorial/MG5_aMC`: code of MadGraph/MadWeight
 - ➔ `Tutorial/MG5_aMC/bin/mg5_amc` : executable
 - ➔ `Tutorial/mg5`: symbolic link to the previous executable

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creation of the dedicated code:

- run the executable
 - `$> cd Tutorial`
 - `$> ./mg5` (or `./MG5_aMC/bin/mg5_amc`) [open a prompt]
 - `mg5_amc> generate p p > t t~, t > b e+ ve, t~ > b~ mu- vm~`
 - `mg5_amc> output madweight MW_TT_FULL_LEPT`

result:

- Now you have a directory in your current directory:
`MW_TT_FULL_LEPT`

comments:

- All LO MG5 syntax are supported (but no NLO)
- For semi-leptonic:
 - `mg5_amc> generate p p > t t~, t > b e+ ve, t~ > b~ j j`
 - `mg5_amc> add process p p > t t~, t > b j j, t~ > b~ e- ve~`

computing the weight:

- `mg5_amc> launch MW_TT_FULL_LEPT`

OR

- `$> cd MW_TT_FULL_LEPT`
- `./bin/madweight.py`

output:

```
Do you want to edit a card (press enter to bypass editing)?
1 / param      : param_card.dat
2 / run        : run_card.dat
3 / madweight  : madweight_card.dat
4 / transfer   : transfer_card.dat
5 / lhco       : input.lhco
you can also
- enter the path to a valid card or banner.
- use the 'set' command to modify a parameter directly.
  The set option works only for param_card and run_card.
  Type 'help set' for more information on this command.
- call an external program (ASperGE/MadWidth/...).
  Type 'help' for the list of available command
- use the 'change_tf' command to set a transfer functions.
[_@, done, 1, param, 2, run, 3, madweight, 4, enter path, ... ][60s to answer]
```

comments:

- param: default model parameter
- run: accelerator/cut (if used)
- madweight: see later
- transfer: definition of TF
- lhco: input data for events

define transfer function:

- answer the question with
- `change_tf all_delta`

output:

- The same question is re-asked.

comments:

- This defines ONLY the functional form of the transfer function
- The numerical coefficient can be changed in `transfer_card.dat`
- There are no numerical coefficients for this transfer function
 - neutrinos are not constrained
 - all other particles are assumed to be perfectly measured
- You can define your own functional form if needed (see backup)

define the input file / number of events /...

- answer the question with
 - madweight

output

- open the madweight_card within an editor (typically vi)

action

- change
 - the number of event to 25 (nb_exp_events)
 - the input path to `././samples/parton.lhco.gz` (inputfile)
 - define a scan over the top mass from 163 to 183 (by step of 5GeV)
 - name of the run to "parton" (name)

comment

- Note that this card de-activate the cut of the run_card. (use_cut)
- you can change each of those parameters with "set name VALUE"

other interesting parameters

- `nb_event_by_node`: for running on cluster. (option for cluster define in `MG5_aMC/input/mg5_configuration.txt`)
- `use_cut`: defines if the cut of the `run_card` are applied or not. (in principle no cut should be applied)
- `permutation`: using only one (or all) jet-parton assignment
- `montecarlo`: using MC over the jet-parton assignment
- more info on the card: see the online wiki.

launch the run

- answer the question with
- done (or just press enter)

What's going on

- The parametrization of the phase-space is chosen
- The actual computation takes places ($N_{\text{event}} \cdot N_{\text{th}}$)

output

- in Events/parton/
 - un-normalized_likelihood.out: $-\ln(L)$ without the cross-section
 - weights.out: column format with the weights
 - output.xml: structured format contains details of the computation
 - The amount of information depends of the “log_level” parameter (can contain weight for each permutation/full log/...)

Compute the normalization factor

- we will use mg5_amc for that
- `$> ./mg5` (or `./MG5_aMC/bin/mg5_amc`) [open a prompt]
- `mg5_amc> generate p p > t t~, t > b e+ ve, t~ > b~ mu- vm~`
- `mg5_amc> output MG_TT_FULL_LEPT`
- `mg5_amc>launch`

output:

```
The following switches determine which programs are run:
1 Run the pythia shower/hadronization:                pythia=OFF
2 Run PGS as detector simulator:                       pgs=OFF
3 Run Delphes as detector simulator:                   delphes=OFF
4 Decay particles with the MadSpin module:             madspin=OFF
5 Add weights to the events based on changing model parameters: reweight=OFF
Either type the switch number (1 to 5) to change its default setting,
or set any switch explicitly (e.g. type 'madspin=ON' at the prompt)
Type '0', 'auto', 'done' or just press enter when you are done.
[0, 1, 2, 3, 4, 5, auto, done, pythia=ON, ... ][60s to answer]
```

output:

- we do not need any of this (shower/...) so just press enter

output:

```
Do you want to edit a card (press enter to bypass editing)?
1 / param      : param_card.dat
2 / run        : run_card.dat
9 / plot       : plot_card.dat
you can also
- enter the path to a valid card or banner.
- use the 'set' command to modify a parameter directly.
  The set option works only for param_card and run_card.
  Type 'help set' for more information on this command.
- call an external program (ASperGE/MadWidth/...).
  Type 'help' for the list of available command
[_0, done, 1, param, 2, run, 9, plot, enter path][60s to answer]
```

action:

- edit the run_card.dat to remove ALL cut. (type run to edit the card)
- then type “set mt scan:range(163,185,5)” [this defines a scan on the top mass]
- then press enter

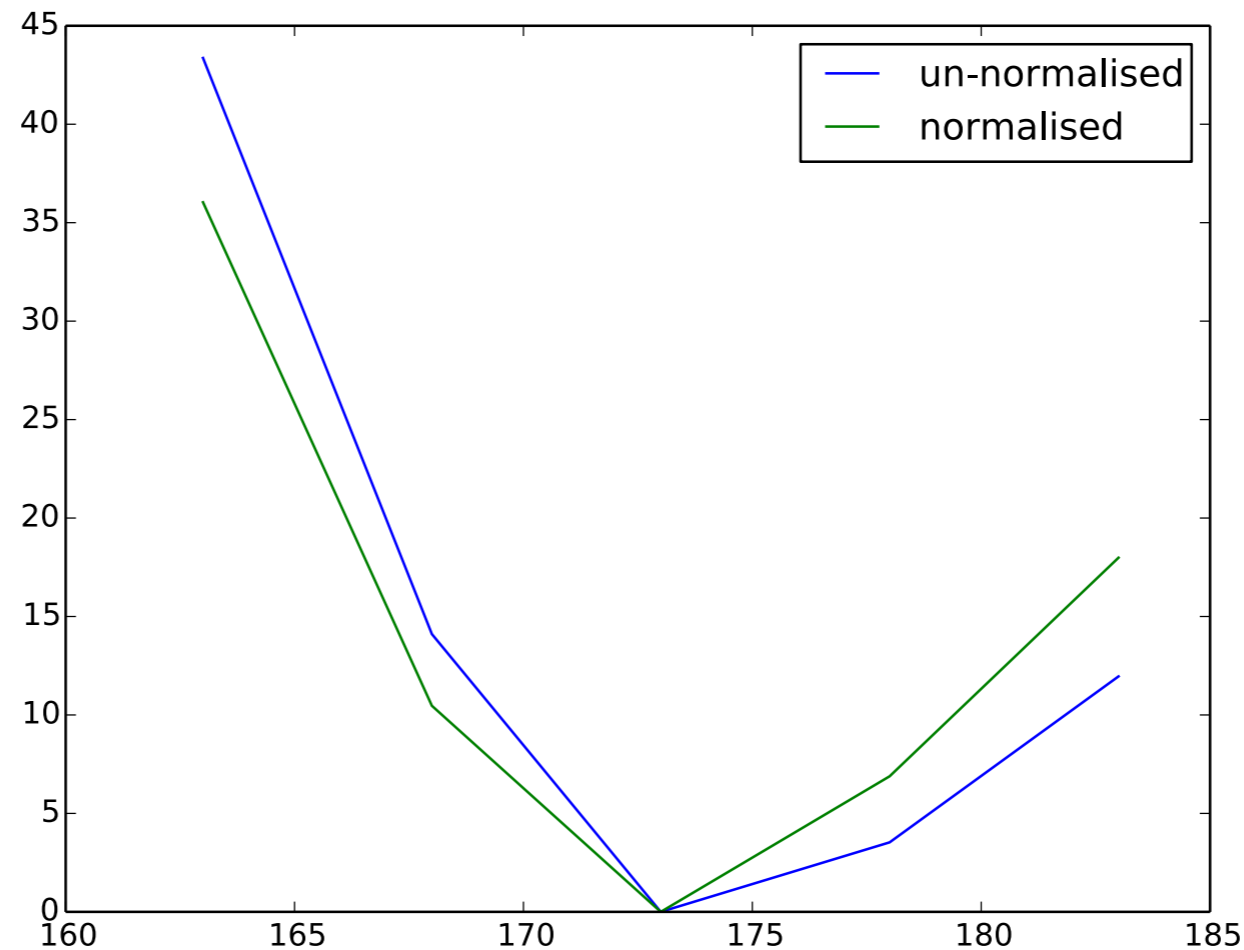
output:

- MG_TT_FULL_LEPT/Events/scan_run_[01-05].txt
 - contains the cross-section for each top mass value

make the plot

- use your favorite program for that:

Expected output



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Exercise 2:

- Same but at detector level
- Show the sensitivity to the transfer function

Exercise 3:

- Add background

creation of the dedicated code:

- `$> cd MW_TT_FULL_LEPT`

comments:

- Since the process is the same (but not the transfer function) we can reuse the same directory/code

configure:

- run `./bin/madweight.py`
- at the question type “change_tf dbl_gauss_pt_jet”

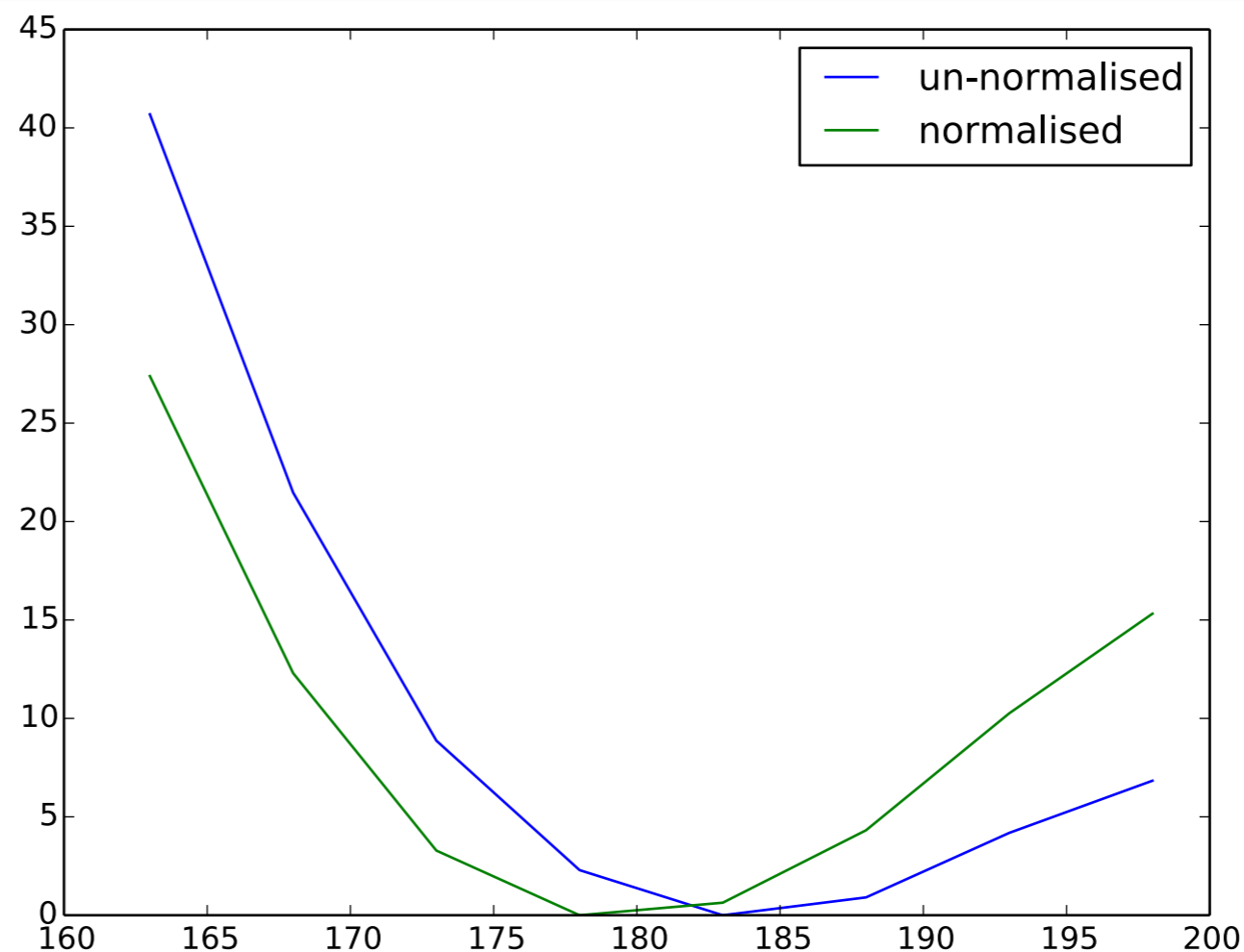
comment:

- the same as all_delta but for jet where the pt is related to a resolution function (double gaussian where each parameter is a function of the pt)
- Now you can edit the value of various parameter in the transfer_card

change the path to the event file:

- pass to ../samples/detector.lhco.gz
- change the name to “detector”

expected output:



question:

- explain the bias (the correct value is 173)

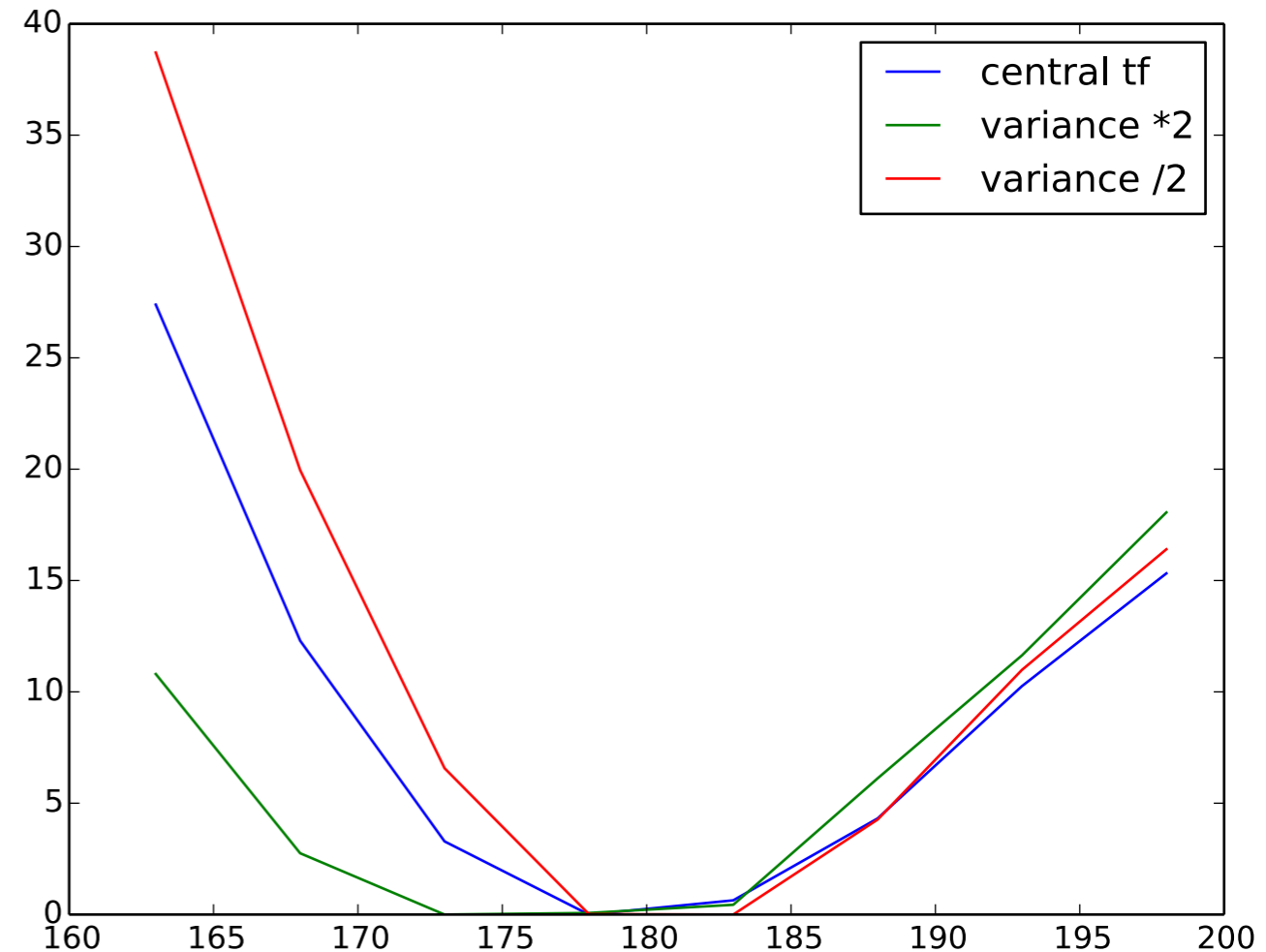
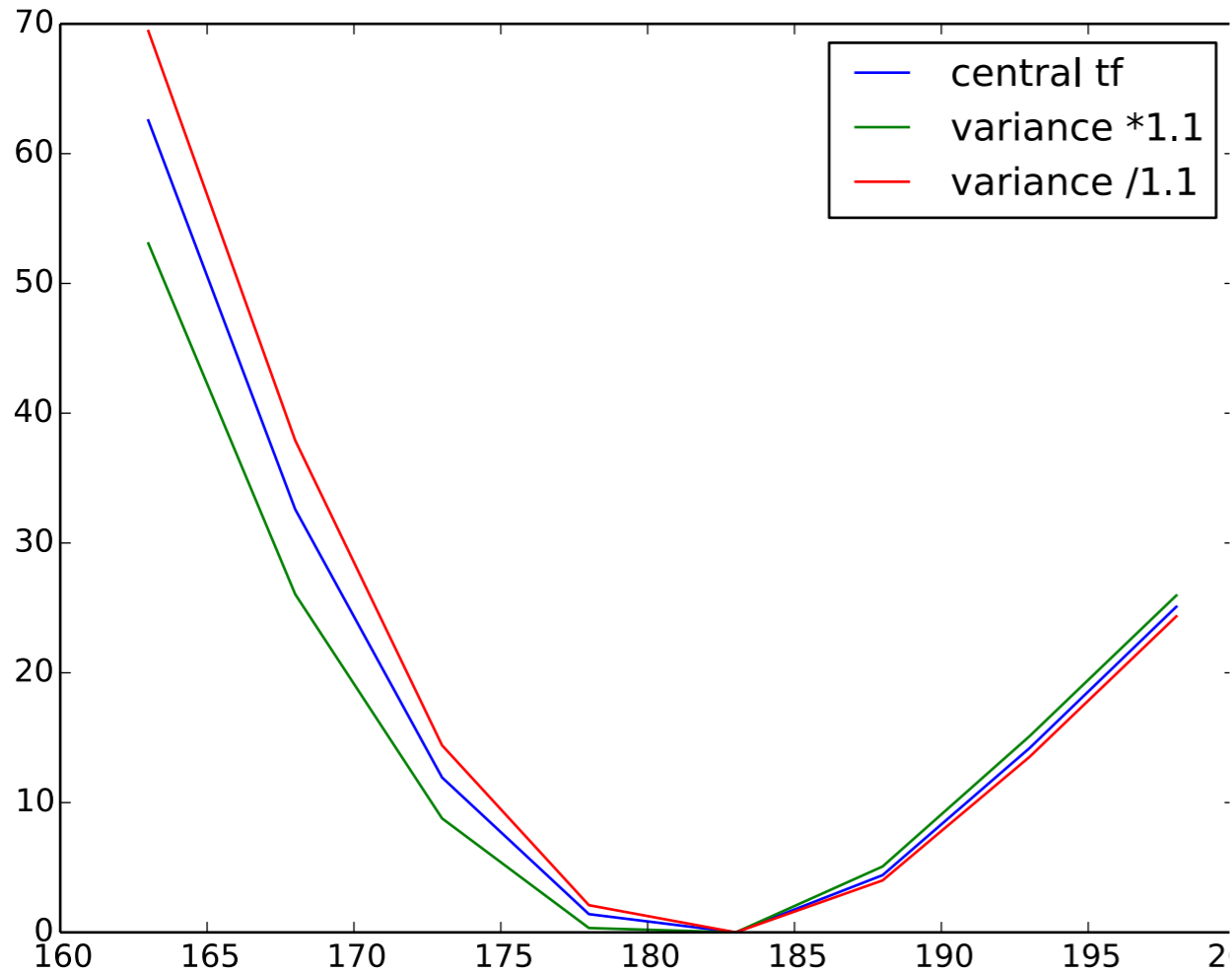
study the impact of transfer function

- rerun the same computation but with the variance of the transfer function divide and multiply by two.
- do it in a single run.

edit transfer_card.dat

- all entry should have three entries:
 - 4 0.9518d0 # sigma first gaussian
 - 4 1.8d0 # for the second tf
 - 4 0.47 # for the third tf
- unmodified parameter should also have three entries.

run and compare the running time



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run and compare the running time

- background is here $w_+ w_- + 2j$
- in samples

- Thanks for your participation, questions and remarks.

```

#####
##                                TF JET                                ##
#####
<block name='jet'>  #name can be anything
<info> doubl gaussian with parameter depending of the energy </info>
<particles> u,d,s,c,b,g </particles>
# this defined when this tf will be used.the letter correspond to the label in
#      particles.dat
<width_type> large </width_type>
# width_type should be thin or large (thin is for energy accurate up to 5-10%)
<variable name='E'>
<tf>
  prov1=(#1+#2*dsqrt(pexp(0))+#3*pexp(0))
  prov2=(#4+#5*dsqrt(pexp(0))+#6*pexp(0))
  prov3=(#7+#8*dsqrt(pexp(0))+#9*pexp(0))
  prov4=(#10+#11*dsqrt(pexp(0))+#12*pexp(0))
  prov5=(#13+#14*dsqrt(pexp(0))+#15*pexp(0))

  tf=(exp(-(p(0)-pexp(0)-prov1)**2/2d0/prov2**2))           !first gaussian
  tf=tf+prov3*exp(-(p(0)-pexp(0)-prov4)**2/2d0/prov5**2)   !second gaussian
  tf=tf*((1d0/dsqrt(2d0*pi))/(prov2+prov3*prov5))         !normalisation
</tf>
<width>
  prov2=(#4+#5*dsqrt(pexp(0))+#6*pexp(0))
  prov5=(#13+#14*dsqrt(pexp(0))+#15*pexp(0))

  width=max(prov2,prov5)
</width>
</variable>

# in this case THETA/PHI are not defined because they are considered
# in delta (=default)
# The same syntax apply
</block>

```