

## Poster location: P13



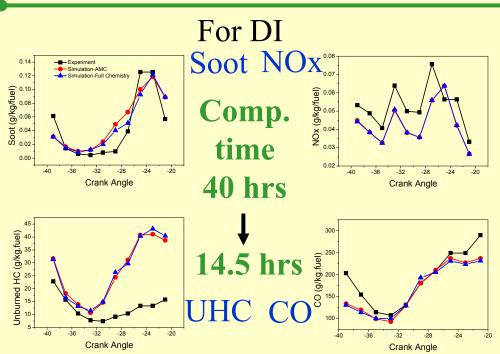
# An adaptive multi-grid chemistry (AMC) model for efficient simulation of HCCI and DI engine combustion

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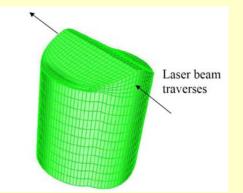
Sponsor: the U.S. Department of Energy

## Summary of the study

Objective: Develop an efficient chemistry model for simulation of HCCI and DI engine combustion within the KIVA3v2 framework Approach: Use an adaptive multi-grid technique to group thermodynamically-similar cells in order to reduce the calling frequency to the chemistry solver

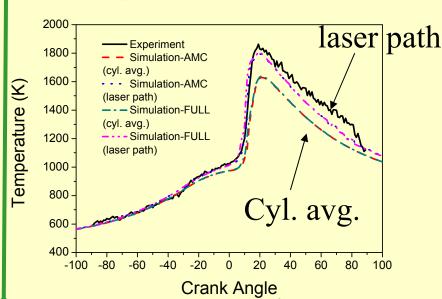


#### For HCCI



Compare with FTIR measurement

### Comp. time 50 hrs $\rightarrow$ 5 hrs





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