

## Background

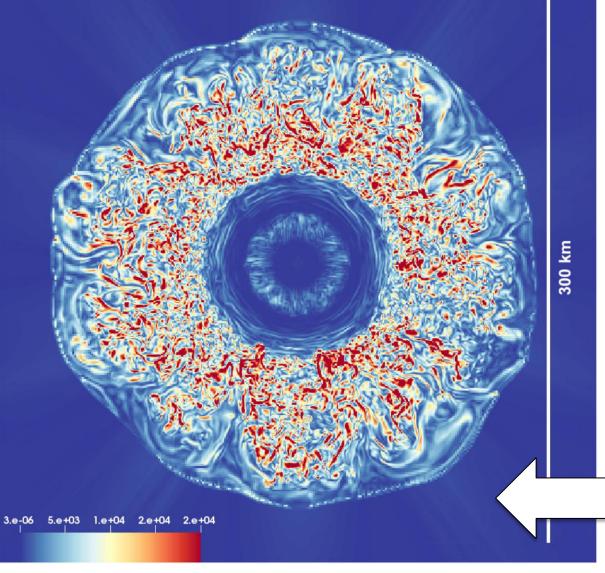
**Core-collapse** supernovae (CCSNe) cosmic are explosions caused by the gravitational collapse of stars with mass greater than eight solar masses (>8  $M_{\odot}$ ). Running computer simulations of supernovae can give valuable insight into the physical processes that cause these phenomena. While three-dimensional (3D) simulations are the most physically accurate models, they are extremely computationally expensive. To save computing power, **1D** simulations are often used instead.

1D simulations cannot naturally exhibit phenomena of **convection** and **turbulence**, which are inherently multidimensional. Instead, we introduce these effects artificially based on the Navier-Stokes Equations. The equations require the correct convective **mixing-length** parameter and turbulent mixing parameters to produce 1D simulations that run smoothly and evolve in a similar way as 3D simulations.

This study tested simulations with different parameters in order to best fit 1D simulations to 3D sample simulations.

# Methods

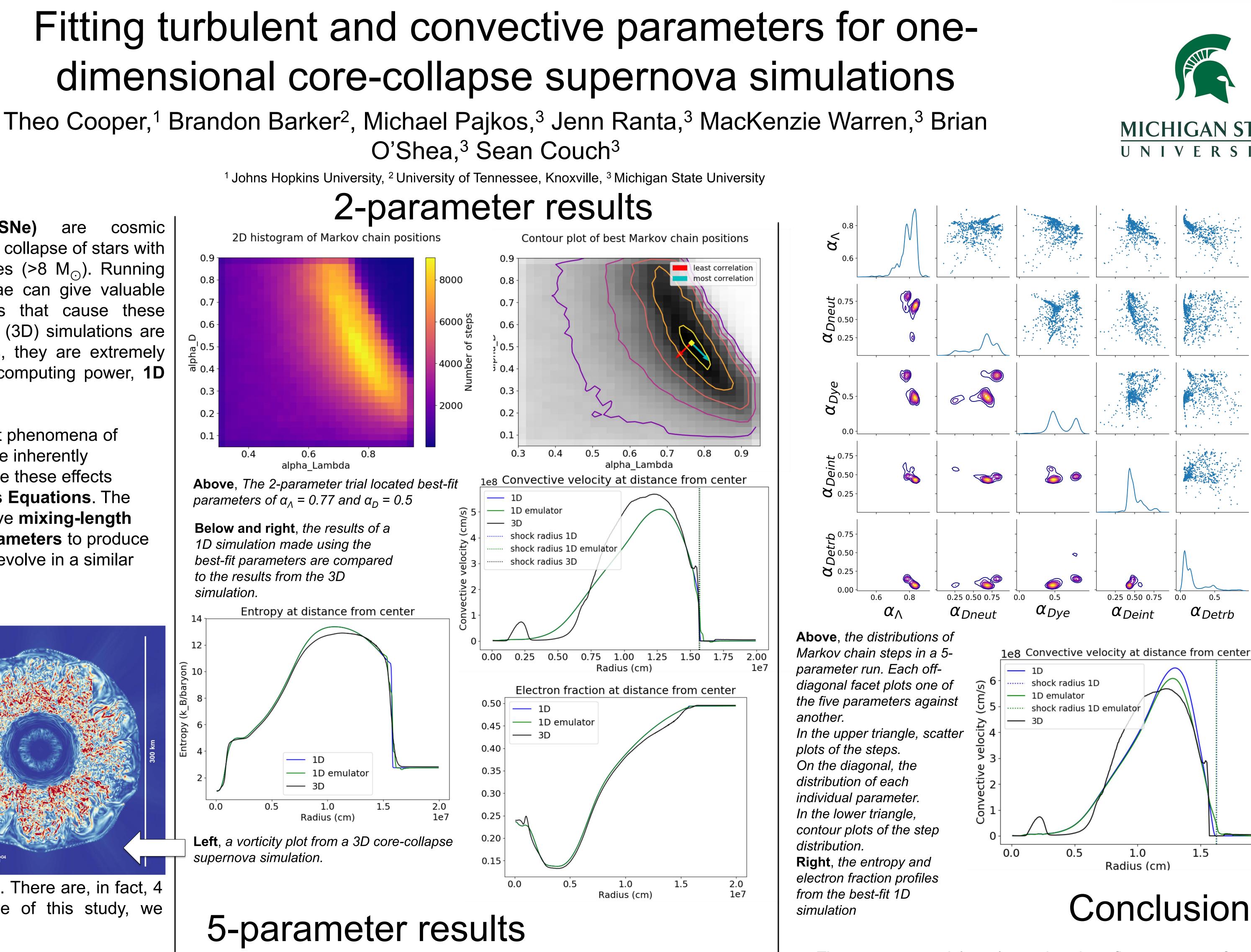
The parameters being varied were the mixinglength parameter  $\alpha_{\Lambda}$ 



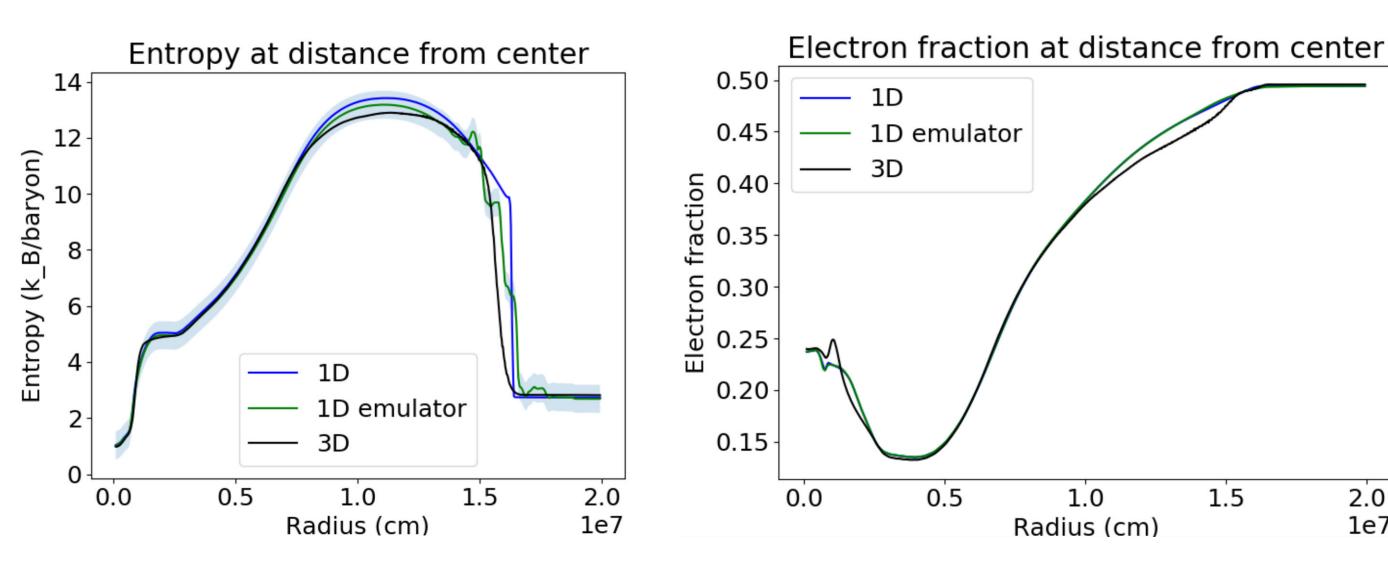
and the turbulent mixing parameter  $\alpha_{\rm p}$ . There are, in fact, 4 separate  $\alpha_{\rm D}$ 's, but in the first phase of this study, we assume they are equal.

This study used Markov Chain Monte Carlo (MCMC) techniques to determine the best-fit parameters. In these methods, one or more "walkers" are placed somewhere in the parameter space, and are allowed to take "steps" in various directions. Each walker bases these steps on the goodness of fit of the new and old sets of parameters. The best fits correspond to the areas where the walkers spend the most steps. The 1D profiles at each step were provided by a Gaussian process emulator that interpolated from sample points.

The first phase of this study used a two-dimensional parameter space, while the second phase used five parameters:  $\alpha_{\Lambda}$ ;  $\alpha_{D, v}$ ;  $\alpha_{D, Ye}$ ;  $\alpha_{D, E int}$ ;  $\alpha_{D, E turb}$ 







**Above**, the entropy and electron fraction profiles from the best-fit 1D simulation from the 5parameter trial, compared to the 3D simulation



The 5-parameter trial results produce best-fit parameters of:

$\alpha_{\wedge}$	α <sub>D, ν</sub>	$a_{D,Ye}$	$\alpha_{D,E}$ int	$a_{D,E}$
0.87	0.83	0.26	0.43	0.06

Under the 2-parameter trial's assumption that all  $\alpha_{\rm D}$  mixing parameters are equal, the best-fit parameters become: •  $\alpha_{\Lambda} = 0.77$ 

•  $a_{\rm D} = 0.5$ 

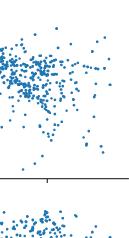
2.0

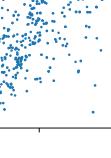
1e7

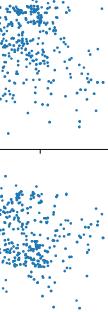
In future trials, this method will be applied to other 3D simulation data to corroborate these results.

### Acknowledgements

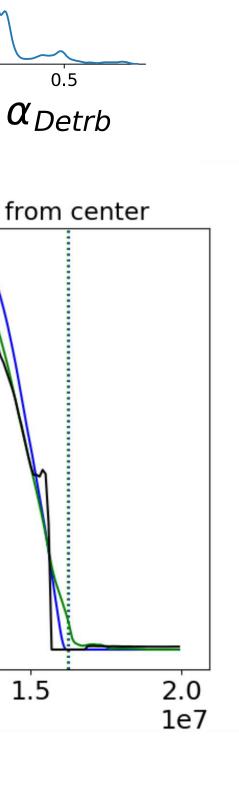
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