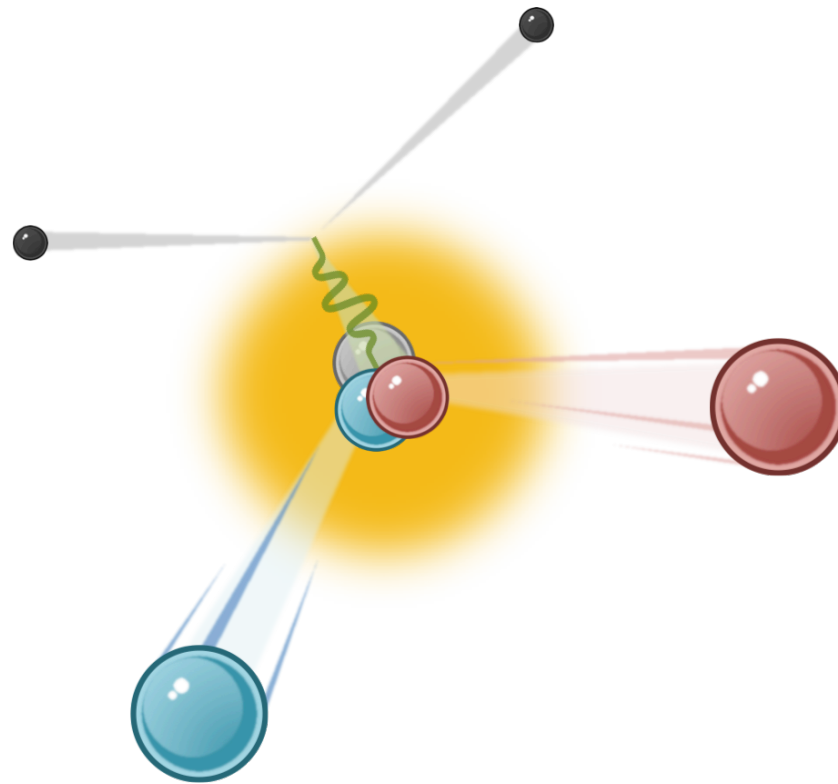
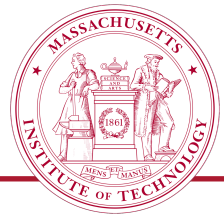


Implementing ^3He and ^3H spectral functions in SIMC



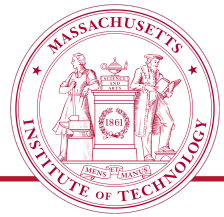
Reynier Cruz Torres,
for the E12-14-011 collaboration
September 19, 2016

Outline



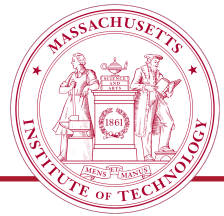
1. SIMC overview
2. ^3He spectral function
3. Implementation
4. Results
5. Next steps

Outline



1. SIMC overview
2. ^3He spectral function
3. Implementation
4. Results
5. Next steps

SIMC overview

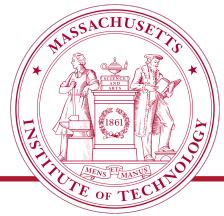


- ❑ SIMC is the standard Hall C Monte Carlo for coincidence reactions (similar to MCEEP) written in FORTRAN (now gfortran compatible).

- ❑ Features:
 - Optics (COSY) and “aperture checking” Monte Carlo of spectrometers
 - Includes radiative effects, multiple scattering, ionization energy loss, particle decay
 - Simple prescriptions available for FSI, Coulomb correction, ...

*Taken from: http://hallaweb.jlab.org/data_reduc/AnaWork2009/simc_overview.pdf

SIMC overview



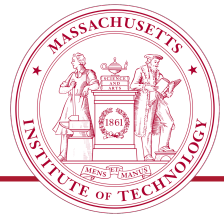
☐ Reactions implemented:



1. Elastic and quasielastic $\rightarrow H(e,e'p), A(e,e'p)$
2. Exclusive pion production $\rightarrow H(e,e'\pi^+)n, A(e,e'\pi^{+/-})$ [quasifree or coherent]
3. Kaon electroproduction $\rightarrow H(e,e'K^+)\Lambda,\Sigma, A(e,e'K^{+/-})$
4. $H(e,e'\pi^{+/-})X, D(e,e'\pi^{+/-})X$ [semi-inclusive]
5. $H(e,e'K^{+/-})X, D(e,e'K^{+/-})X$ [semi-inclusive]
6. $H(e,e'\rho \rightarrow \pi^+\pi^-)p, D(e,e'\rho \rightarrow \pi^+\pi^-)$ [diffractive ρ]

*Taken from: http://hallaweb.jlab.org/data_reduc/AnaWork2009/simc_overview.pdf

Outline



1. SIMC overview
2. ^3He spectral function
3. Implementation
4. Results
5. Next steps

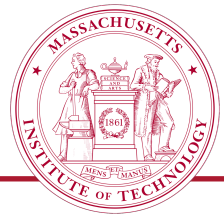
^3He spectral function



- ❑ File is already in the SIMC main directory (benharsf_3mod.dat)
- ❑ It is a lookup table

9.866	5.600	0.1261E-01	0.0000E+00	19.733	1.000
9.866	10.300	0.2382E-02	0.3563E-02	19.733	5.000
9.866	15.300	0.7343E-03	0.1098E-02	19.733	5.000
9.866	20.300	0.2631E-03	0.3935E-03	19.733	5.000
9.866	25.300	0.1097E-03	0.1641E-03	19.733	5.000
...

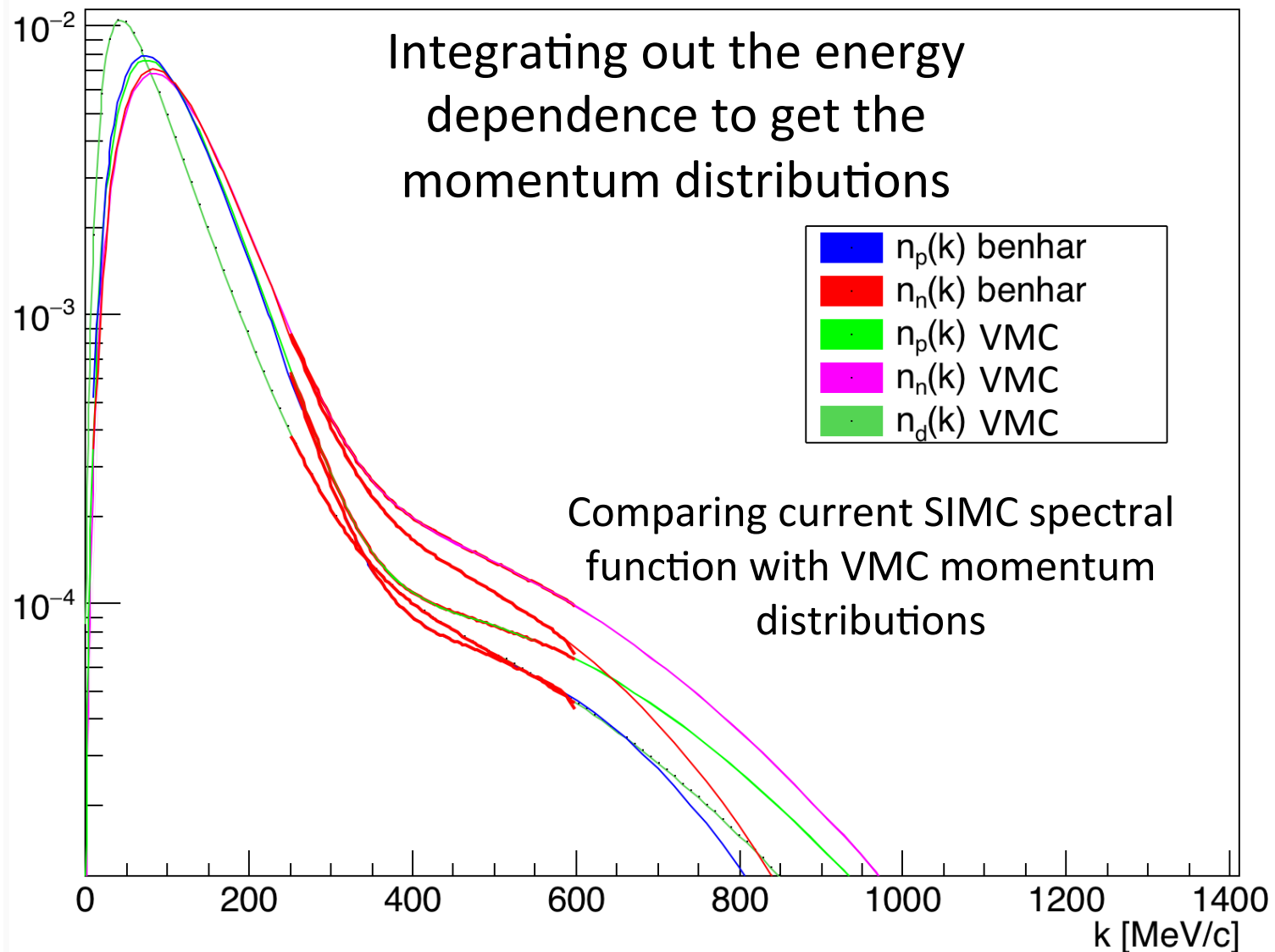
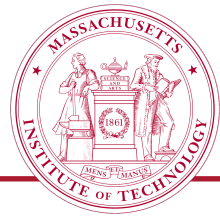
^3He spectral function



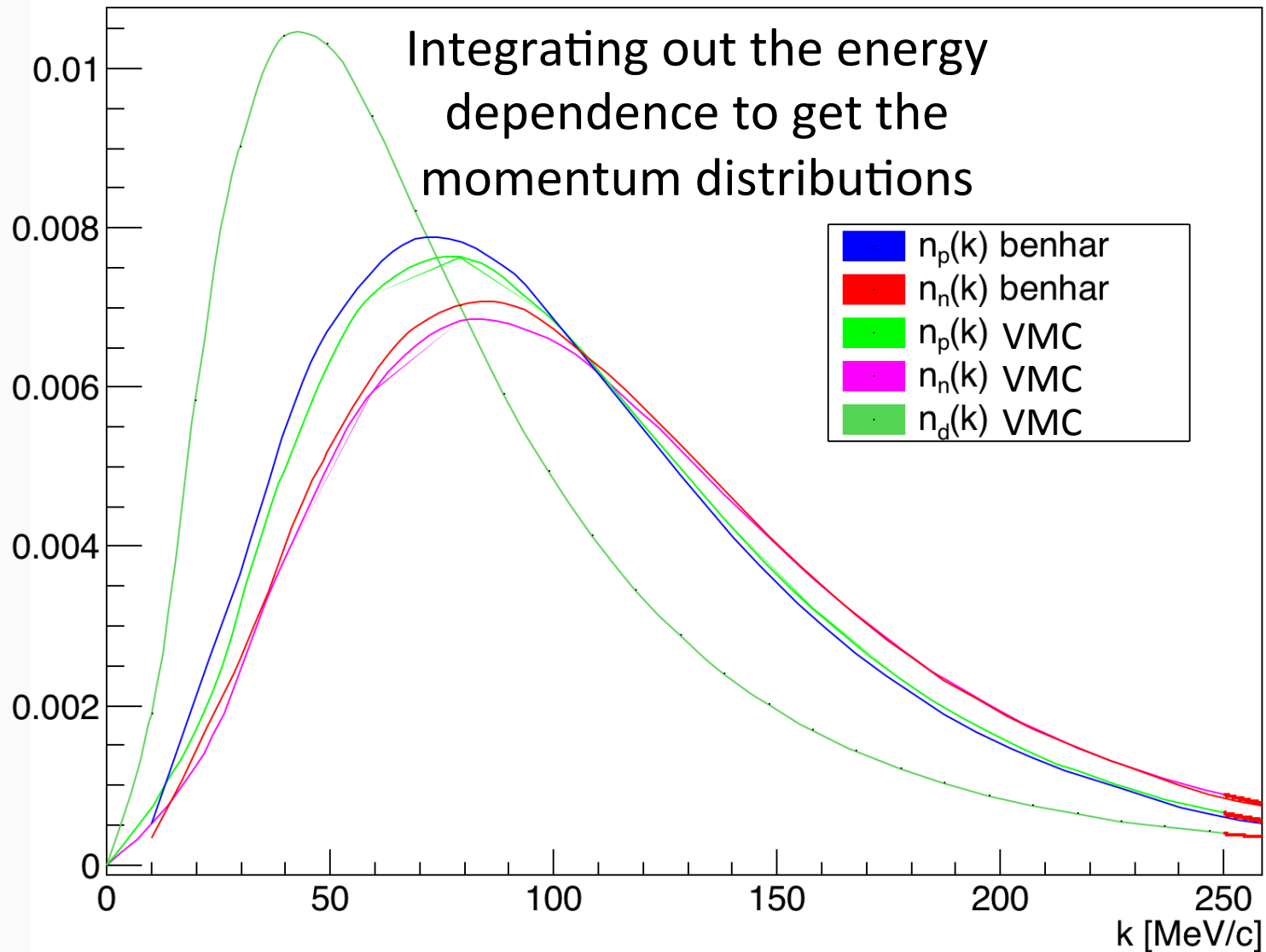
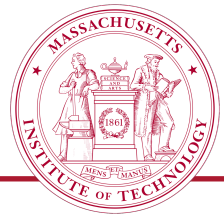
- ❑ File is already in the SIMC main directory (benharsf_3mod.dat)
- ❑ It is a lookup table

9.866	5.600	0.1261E-01	0.0000E+00	19.733	1.000
9.866	10.300	0.2382E-02	0.3563E-02	19.733	5.000
9.866	15.000	0.3443E-03	0.1098E-02	19.733	5.000
9.866	20.300	0.2531E-03	0.3935E-03	19.733	5.000
9.866	25.300	0.1097E-03	0.1641E-03	19.733	5.000
...

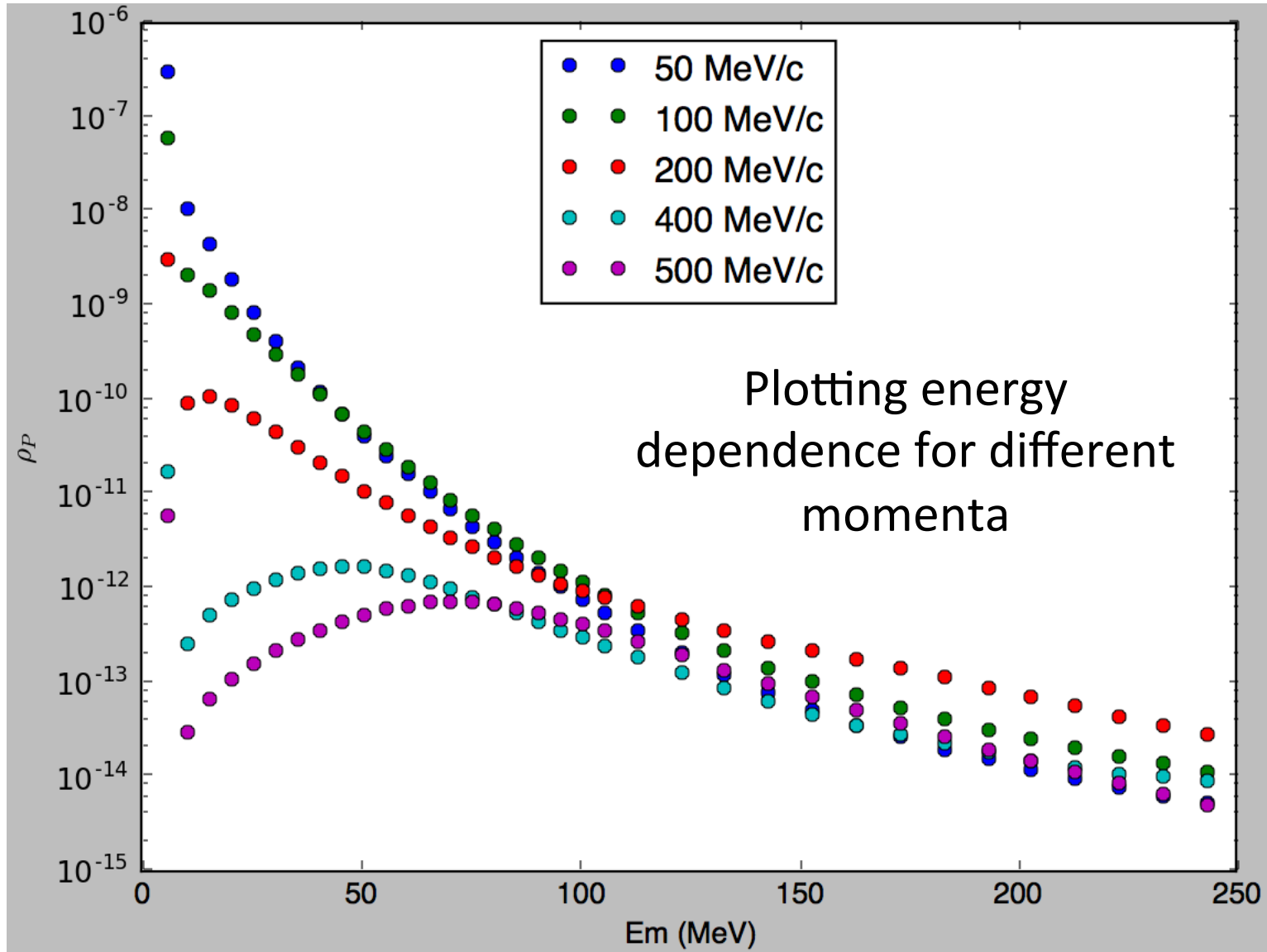
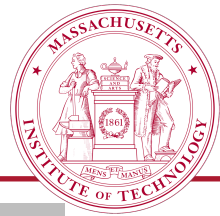
^3He spectral function



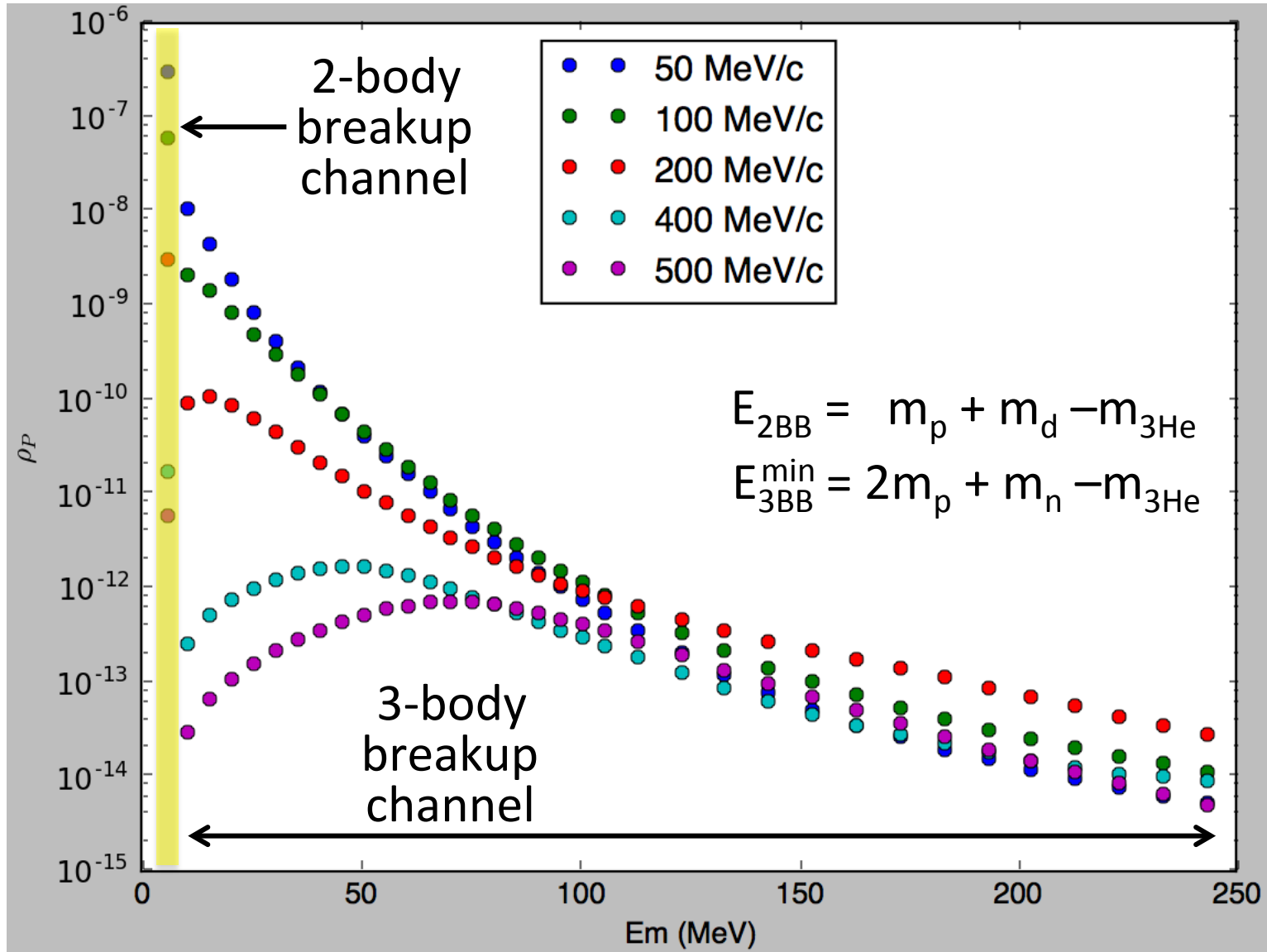
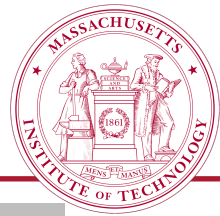
^3He spectral function



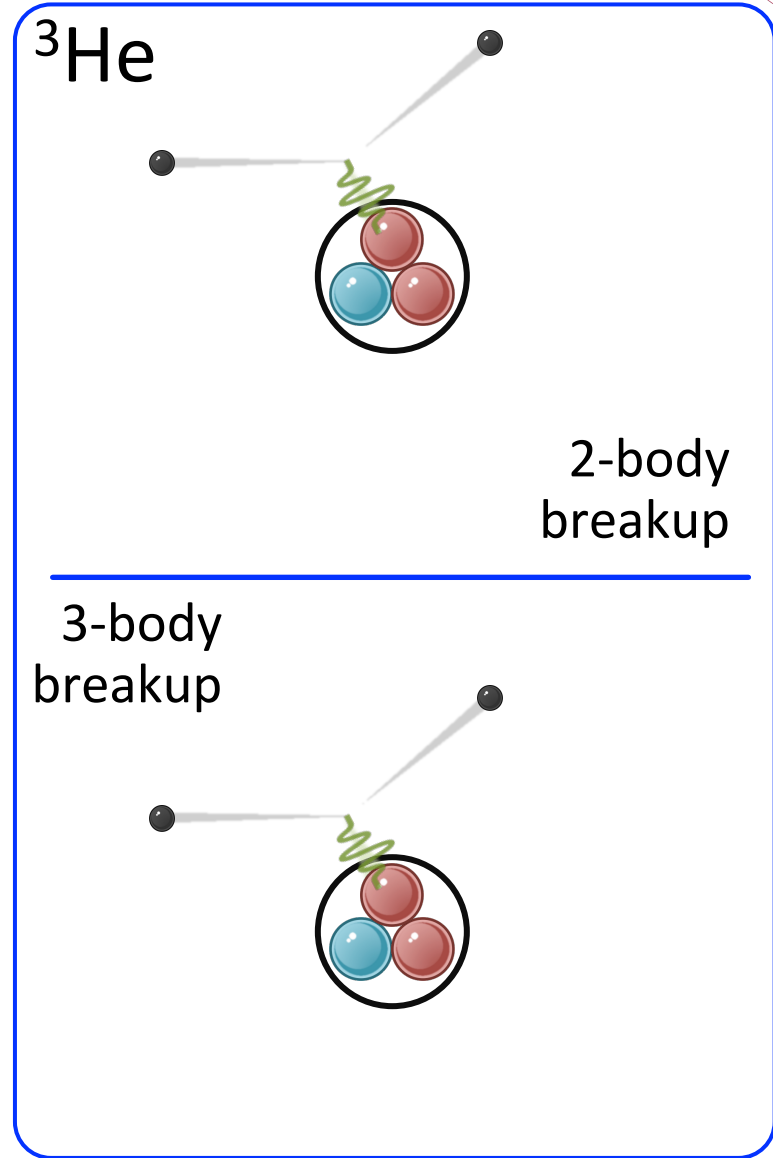
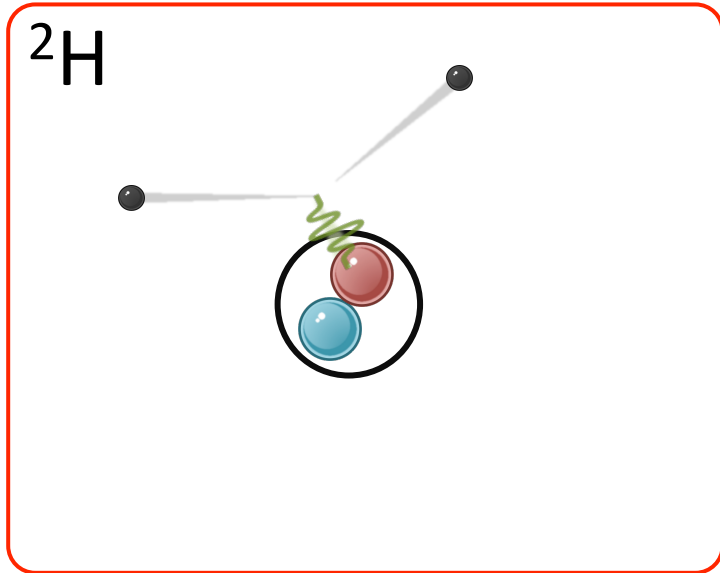
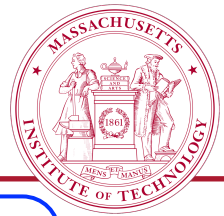
^3He spectral function



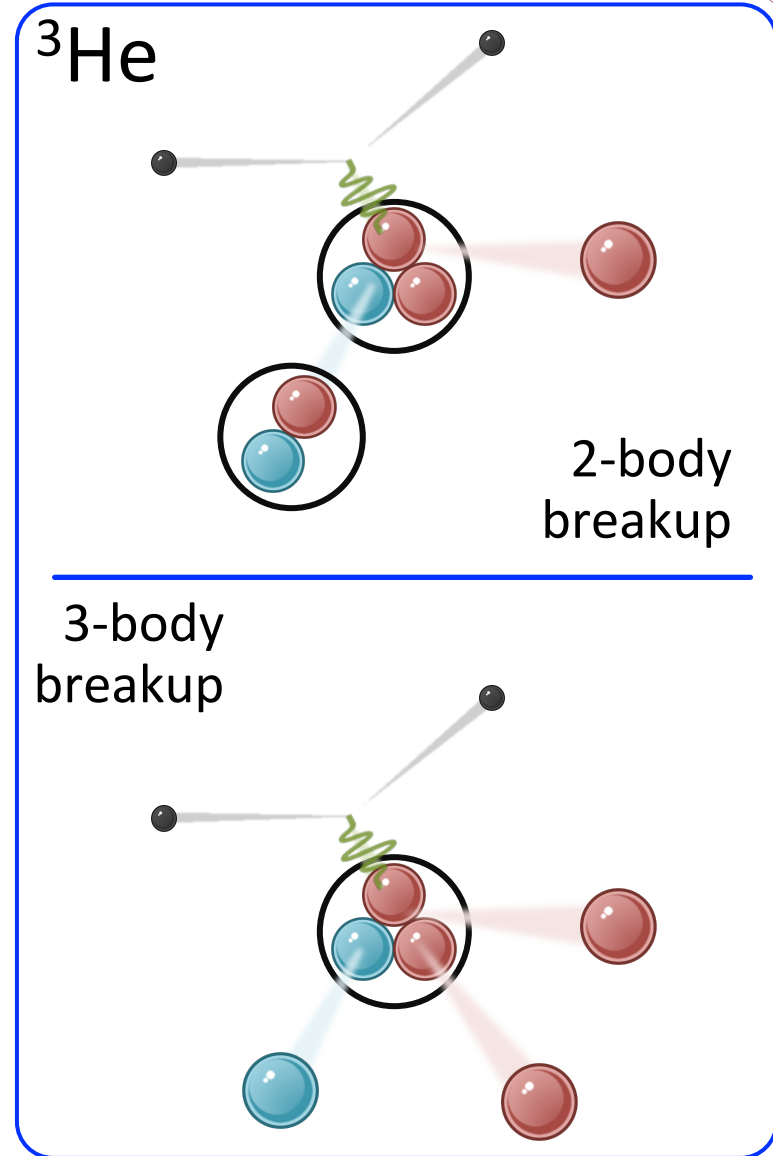
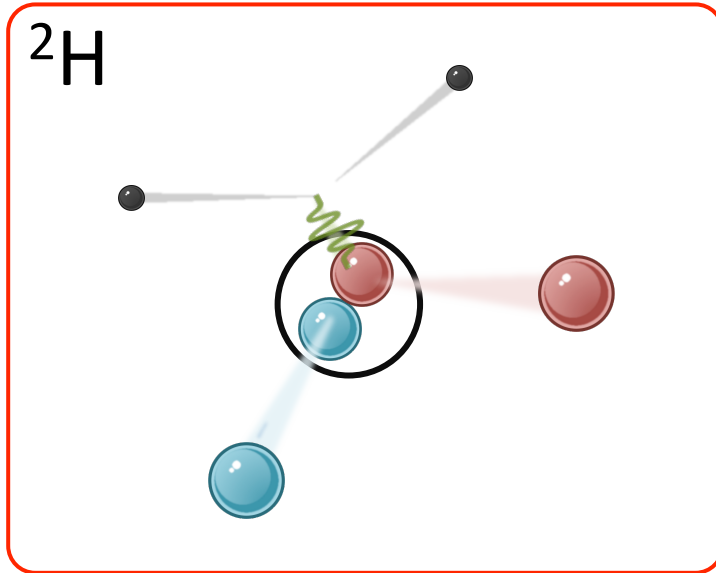
^3He spectral function



2- and 3-body breakup channels



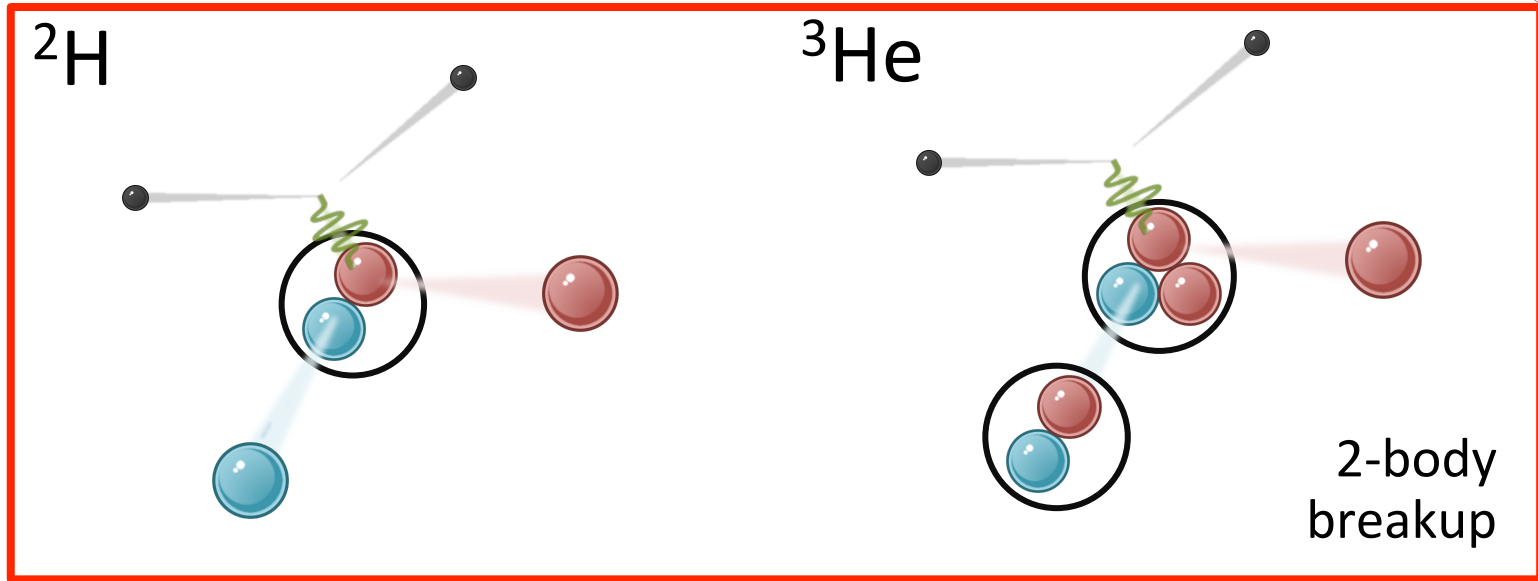
2- and 3-body breakup channels



□ Kinematics in ${}^3\text{He}$ 2-body breakup and 3-body breakup channels are not the same.

□ ${}^3\text{He}$ 3-body breakup channel has one extra dof.

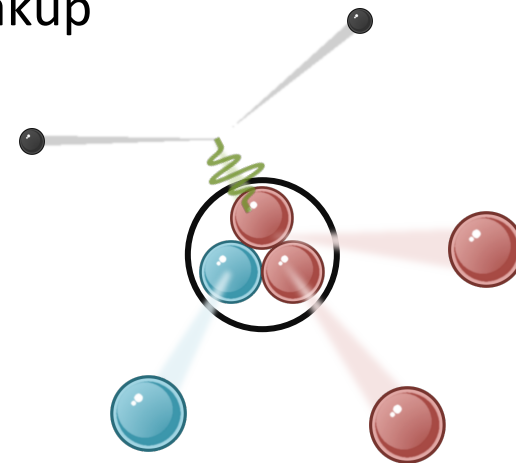
2- and 3-body breakup channels



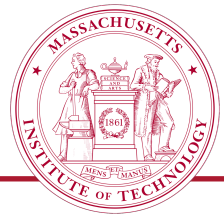
❑ Kinematics in ${}^3\text{He}$ 2-body breakup and ${}^2\text{H}$ channels are the same.

❑ Need to take this into account in SIMC

3-body breakup

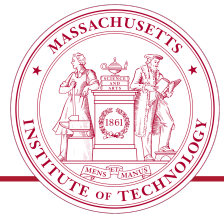


Outline



1. SIMC overview
2. ^3He spectral function
3. Implementation
4. Results
5. Next steps

Implementation



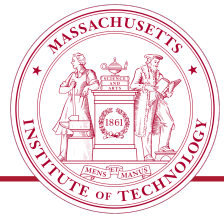
Started using Werner Boeglin's version of SIMC (slightly different from the standard version).

1. Removed a few lines of code that were blocking the spectral function in event.f
2. Created a new (boolean) parameter in the input file called "doing_bound" that only affects the ^3He sections of the code:

- doing_bound = T \rightarrow 2-body breakup channel
- doing_bound = F \rightarrow 3-body breakup channel

* To run SIMC using the spectral function, run twice: once with doing_bound = T, and once with doing_bound = F. Then combine the output root files.

Implementation



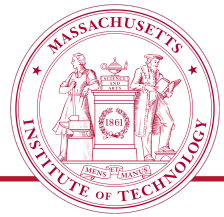
3. Implemented the following changes (since the kinematics in the kinematics in the ^3He 2-body breakup channel are deuteron-like):

from:	to:
doing_deuterium	doing_deuterium or (doing_heavy and doing_bound)
doing_heavy	doing_heavy and <u>NOT</u> doing_bound

* This last change is not true for every single doing_deuterium/doing_heavy statement.

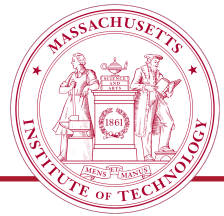
** Please, contact me for the specifics (reynier@mit.edu)

Outline



1. SIMC overview
2. ^3He spectral function
3. Implementation
4. Results
5. Next steps

Results



- ✓ Ran SIMC in the two kinematical configurations that we have in our experiment for deuteron (standard way) and ^3He (using the implemented spectral function):

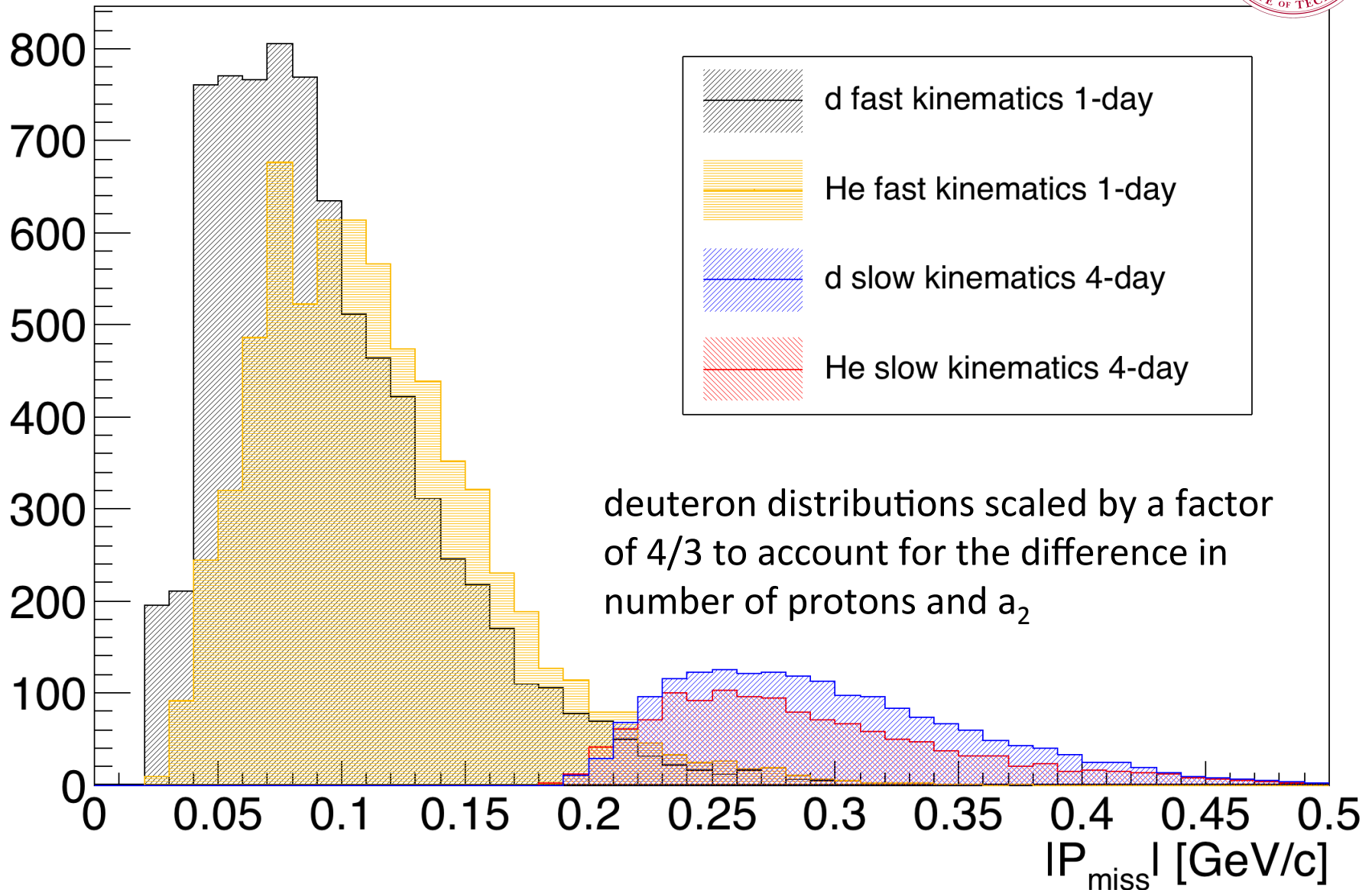
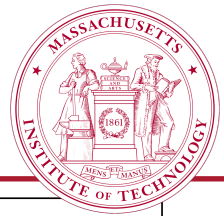
	$\langle p_m \rangle$ (MeV/c)	x	E_e (GeV)	θ_e	p_p (GeV/c)	θ_p	total beam time (days)
“Fast kinematics”	100	1.15	3.47	20.86°	1.607	48.67°	1
“Slow kinematics”	300	1.41	3.64	20.35°	1.352	58.55°	8

- ✓ Other parameters:

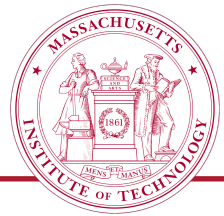
1. Beam current $I = 25 \mu\text{A}$
2. Luminosity $L = 1.88 \mu\text{A}\cdot\text{g}/\text{cm}^2$
3. Target thickness $t = 75 \text{ mg}/\text{cm}^2$
4. Target density $\rho = 0.003 \text{ g}/\text{cm}^3$
5. Target length $l = 25 \text{ cm}$.

- ✓ Cuts applied in the analysis stage: $\theta_{nq} < 40^\circ$ to reduce FSI

Results



Outline

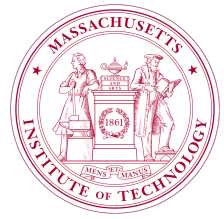


1. SIMC overview
2. ^3He spectral function
3. Implementation
4. Results
5. Next steps

Next steps



- ✓ Fix the normalizations:
 - Currently working on it
- ✓ Add tritium spectral function:
 - Also working on this. For the moment we will be using the n-distribution from ${}^3\text{He}$ which (from isospin symmetry) is effectively the same as the p-distribution for tritium
- ✓ Add other versions of the spectral function
- ✓ Kinematics optimization
- ✓ Calibration
- ✓ Any others?



Thank you