

MG5aMC tutorial

NLO / Loop induced / MLM

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MadGraph5

- Cross-section @LO
- event generation @LO
 - including loop-induced
- Interface to
 - Pythia6 (8 in progress)
 - Delphes
 - MadSpin
- Any BSM model supported

—

aMC@NLO

- Cross-section @NLO
- event generation @NLO+PS
- Interface to
 - PY6/PY8/HW6/HW++ (HW7 in progress)
 - MadSpin
- BSM in quick expansion
 - Framework ready

- <https://launchpad.net/madgraph5>
 - untar it (`tar -xzipvf MG5_XXX.tgz`)
 - launch it (`$./bin/mg5_amc`)
- **learn** it!
 - ➔ Type **tutorial** and follow instructions
- install external package
 - ➔ install pythia-pgs
 - ➔ install MadAnalysis
- put MC4BSM UFO model from indico in the models directory

[Home page](#) [Wiki](#)

The MadGraph Matrix Element Generator version 5

[Overview](#) [Code](#) [Bugs](#) [Blueprints](#) [Translations](#) [Answers](#)

Registered 2009-09-15 by [Michel Herquet](#)

The version 5 of the MadGraph Matrix Element Generator for the simulation of parton-level events for decay and collision processes at high energy colliders. Allows matrix element generation and event generation for any model that can be written as a Lagrangian, using the output of the FeynRules Feynman rule calculator. Provides output in multiple formats and languages, including Fortran MadEvent, Fortran Standalone matrix elements, C++ matrix elements, and Pythia 8 process libraries.

Note that process generation can also be done directly online at <http://madgraph.phys.ucl.ac.be> or <http://madgraph.hep.uiuc.edu>.
If you use MadGraph 5, please cite JHEP 1106(2011)128, arXiv:1106.0522 [hep-ph].

Installation:
MadGraph 5 needs Python version 2.6 or 2.7. The latest stable release is in the trunk, which can be branched using the Bazaar versioning system:
`bzr branch lp:madgraph5`
or be downloaded as a tar.gz package to the right. This release contains everything needed for process generation in multiple models, as well as event generation through MadEvent, and standalone matrix element evaluation for Fortran or C++ output.
In order to use the process library output for Pythia 8, you need Pythia 8.150 or later installed.

Getting started:
Run `bin/mg5` and type "help" to learn how to run MadGraph 5 using the command interface, or run the interactive quick-start tutorial by typing "tutorial".
Or copy the Template, edit the `Cards/proc_card_mg5.dat` and run `bin/newprocess_mg5`.

Examples of process generation syntax:
`pp > w+ jj`
`pp > t t-, t > b jj, t- > b- l- vl-`
`e+ e- > z > n2 n2, (n2 > x1+ w-, x1+ > l+ vl n1, w- > l- vl-), n2 > jj n1`

To output model files for MadGraph 5 with FeynRules, use version 1.6 or later, and use the WriteUFO command.

Get Involved

[Report a bug](#) →

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[Help translate](#) →

Downloads

Latest version

MadGraph5_v1.5.10.tar.gz ↓

MadGraph5_v...eta3.tar.gz ↓

released on 2012-09-29

[All downloads](#)

Announcements

aMC@NLO in MadGraph5 on 2012-11-08

On Nov 8th 2012, version 2.0 beta of MadGraph5 has been released. This is a m...

[Read all announcements](#)

Project information

Maintainer: Find:

Driver: Next Previous Highlight all Match case

Series and milestones

[View full history](#)

trunk

- Ask me
- Use the command “help” / “help XXX”
 - ➔ “help” tell you the next command that you need to do.
- Launchpad:
 - ➔ <https://answers.launchpad.net/madgraph5>
 - ➔ FAQ: <https://answers.launchpad.net/madgraph5/+faq>

- How do you change
 - ➔ top mass
 - ➔ top width
 - ➔ W mass
 - ➔ beam energy
 - ➔ pt cut on the lepton

Goal ● What's the default choice for QED/QCD order

Learn ● What's the difference between

→ $pp > tt$

→ $pp > tt$ QED ≤ 2

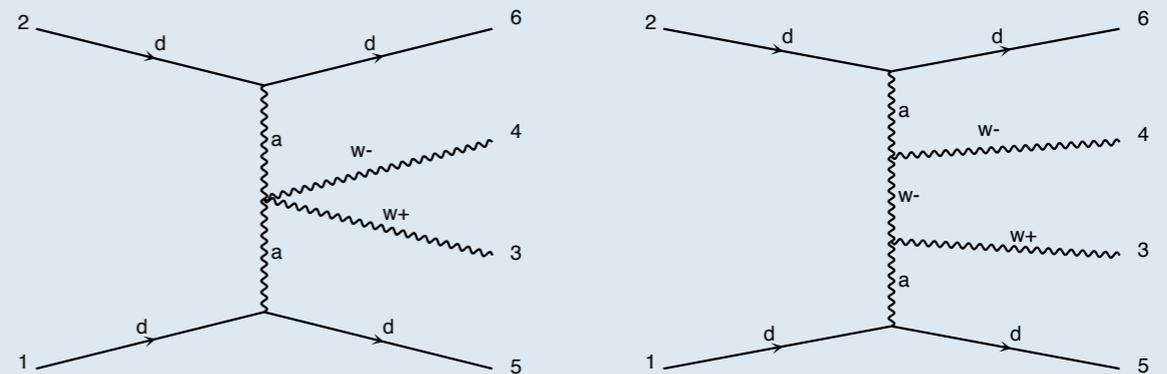
→ $pp > tt$ QED=0

- Compute the cross-section for each of those and check the diagram

Check

- Generate VBF process (two jet + two W in final state) **only the diagram!**

- check that you have the QED diagram that you want:



- What's the meaning of the order QED/QCD
 - ➔ By default MG5 takes the lowest order in QED!

INFO: Trying coupling order WEIGHTED<=2: WEIGHTED IS 2*QED+QCD

- ➔ $p p > t t^{\sim}$ IS the same as $p p > t t^{\sim}$ QED=0
- ➔ $p p > t t^{\sim}$ QED=2 has additional diagrams (photon/z exchange)

$p p > t t^{\sim}$

Cross section (pb)

555 ± 0.84

$p p > t t^{\sim}$ QED=2

Cross section (pb)

555.8 ± 0.91

No significant QED contribution

- generate $pp \rightarrow w^+ w^- jj$
 - 76 processes
 - 1432 diagrams
 - None of them are VBF

- generate $pp \rightarrow w^+ w^- jj$ QED = 2
 - 76 processes
 - 1432 diagrams
 - None of them are VBF

- generate $pp \rightarrow w^+ w^- jj$ QED = 4
 - 76 processes
 - 5332 diagrams
 - VBF present! + those not VBF

- generate $pp \rightarrow w^+ w^- jj$ QCD = 0
 - 60 processes
 - 3900 diagrams
 - VBF present!

- generate $pp \rightarrow w^+ w^- jj$ QCD = 2
 - 76 processes
 - 5332 diagrams

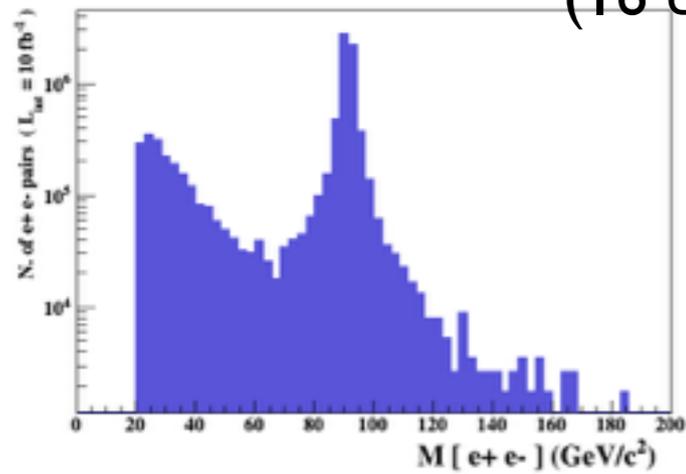
- generate $pp \rightarrow w^+ w^- jj$ QCD = 4
 - 76 processes
 - 5332 diagrams

- Generate the cross-section and the distribution (invariant mass) for
 - $pp > e^+ e^-$
 - $pp > z, z > e^+ e^-$
 - $pp > z > e^+ e^-$
 - $pp > e^+ e^- \$ z$
 - $pp > e^+ e^- / z$

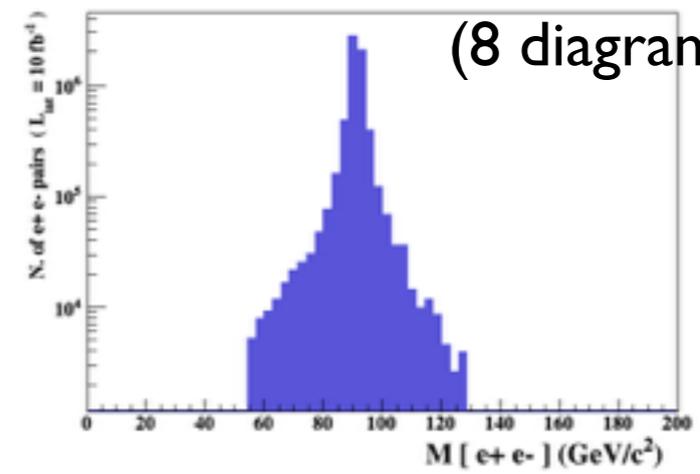
Hint :To plot automatically distributions:
`mg5> install MadAnalysis`

- Use the invariant mass distribution to determine the meaning of each syntax.

$pp \rightarrow e^+ e^-$
(16 diagrams)

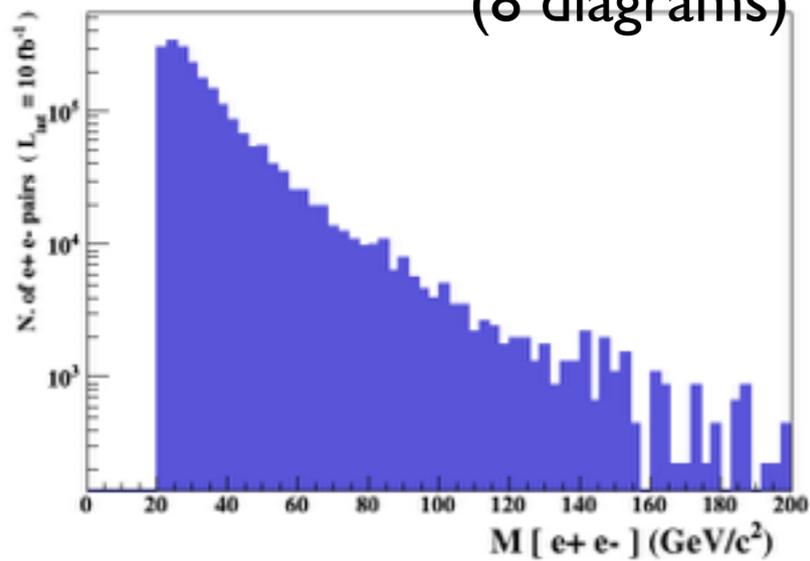


$pp \rightarrow z, z \rightarrow e^+ e^-$
(8 diagrams)



$pp \rightarrow e^+ e^- / z$

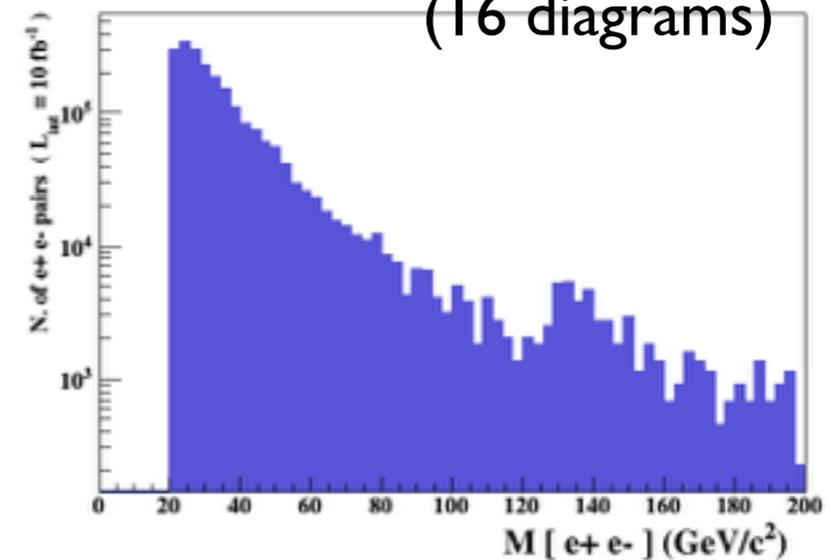
(8 diagrams)



No Z

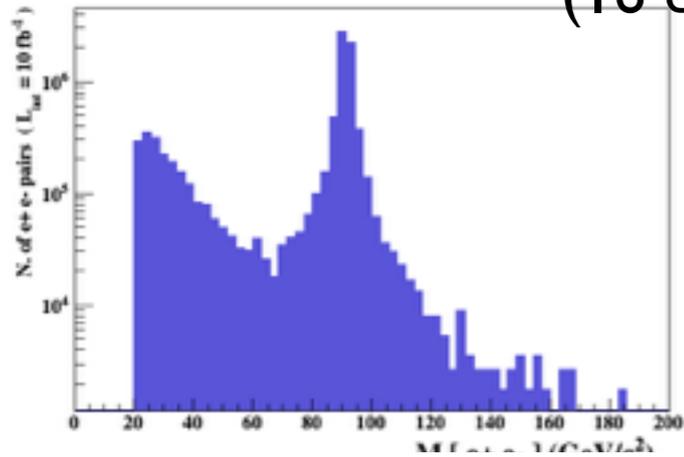
$pp \rightarrow e^+ e^- \cancel{z}$

(16 diagrams)



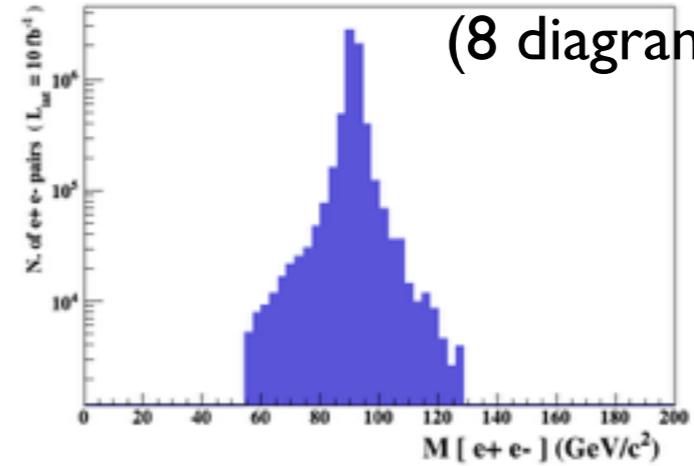
Z- onshell veto

$pp \rightarrow e^+ e^-$
(16 diagrams)



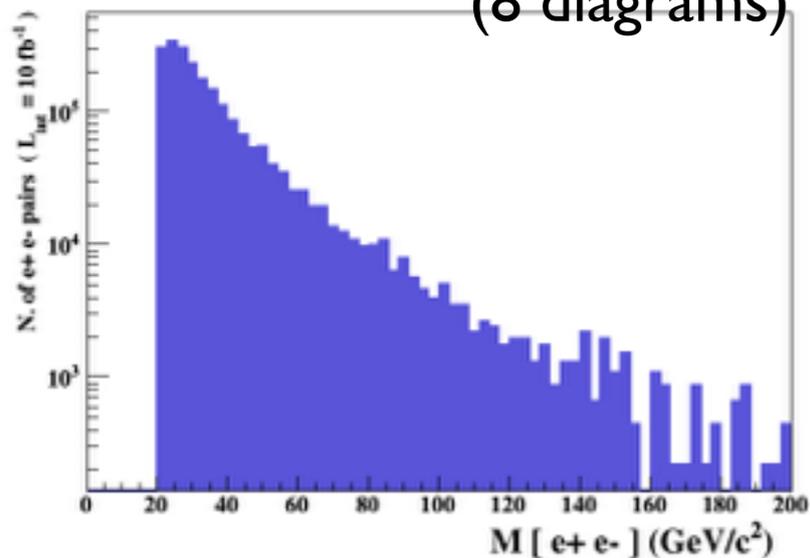
Correct Distribution

$pp \rightarrow z, z \rightarrow e^+ e^-$
(8 diagrams)



$pp \rightarrow e^+ e^- / z$

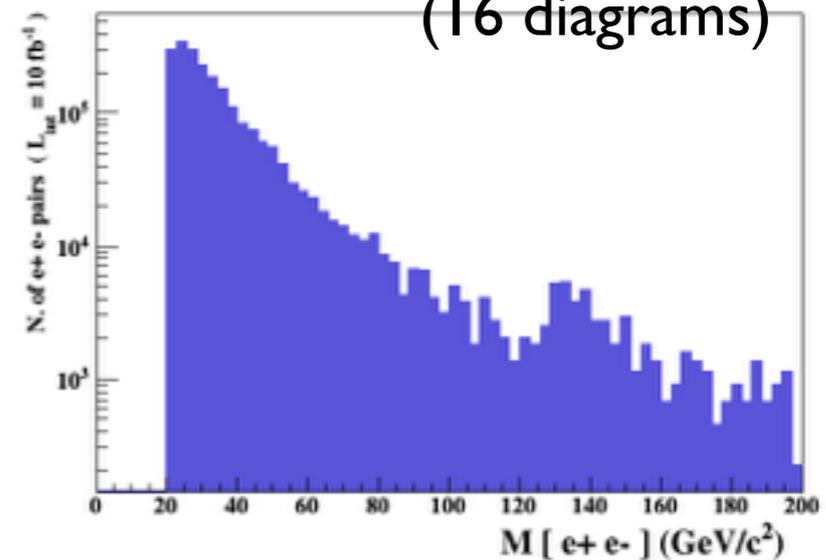
(8 diagrams)



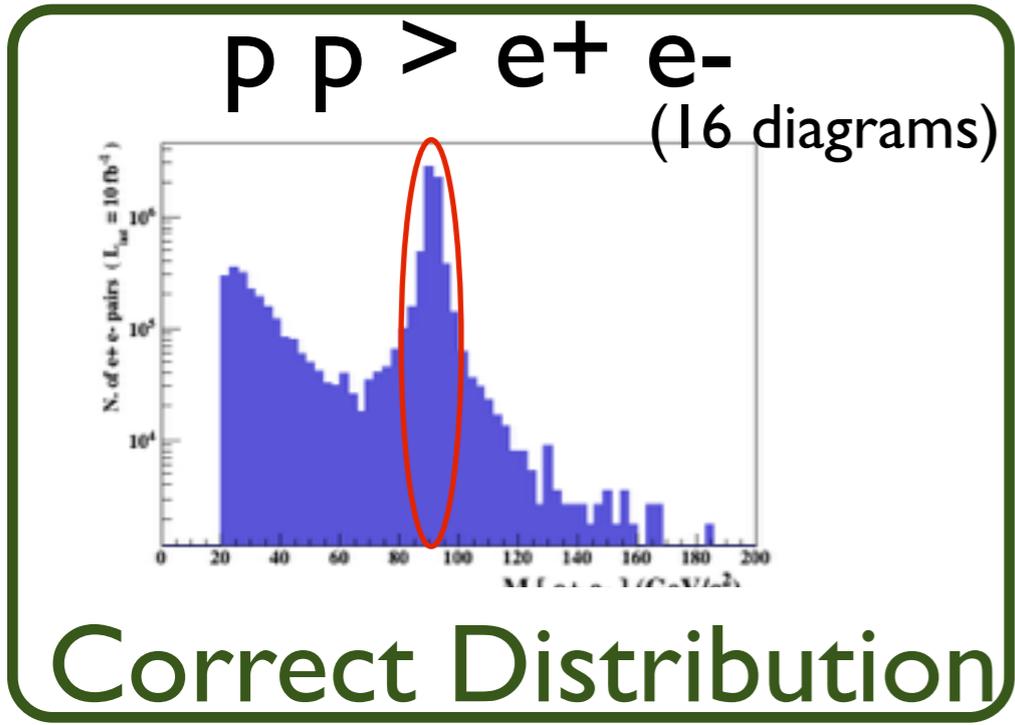
No Z

$pp \rightarrow e^+ e^- \cancel{z}$

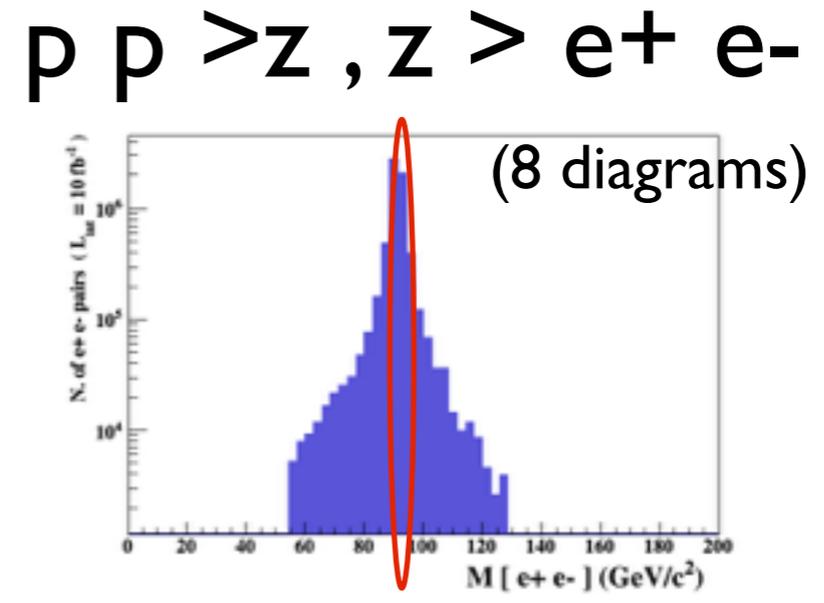
(16 diagrams)



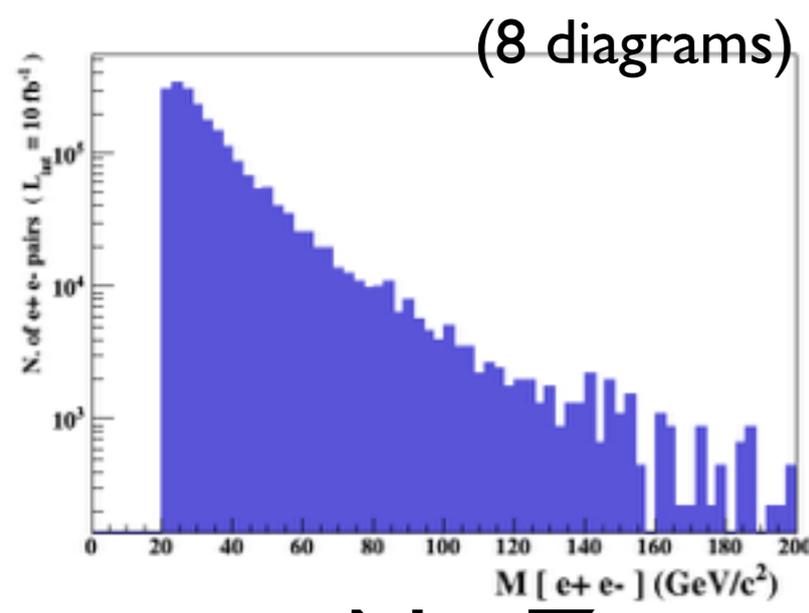
Z- onshell veto



Z Peak



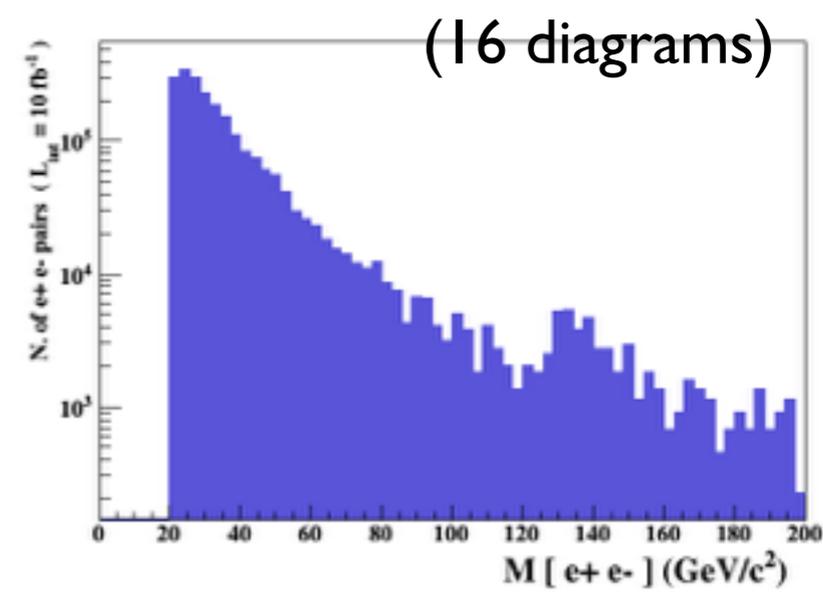
$p p \rightarrow e^+ e^- / z$



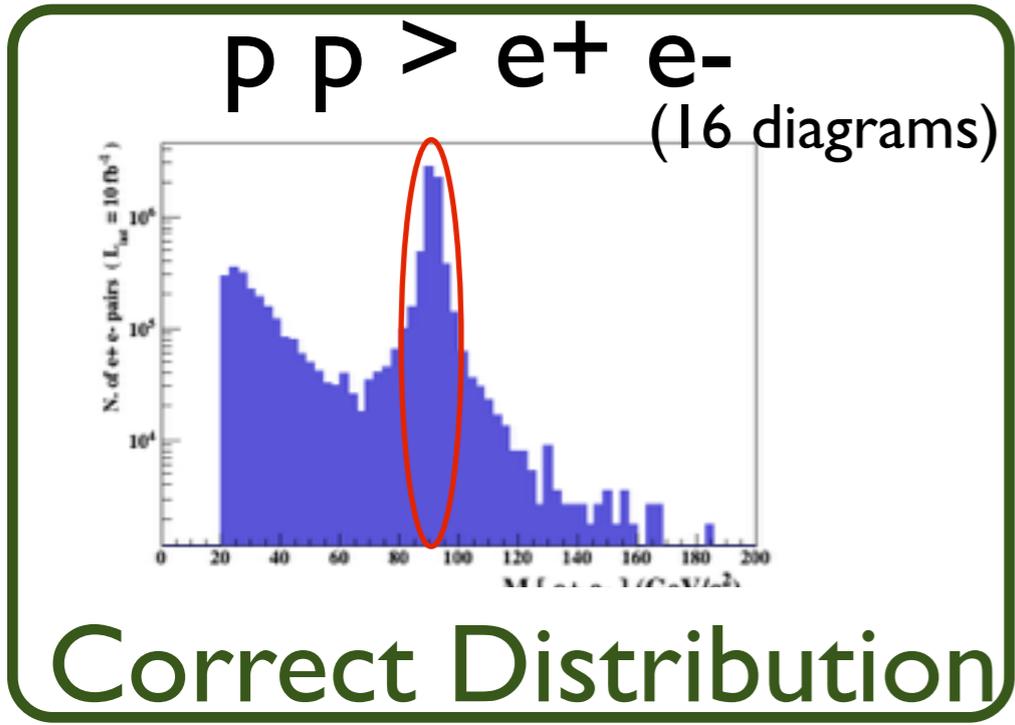
NO Z Peak

No Z

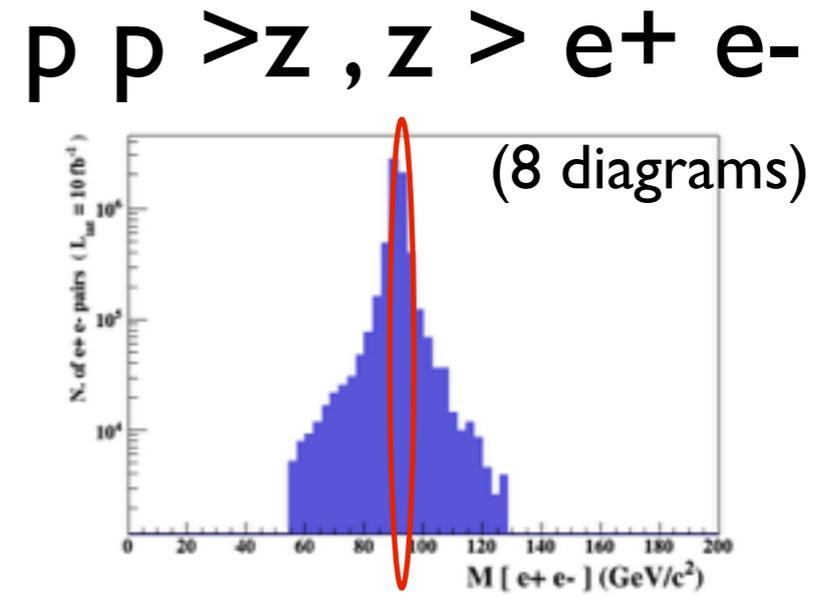
$p p \rightarrow e^+ e^- \cancel{z}$



Z- onshell veto

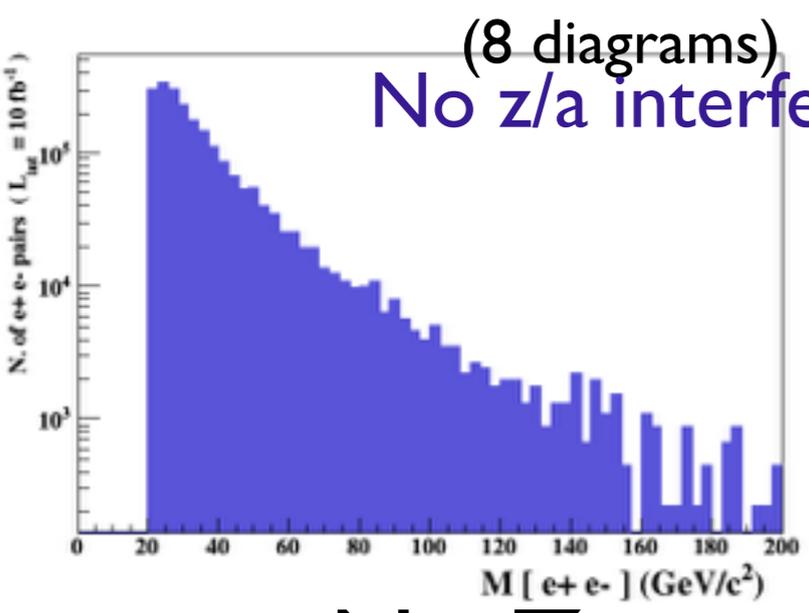


Z Peak



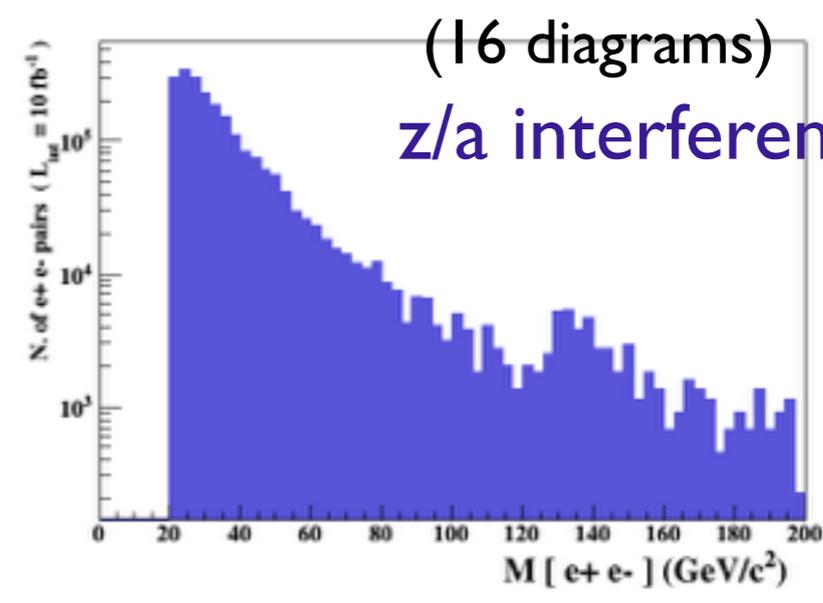
$p p \rightarrow e^+ e^- / z$

$p p \rightarrow e^+ e^- \cancel{z}$

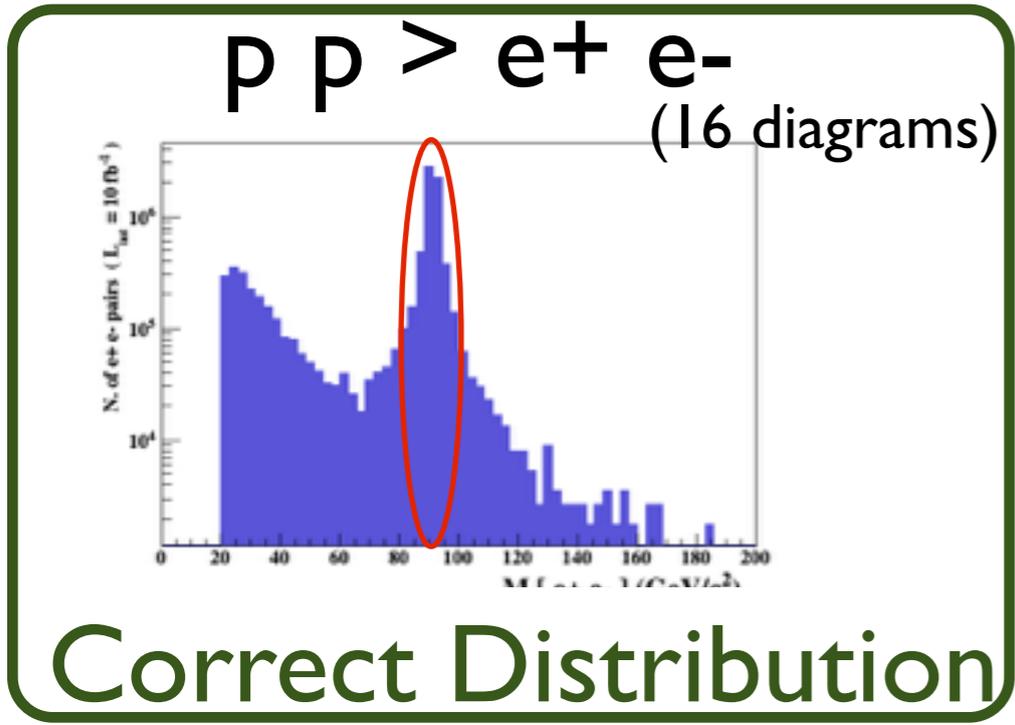


NO Z Peak

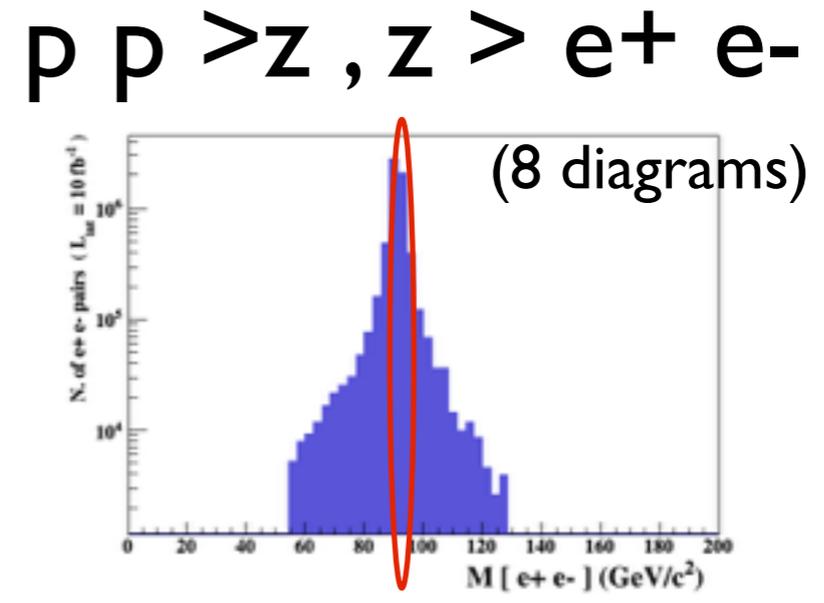
No Z



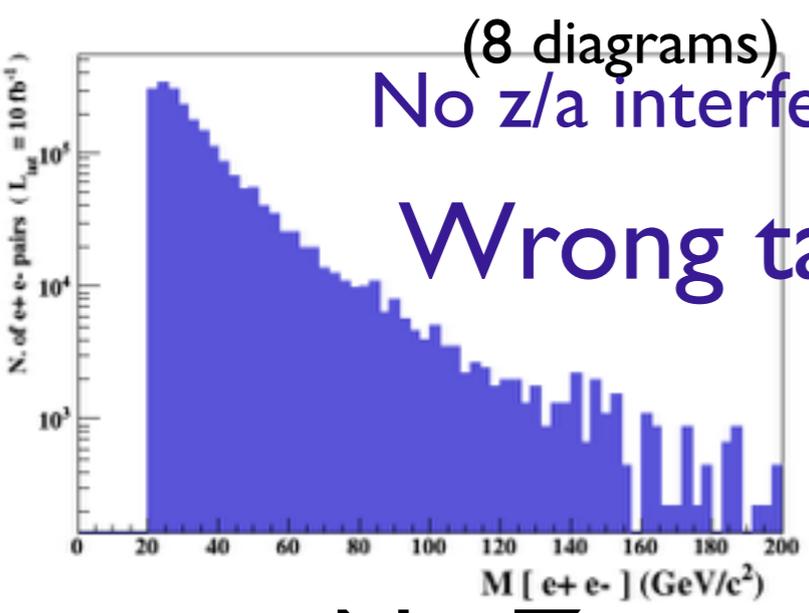
Z- onshell veto



Z Peak



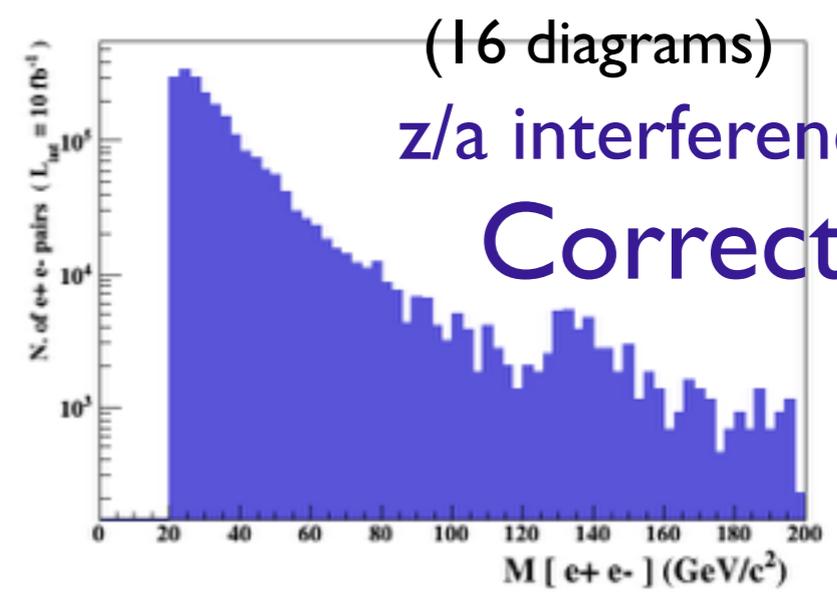
$p p \rightarrow e^+ e^- / z$



NO Z Peak

No Z

$p p \rightarrow e^+ e^- \cancel{z}$

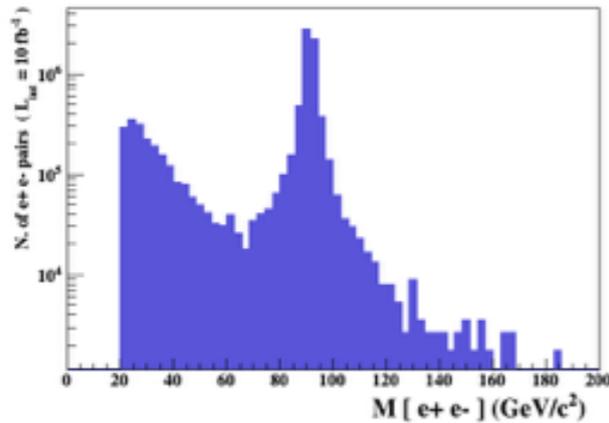


Z- onshell veto

$p p \rightarrow e^+ e^-$

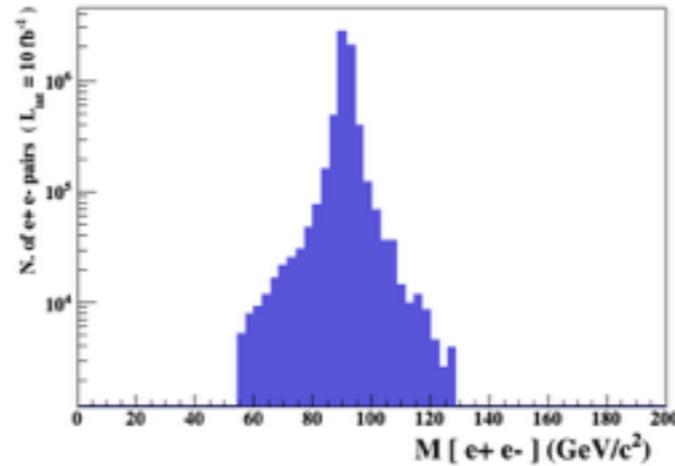
$p p \rightarrow Z, Z \rightarrow e^+ e^-$

$p p \rightarrow e^+ e^- \text{ \$ } Z$



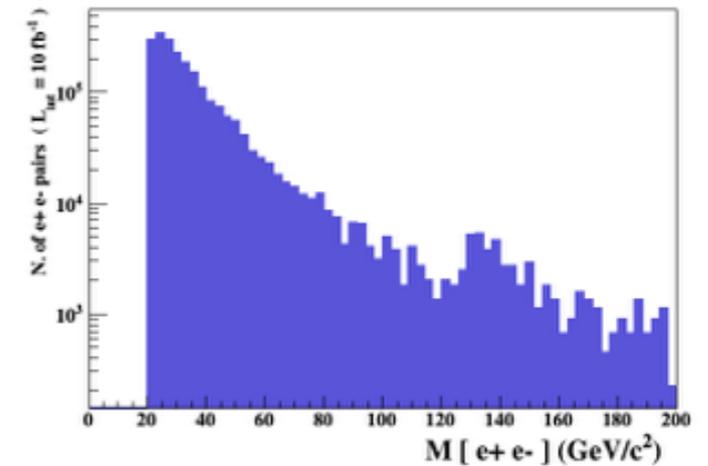
(16 diagrams)

=



(8 diagrams)

+



(16 diagrams)

Onshell cut: BW_cut

$$|M^* - M| < BW_{cut} * \Gamma$$

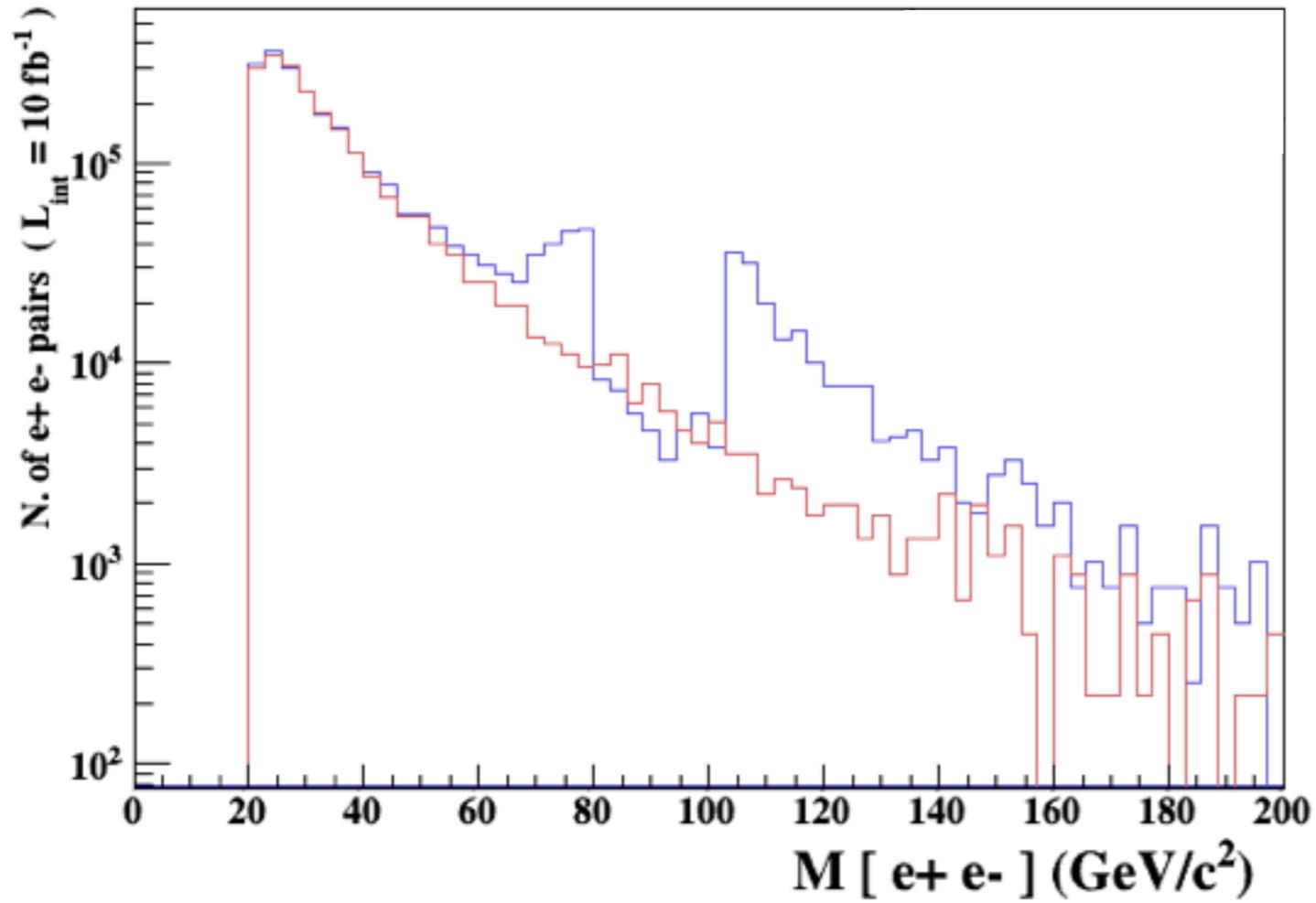
- The Physical distribution is (very close to) exact sum of the two other one.
- The “\$” forbids the Z to be onshell but the photon invariant mass can be at MZ (i.e. on shell subtraction).
- The “/” is to be avoid if possible since this leads to violation of gauge invariance.

- NEXT SLIDE is generated with `bw_cut = 5`
- This is **TOO SMALL** to have a physical meaning (15 the default value used in previous plot is better)
- This was done to **illustrate** more in detail how the “\$” syntax works.

$$p p > e^+ e^- / Z$$

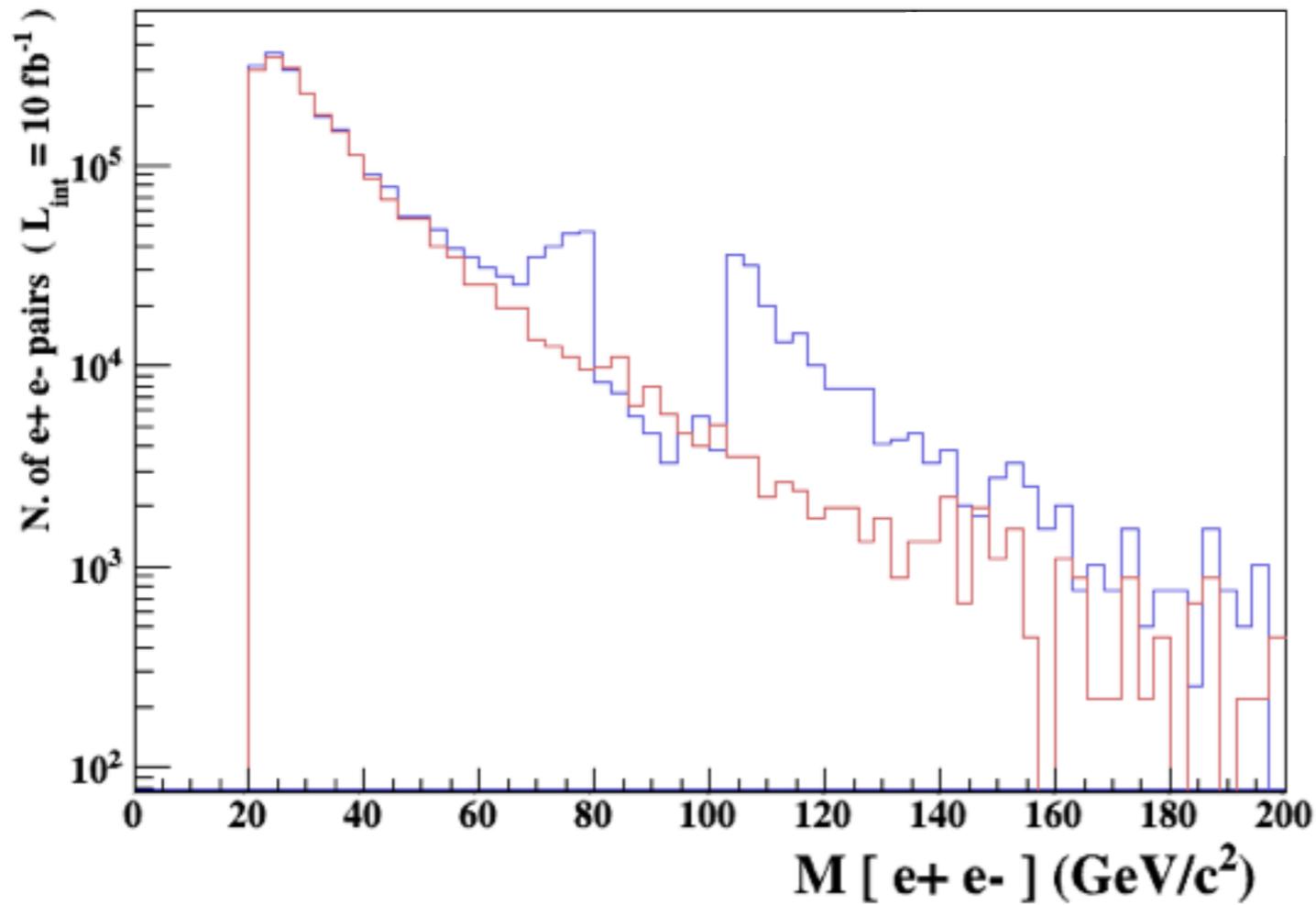
(red curve)

(blue curve)



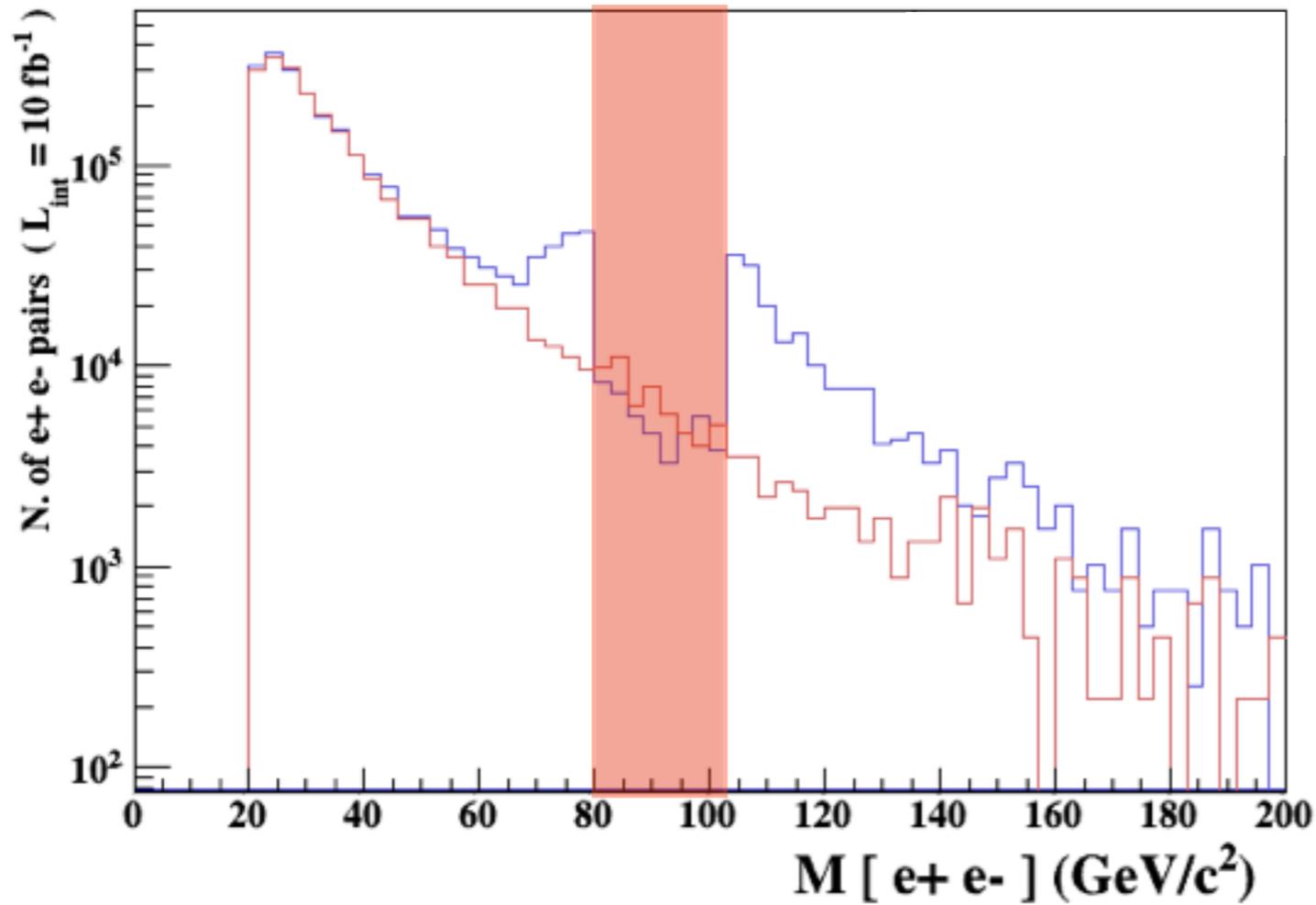
$p p \rightarrow e^+ e^- / Z$
(red curve)

adding $p p \rightarrow e^+ e^- \text{ } \$ Z$
(blue curve)



$p p > e^+ e^- / Z$
(red curve)

adding $p p > e^+ e^- \text{ } \$ Z$
(blue curve)

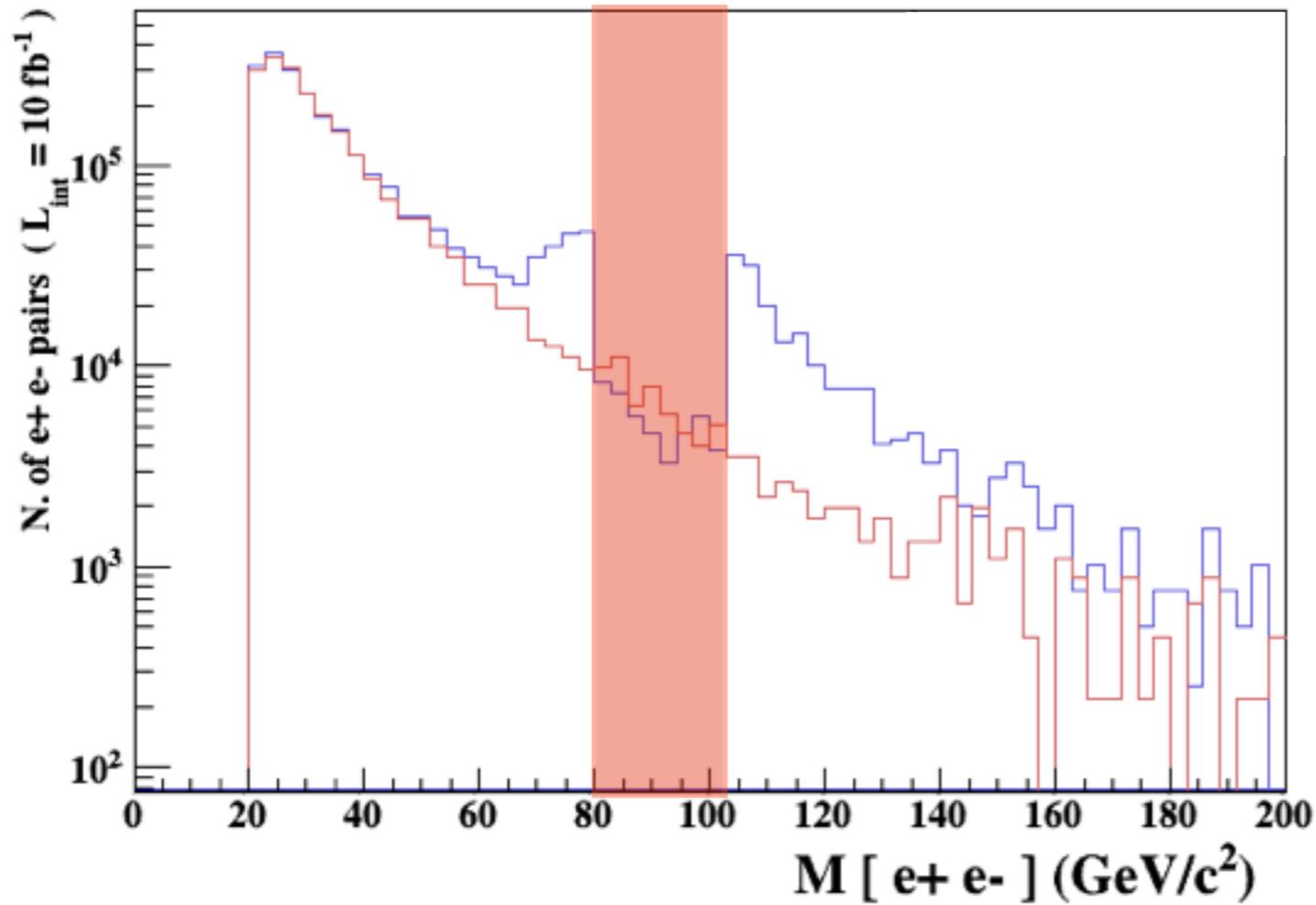


- Z onshell veto

5 times width area

$p p \rightarrow e^+ e^- / Z$
(red curve)

adding $p p \rightarrow e^+ e^- \gamma Z$
(blue curve)

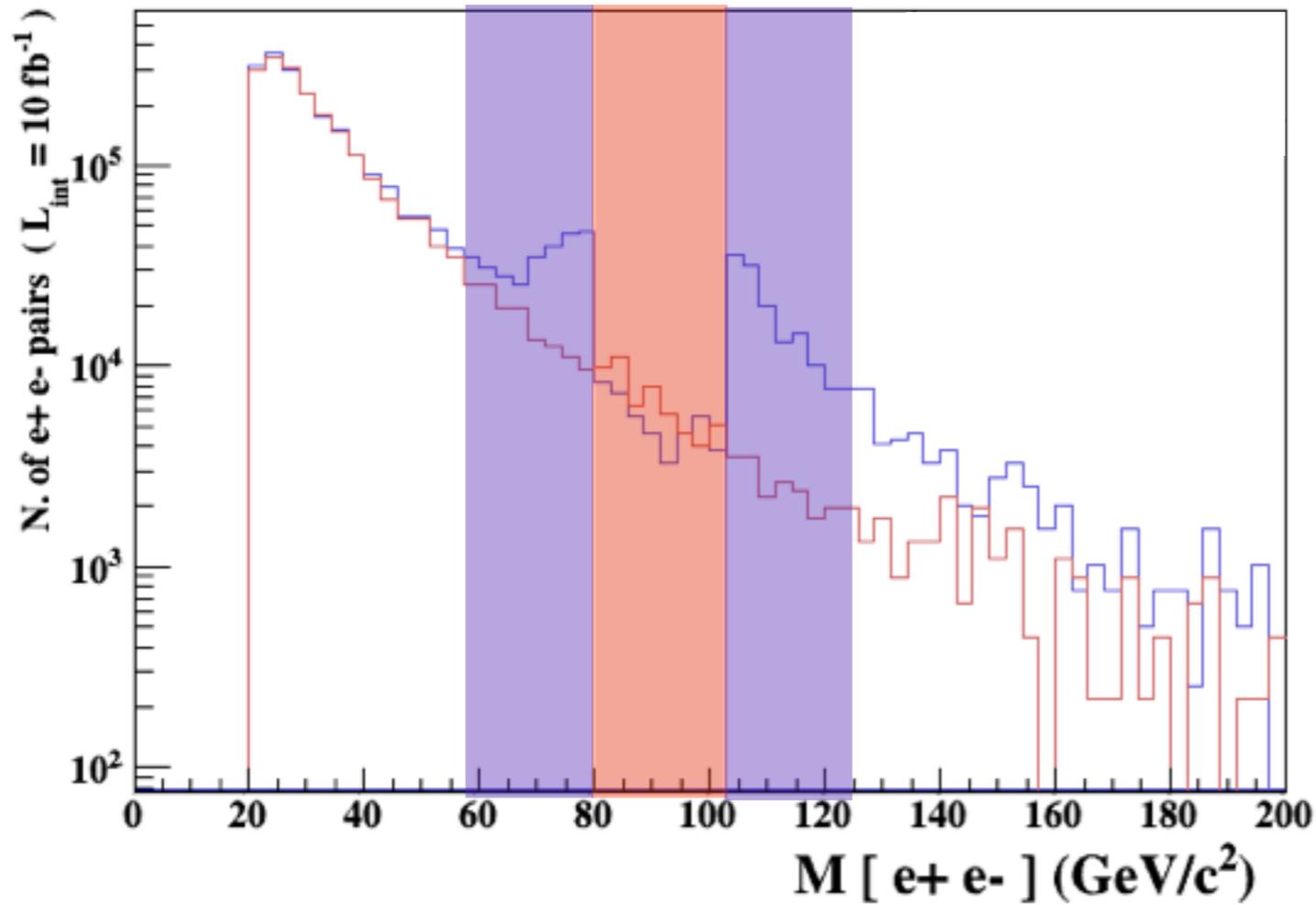


- Z onshell veto
- In veto area only photon contribution

5 times width area

$p p > e^+ e^- / Z$
(red curve)

adding $p p > e^+ e^- \text{ } \$ Z$
(blue curve)



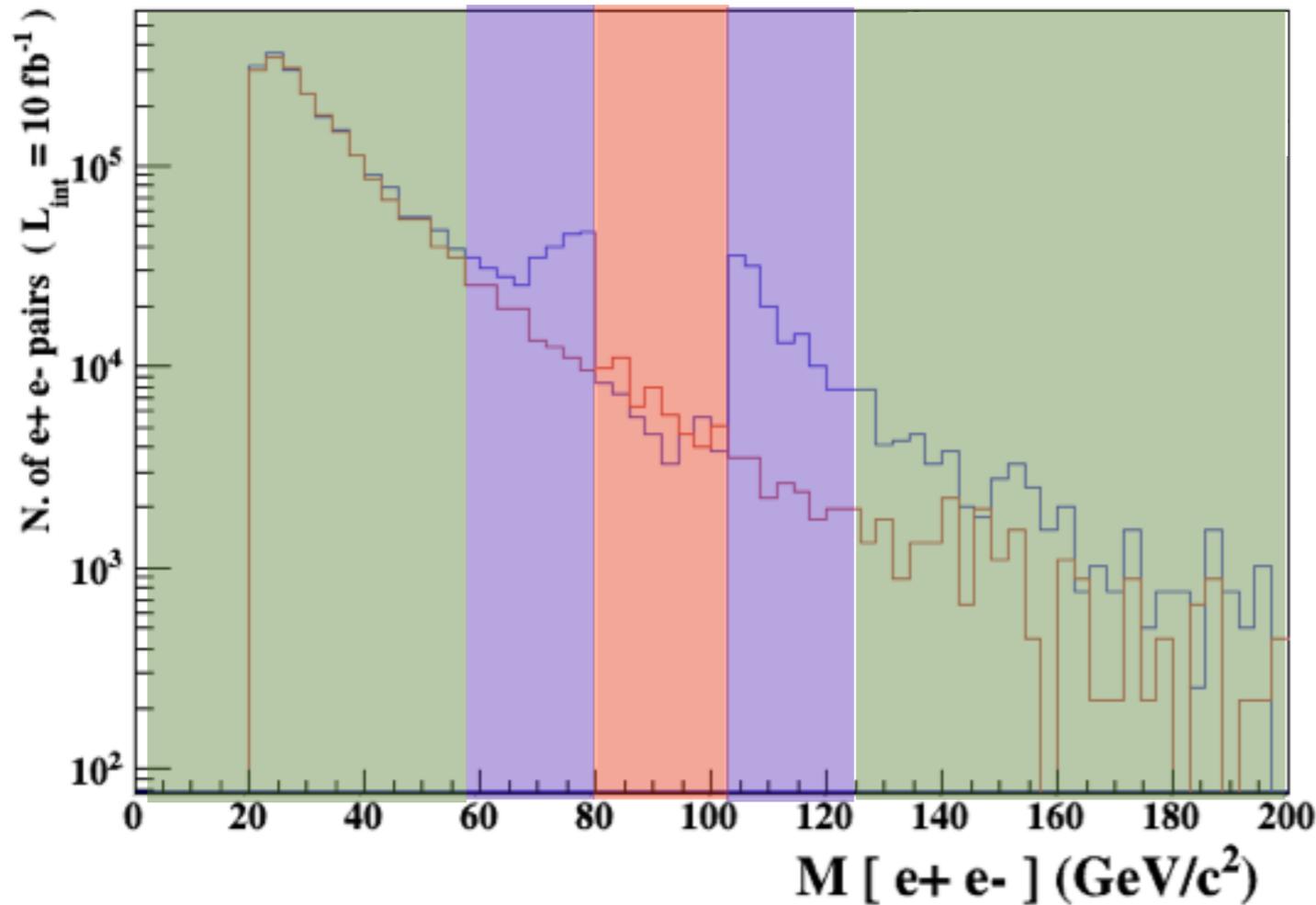
- Z onshell veto
- In veto area only photon contribution
- area sensitive to z-peak

5 times width area

15 times width area

$p p > e^+ e^- / Z$
(red curve)

adding $p p > e^+ e^- \text{ } \$ Z$
(blue curve)



- Z onshell veto
- In veto area only photon contribution
- area sensitive to z-peak
- very off-shell Z, the difference between the curve is due to interference which are need to be **KEPT** in simulation.

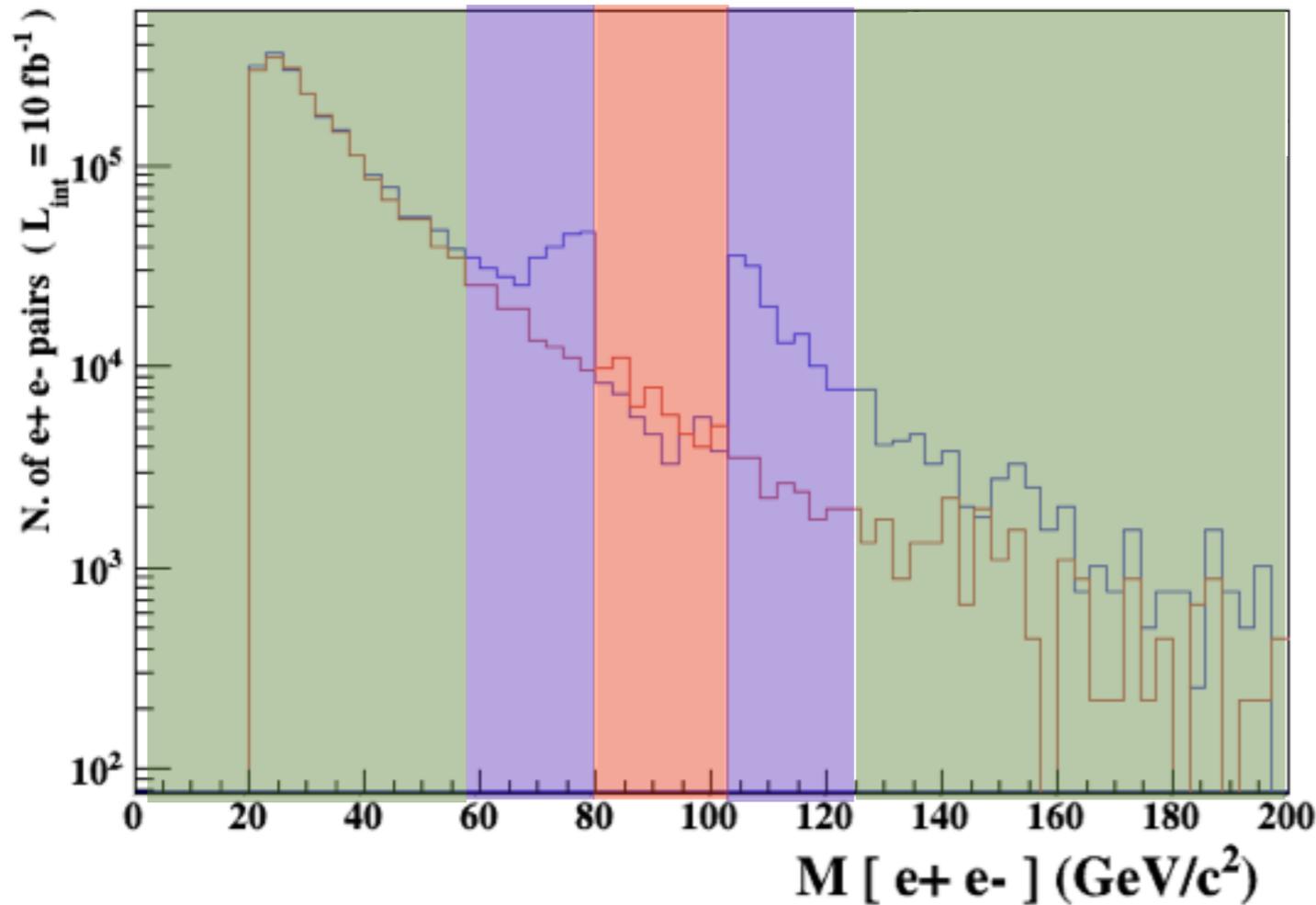
5 times width area

15 times width area

> 15 times width area

$p p > e^+ e^- / Z$
(red curve)

adding $p p > e^+ e^- \$ Z$
(blue curve)



- Z onshell veto
- In veto area only photon contribution
- area sensitive to z-peak
- very off-shell Z, the difference between the curve is due to interference which are need to be **KEPT** in simulation.

5 times width area

15 times width area

> 15 times width area

The “\$” can be use to split the sample in BG/SG area

- Syntax Like

→ $p p \rightarrow z \rightarrow e^+ e^-$

(ask one S-channel z)

→ $p p \rightarrow e^+ e^- / z$

(forbids any z)

→ $p p \rightarrow e^+ e^- \text{ $$ } z$

(forbids any z in s-channel)

- ARE NOT GAUGE INVARIANT !
- forgets diagram interference.
- can provides un-physical distributions.

- Syntax Like

- $p p \rightarrow z \rightarrow e^+ e^-$

(ask one S-channel z)

- $p p \rightarrow e^+ e^- / z$

(forbids any z)

- $p p \rightarrow e^+ e^- \text{ $$ } z$

(forbids any z in s-channel)

- ARE NOT GAUGE INVARIANT !
- forgets diagram interference.
- can provides un-physical distributions.

Avoid Those as much as possible!

- Syntax Like

→ $p p \rightarrow z \rightarrow e^+ e^-$ (ask one S-channel z)

→ $p p \rightarrow e^+ e^- / z$ (forbids any z)

→ $p p \rightarrow e^+ e^- \text{ $$ } z$ (forbids any z in s-channel)

- ARE NOT GAUGE INVARIANT !
- forgets diagram interference.
- can provides un-physical distributions.

Avoid Those as much as possible!

check physical meaning and gauge/Lorentz invariance if you do.

- Syntax like
 - $p p \rightarrow z, z \rightarrow e^+ e^-$ (on-shell z decaying)
 - $p p \rightarrow e^+ e^- \cancel{z}$ (forbids s-channel z to be on-shell)
- Are linked to cut $|M^* - M| < BW_{cut} * \Gamma$
- Are more safer to use
- **Prefer** those syntax to the previous slides one

Goal ● Handling model

Import ● `type:` `import model MC4BSM`

- Automatic switch to 4/5 flavor computation according to the mass of the b quark
- MGaMC renames sm/susy particles such that their names are identical for any model. This can be avoided by `import model MC4BSM --modelname`

Check ● `type:` `define bsm = uv uv~ ev ev~ p1 p2`
`check p p > bsm bsm`

- checks internal validity of the BSM part and consistency of the model (lorentz/gauge)

Goal • understanding decay-chain handling

Exercise

- Compare the cross-section for

```
define evdec = bsm / ev ev~
generate p p > ev ev~
output; launch
```

```
generate p p > ev ev~, ev > evdec all
output; launch
```

This is called the decay chain syntax

```
generate p p > ev > evdec all ev~
output; launch
```

- Use Automatic width computation (for all 3 cases)

```
set width wev Auto
```

To enter at the time of the edition of the cards

- Change the “cut_decays” parameter

```
set cut_decays T
```

Goal • understanding decay-chain handling

```
define bsm = bsm / ev ev~
generate p p > ev ev~
output; launch
```

```
generate p p > ev ev~, ev > bsm all
output; launch
```

```
generate p p > ev > bsm all ev~
output; launch
```

Default	Correct width	+cut_decays=T
19.7 pb	19.6 pb	19.7 pb
0.1 pb	19.3 pb	11.8 pb
0.07 pb	11.9 pb	11.9 pb

Remember

- We do not use the BR information. The cross-section depends of the total width
- particle from on shell decay do not have cut by default

Goal • present the various way to compute the width

• Check with MG the width computed with FR:

- ➔ generate uv > all all; output; launch
- ➔ generate ev > all all; output; launch
- ➔ generate p1 > all all; output; launch
- ➔ generate p2 > all all; output; launch

FR Number

0.0706 GeV

0.00497 GeV

0 GeV

0.0224 GeV

• Compare with `compute_widths bsm`

• Why the width of uv is zero here? Function called when width on Auto

• $M_{uv} = 400 \text{ GeV}$ $M_{ev} = 50 \text{ GeV}$ $\lambda = 0.1$

• $m_1 = 1 \text{ GeV}$ $m_2 = 100 \text{ GeV}$ $m_{12} = 0.5 \text{ GeV}$

Goal ● script and scan

Parameter scan:

- compute the cross-section for a couple of mass

```
generate p p > ev ev~
```

- for that you can enter for the ev mass:

```
set mev scan:[100,200, 300]
```

```
set mev scan:[100*i for i in range(1,4)]
```

Any python syntax is valid!!

scripting/ other scan:

- write in a file (./MYFILE)
- run it as ./bin/mg5_aMC ./MYFILE

```
import model MC4BSM
generate p p > ev ev~
output TUTO
launch
set nevents 5000
set LHC 13
launch
set LHC 14
```

Goal ● script and scan

Parameter scan:

- compute the cross-section for a couple of mass

```
generate p p > ev ev~
```

- for that you can enter for the ev mass:

```
set mev scan:[100,200, 300]
```

```
set mev scan:[100*i for i in range(1,4)]
```

Any python syntax is valid!!

Comment:

- ONLY for param_card entry!! Use scripting for other type of parameters (run_card,...)
- synchronized scan can be done via

```
set mev scan1:[100,200, 300]
set muv scan1:[200,300,400]
```

Three value will be computed!!

scripting/ other scan:

- write in a file (./MYFILE)
- run it as ./bin/mg5_aMC ./MYFILE

```
import model MC4BSM
generate p p > ev ev~
output TUTO
launch
  set nevents 5000
  set LHC 13
launch
  set LHC 14
```

Comment on scripting

- Do not use ./bin/mg5_aMC < ./MYFILE
- If an answer to a question is not present: **Default is taken** automatically
- **EVERYTHING** that you type can be put in the entry file

Goal ● Learn MadSpin for Onshell Decay

What is MadSpin

arXiv:1212.3460

- Program to decay on-shell particles
 - Use the NWA and the Branching-ratio
 - keep full spin-correlation
 - keep off-shell effect (up to cut-off)
 - keep unweighted event

Exercise

- generate all decay from e^+e^- pair production via MadSpin (and compare with decay-chain syntax)

Goal ● Learn MadSpin for Onshell Decay

How to

```
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=NOT INSTALLED
4 Decay particles with the MadSpin module:      madspin=OFF
5 Add weight to events based on coupling 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, 4, 5, auto, done, pythia=ON, pythia=OFF, ... ][60s to answer]
```

When you see this text, type madspin=ON

- Then edit the madspin_card and include

```
decay ev > all all
```

- You are done

Note

- Also valid for NLO processes
- sometimes faster/slower than decay-chain

Goal ● Learn loop-induced syntax

Ex. ● Compare Large stop limit and full loop

```
import model heft
generate g g > h
output; launch
```

```
import model sm
generate g g > h [QCD]
output; launch
```

```
import model sm-no_b_mass
generate g g > h [QCD]
output; launch
```

Note

- Interface fully identical to LO one
- No decay-chain/MadSpin allowed

Goal • Learn NLO syntax

Ex. • Run the pair-production at NLO

```
import model MC4BSM
generate p p > ev ev~ [QCD]
output; launch
```

Note

- Interface close but different to LO one
 - different options
 - different cuts
- No decay-chain but MadSpin allowed
- Need dedicated model (not all model valid@NLO)

The following switches determine which operations are executed:

1 Perturbative order of the calculation:	order=NLO
2 Fixed order (no event generation and no MC@[N]LO matching):	fixed_order=OFF
3 Shower the generated events:	shower=ON
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 'order=L0' at the prompt)
Type '0', 'auto', 'done' or just press enter when you are done.
[0, 1, 2, 3, 4, 5, auto, done, order=L0, ...][60s to answer]

order=L0 / order=NLO

- Use this switch to compute K-factor with the exact same settings

fixed_order=ON / fixed_order=OFF

- if ON, we perform a pure NLO computation of the cross-section — no event generation—
- if OFF, we run NLO+PS, with the MC counter-term for a given parton shower —with event generation

- I. Generate $p p > w+$ with 0 jets, 0,1 jets and 0,1,2 jets
(Each on different computers - use the most powerful computer for 0,1,2 jets)
 - a. Generate 20,000 events for a couple of different $xqcut$ values.
 - b. Compare the distributions (before and after Pythia) and cross sections (before and after Pythia) between the different processes, and between the different $xqcut$ values.
 - c. Summarize: How many jets do we need to simulate? What is a good $xqcut$ value? How are the distributions affected?

- generate the diagram with
 - ➔ generate
 - ➔ add process
- output
- launch
 - ➔ ask to run pythia
 - ➔ In run_card: put icckw=1
 - ◆ set the value for xqcut
 - ➔ In pythia_card set a value for qcut

- Qcut is the matching scale (the separation between the shower and the matrix element)
- xqcut should be strictly lower (by at least 10-15GeV) than qcut

Solution MLM

	w+0j	w+1j	w+2j	w+3j
no matching	8.35E+04	1.58E+04	8.7E+03	3.5E+03

	1GeV	10GeV	20GeV	50GeV	100GeV	500GeV
w+0	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04
0+1	1.07E+05	9.09E+04	8.91E+04	8.61E+04	8.40E+04	8.35E+04
0+1+2	1.12E+05	9.29E+04	9.03E+04	8.66E+04	8.44E+04	8.35E+04
0+1+2+3	1.20E+05	9.47E+04	9.07E+04	8.68E+04	8.40E+04	8.35E+04

	w+0j	w+1j	w+2j	w+3j
no matching	8.35E+04	1.58E+04	8.7E+03	3.5E+03

	1GeV	10GeV	20GeV	50GeV	100GeV	500GeV
w+0	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04
0+1	1.07E+05	9.09E+04	8.91E+04	8.61E+04	8.40E+04	8.35E+04
0+1+2	1.12E+05	9.29E+04	9.03E+04	8.66E+04	8.44E+04	8.35E+04
0+1+2+3	1.20E+05	9.47E+04	9.07E+04	8.68E+04	8.40E+04	8.35E+04

Slow

Fast

low efficiency

High efficiency

	w+0j	w+1j	w+2j	w+3j
no matching	8.35E+04	1.58E+04	8.7E+03	3.5E+03

	1GeV	10GeV	20GeV	50GeV	100GeV	500GeV
w+0	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04
0+1	1.07E+05	9.09E+04	8.91E+04	8.61E+04	8.40E+04	8.35E+04
0+1+2	1.12E+05	9.29E+04	9.03E+04	8.66E+04	8.44E+04	8.35E+04
0+1+2+3	1.20E+05	9.47E+04	9.07E+04	8.68E+04	8.40E+04	8.35E+04

- No effect of the matching for 0 jet sample.

	w+0j	w+1j	w+2j	w+3j
no matching	8.35E+04	1.58E+04	8.7E+03	3.5E+03

	1GeV	10GeV	20GeV	50GeV	100GeV	500GeV
w+0	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04
0+1	1.07E+05	9.09E+04	8.91E+04	8.61E+04	8.40E+04	8.35E+04
0+1+2	1.12E+05	9.29E+04	9.03E+04	8.66E+04	8.44E+04	8.35E+04
0+1+2+3	1.20E+05	9.47E+04	9.07E+04	8.68E+04	8.40E+04	8.35E+04

- matching scale too high only the 0 jet sample contributes => all radiations are from pythia

	w+0j	w+1j	w+2j	w+3j
no matching	8.35E+04	1.58E+04	8.7E+03	3.5E+03

	1GeV	10GeV	20GeV	50GeV	100GeV	500GeV
w+0	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04
0+1	1.07E+05	9.09E+04	8.91E+04	8.61E+04	8.40E+04	8.35E+04
0+1+2	1.12E+05	9.29E+04	9.03E+04	8.66E+04	8.44E+04	8.35E+04
0+1+2+3	1.20E+05	9.47E+04	9.07E+04	8.68E+04	8.40E+04	8.35E+04

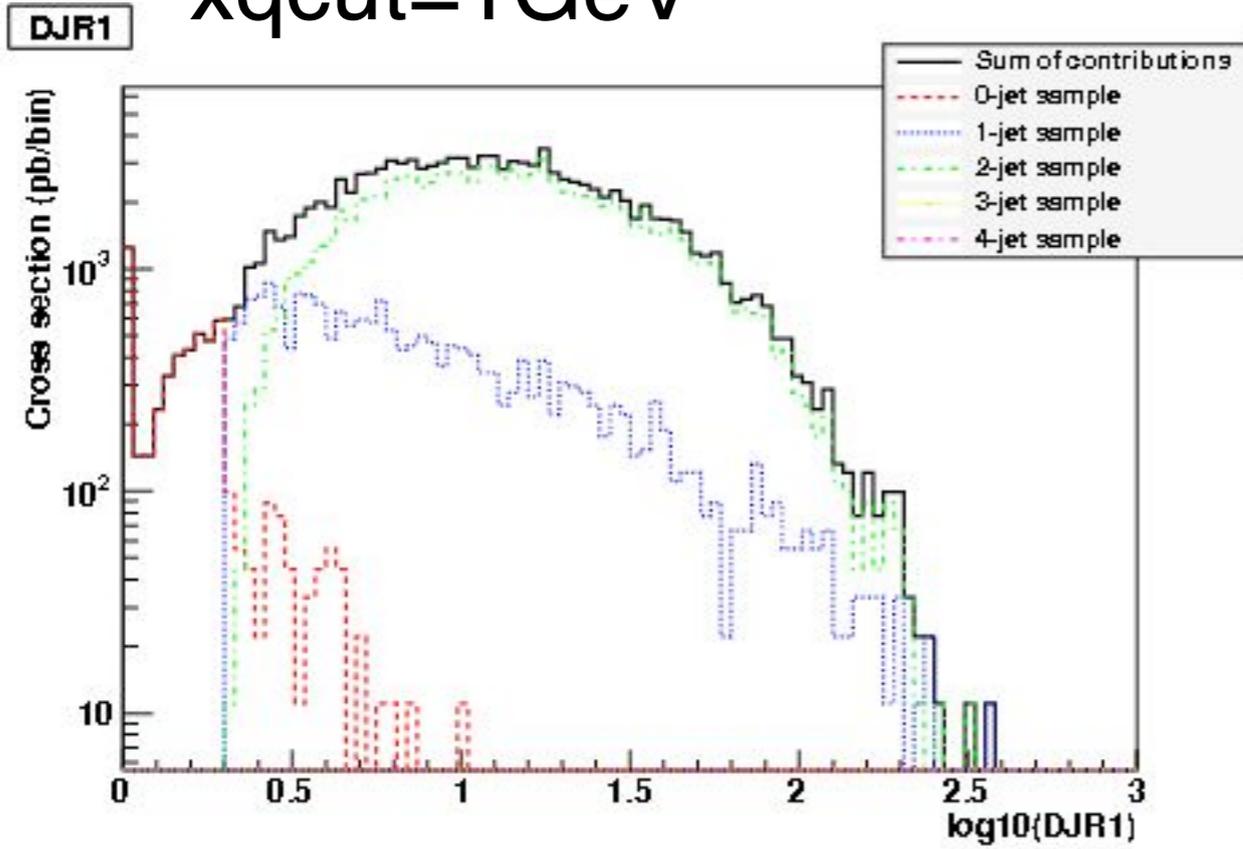
- matching scale too low. Only highest multiplicity sample contributes and low efficiency

	w+0j	w+1j	w+2j	w+3j
no matching	8.35E+04	1.58E+04	8.7E+03	3.5E+03

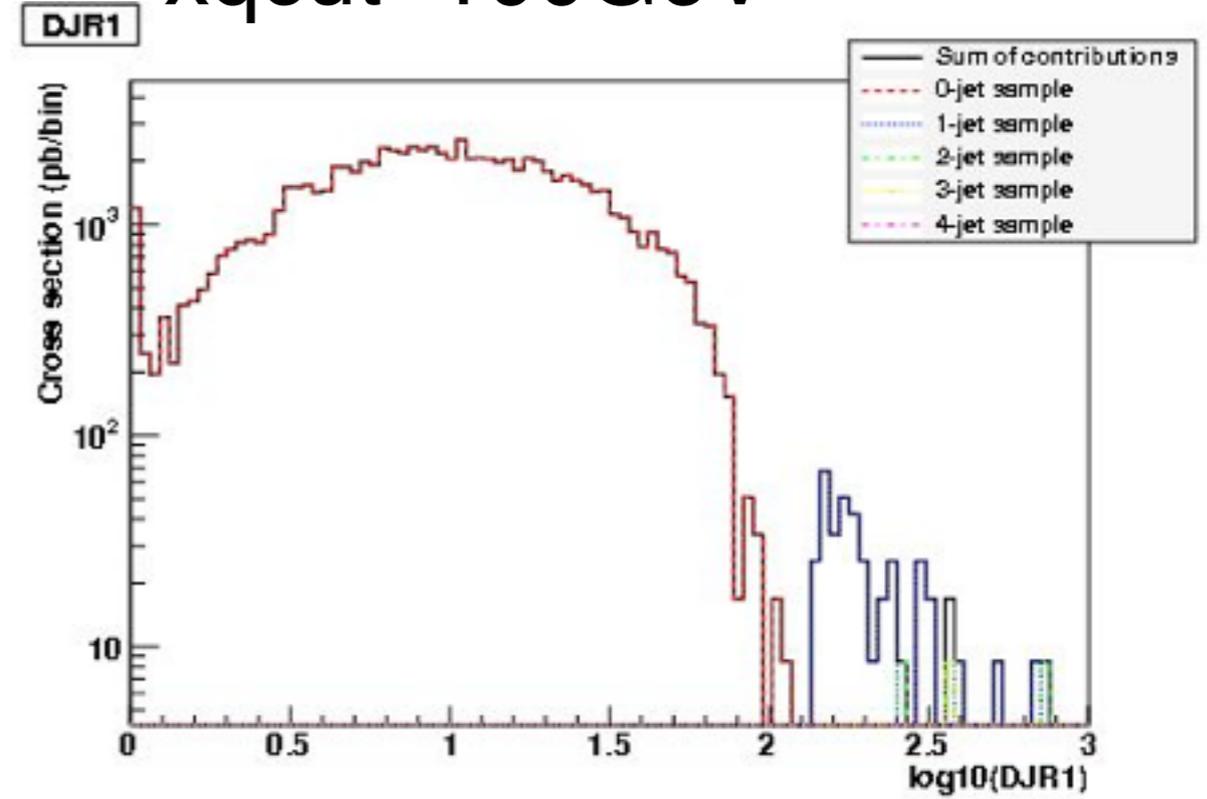
	1GeV	10GeV	20GeV	50GeV	100GeV	500GeV
w+0	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04
0+1	1.07E+05	9.09E+04	8.91E+04	8.61E+04	8.40E+04	8.35E+04
0+1+2	1.12E+05	9.29E+04	9.03E+04	8.66E+04	8.44E+04	8.35E+04
0+1+2+3	1.20E+05	9.47E+04	9.07E+04	8.68E+04	8.40E+04	8.35E+04

- Wrong differential rate plot. so to discard.

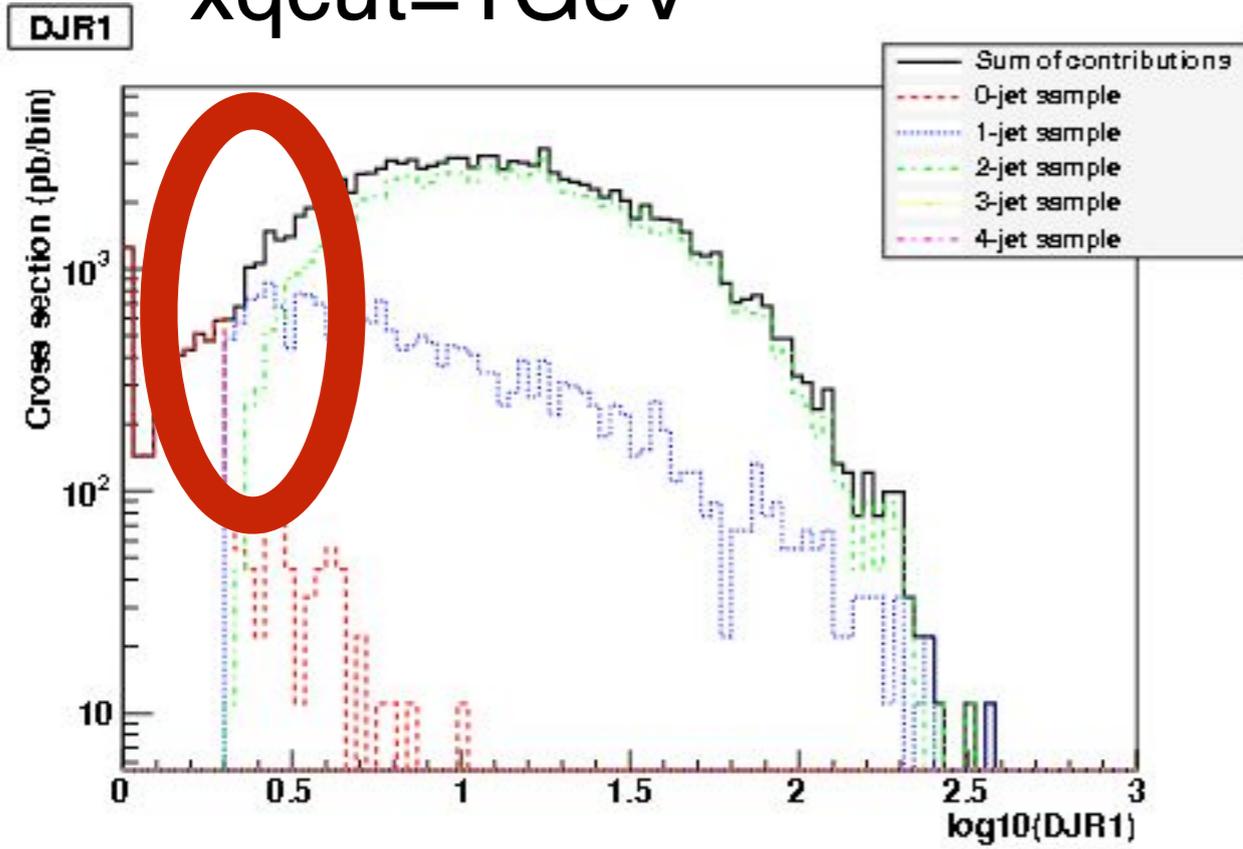
xqcut=1 GeV



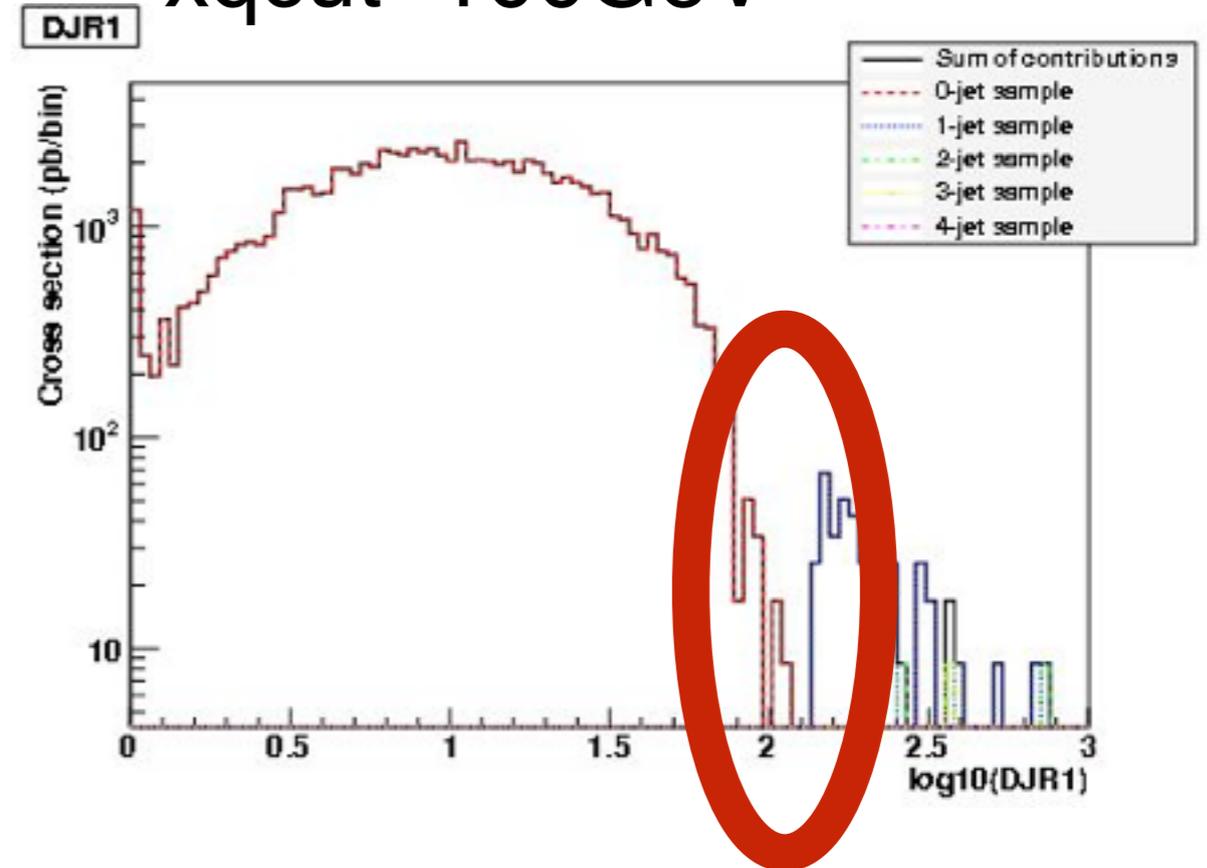
xqcut=100 GeV



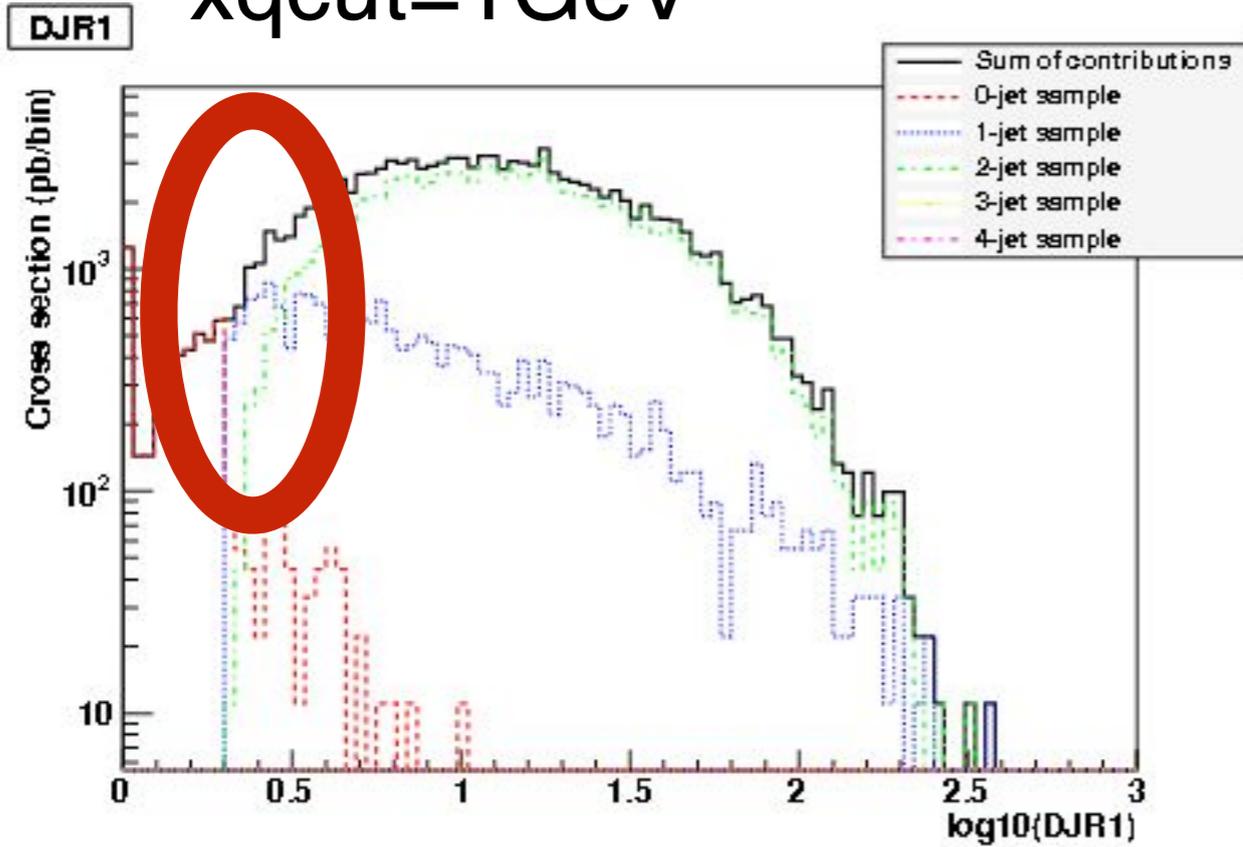
xqcut=1 GeV



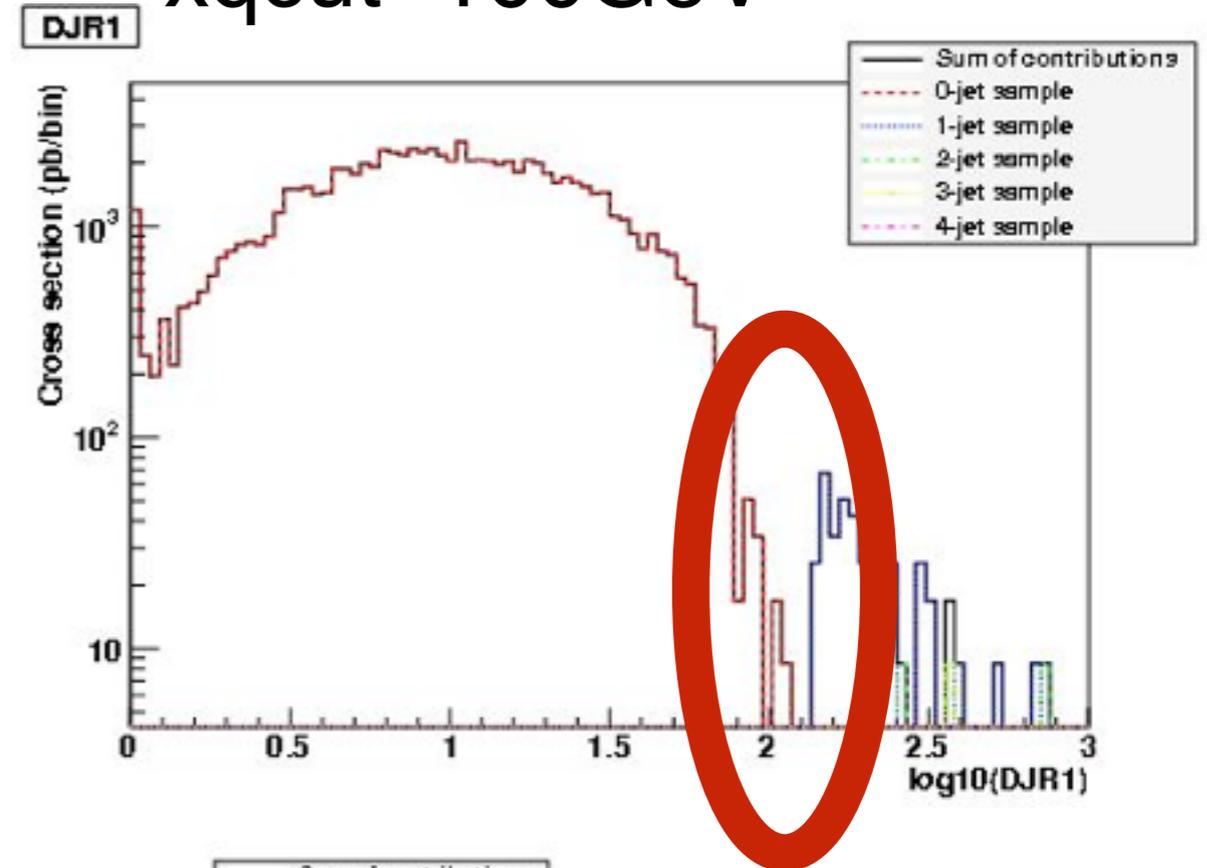
xqcut=100 GeV



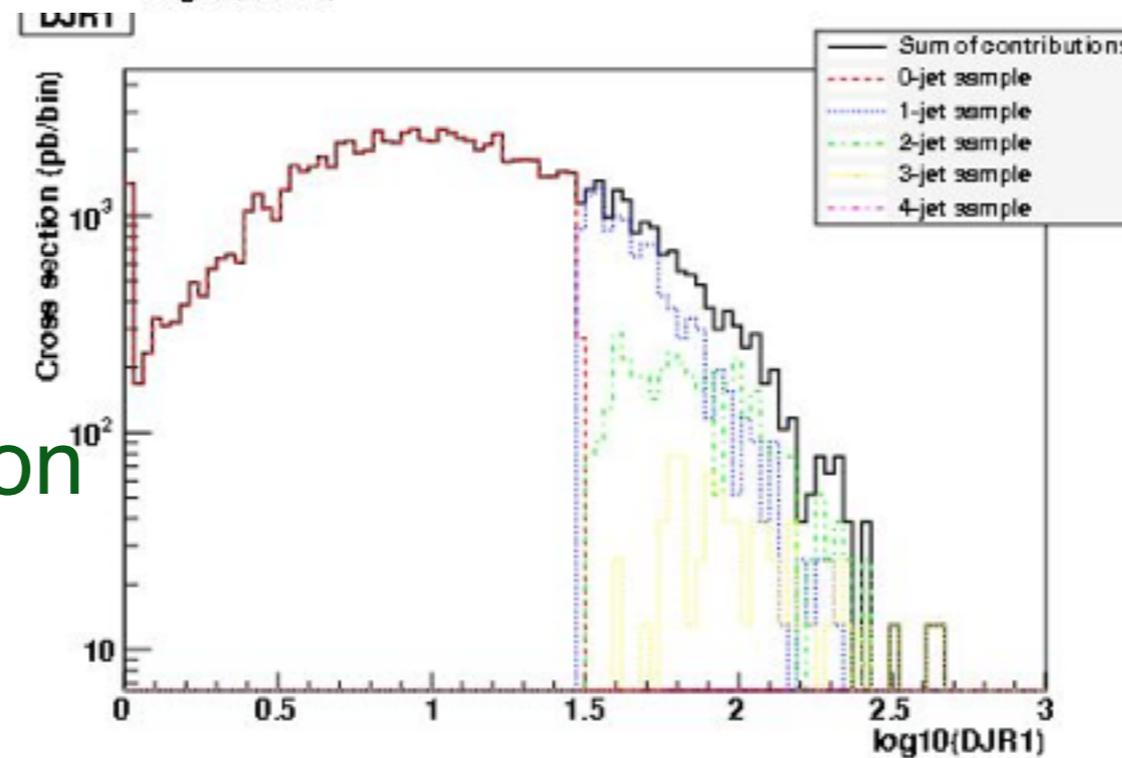
xqcut=1 GeV



xqcut=100 GeV



xqcut=20 GeV
 smooth transition



	w+0j	w+1j	w+2j	w+3j
no matching	8.35E+04	1.58E+04	8.7E+03	3.5E+03

	1GeV	10GeV	20GeV	50GeV	100GeV	500GeV
w+0	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04	8.35E+04
0+1	1.07E+05	9.59E+04	8.91E+04	8.61E+04	8.40E+04	8.35E+04
0+1+2	1.12E+05	9.29E+04	9.03E+04	8.66E+04	8.44E+04	8.35E+04
0+1+2+3	1.20E+05	9.17E+04	9.07E+04	8.68E+04	8.40E+04	8.35E+04

- Relatively stable cross-section! Important check.
- Close to the unmatched 0j cross-section