

Spectral Analysis of Al/MgF₂ Foils Heated by Z-pinch Radiation

October 30, 2001

Gregory A. Rochau, J. E. Bailey
Sandia National Laboratories, Albuquerque, NM
J. J. MacFarlane
Prism Computational Sciences, Madison, WI



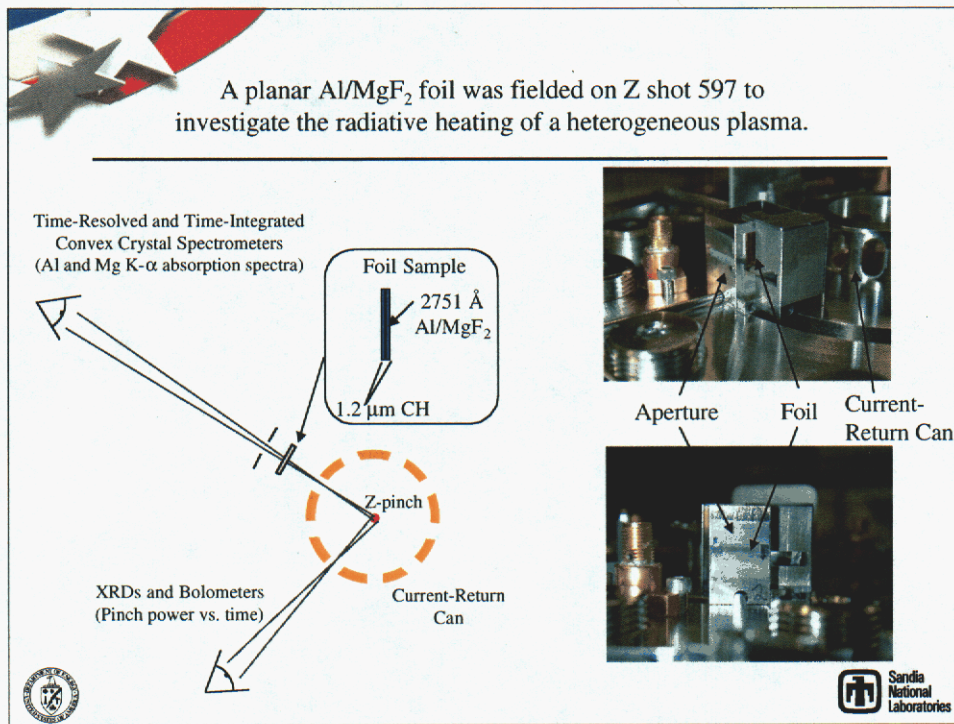
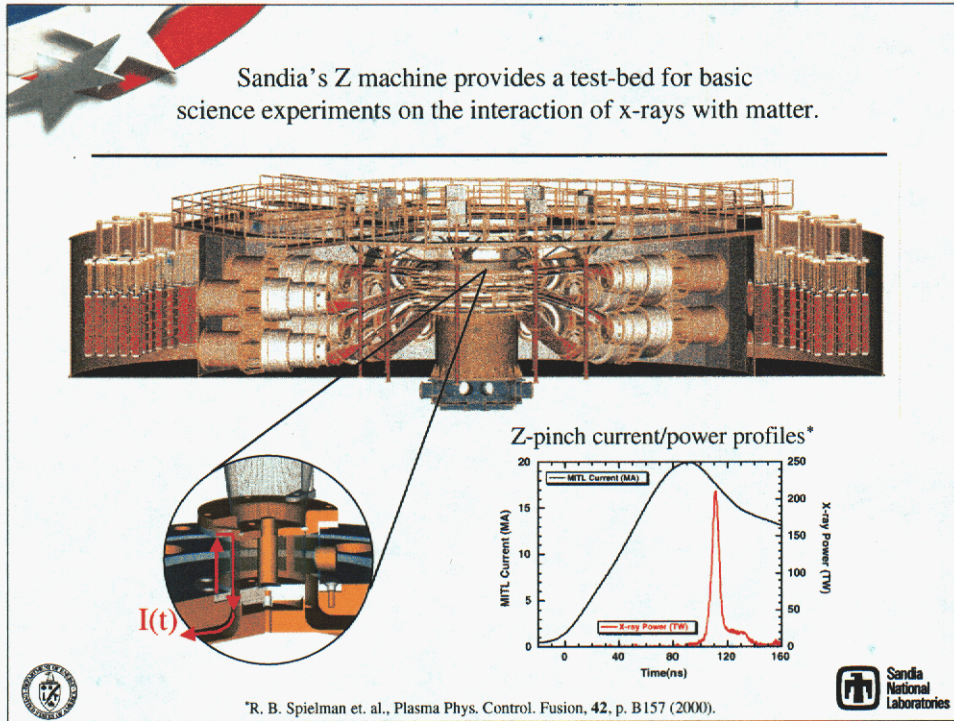
Sandia is a multiprogram laboratory operated by Sandia Corporation, a Lockheed Martin Company, for the United States Department of Energy under contract DE-AC04-94AL85000.

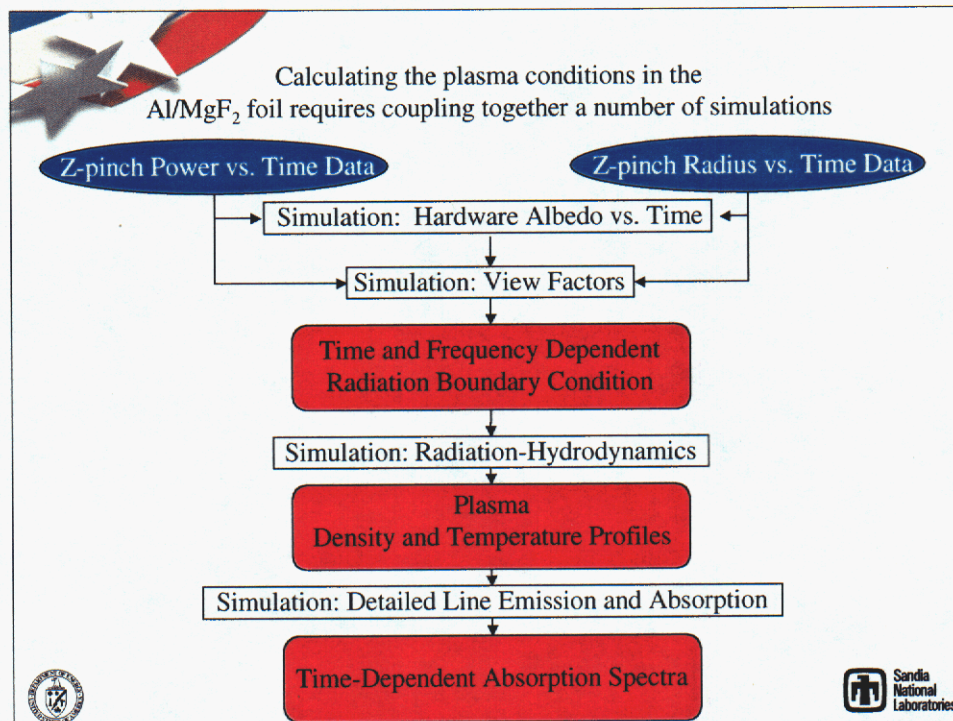
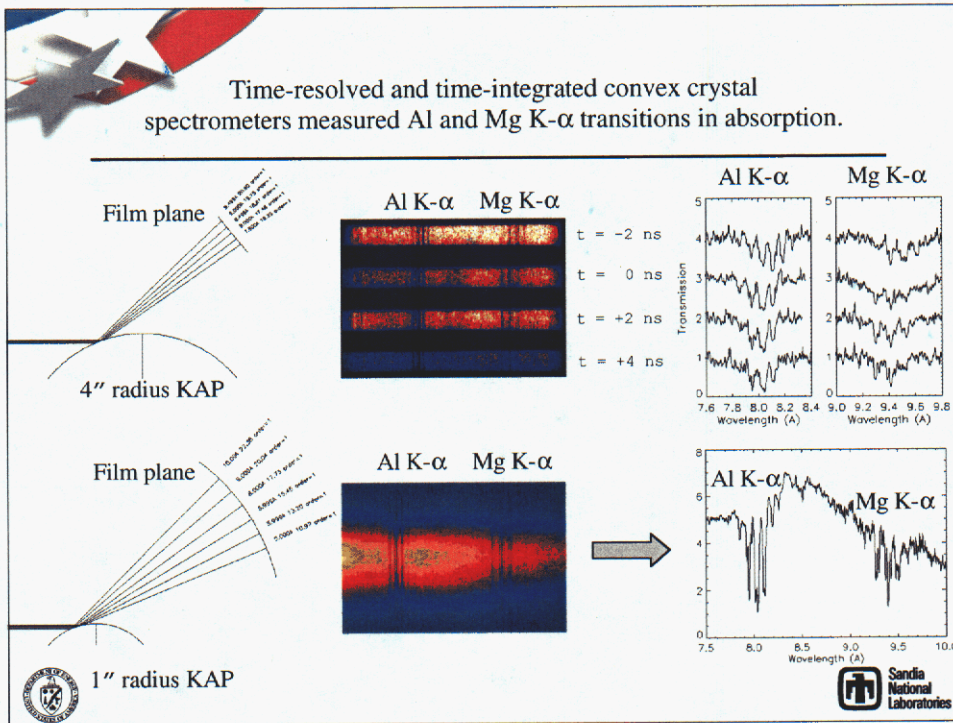


Abstract

Sandia National Laboratories' Z machine provides a good test bed for conducting basic plasma research on the interaction of x-rays with matter. In particular, recent experiments have been conducted that irradiate thin Al/MgF₂ metal foils by the radiation pulse from the side-on emission of a fast z-pinch. In these experiments, time-resolved and time-integrated spectra of K-shell absorption lines, backlit by the high-energy tail of the z-pinch radiation, are the primary diagnostic of the foil plasma conditions. The experiments are simulated by 1-D radiation-hydrodynamics calculations using a time- and frequency-dependent radiation boundary condition determined by 3-D view factor simulations of the z-pinch diode region. The calculated plasma conditions are then utilized in a detailed configuration accounting model of the atomic level populations to determine the relative amplitude of absorption features over the spectral range of interest. These calculations, and their comparison to the experimental data, will be presented and discussed.



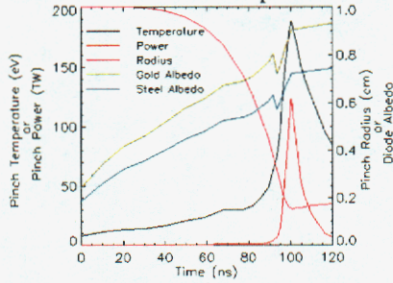




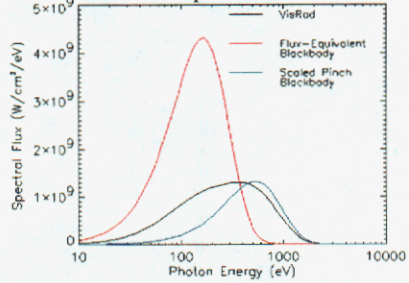


3-D simulations using the VISRAD view-factor code are utilized to calculate time-dependent, non-thermal drive spectra on the sample surface.

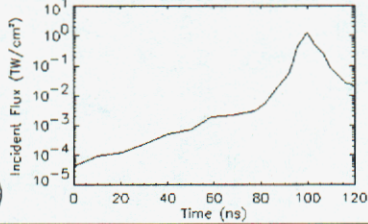
View-Factor Inputs



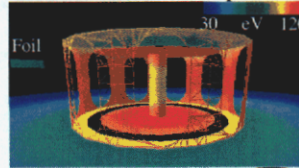
Incident Spectra at t = 100 ns



Calculated Drive Flux

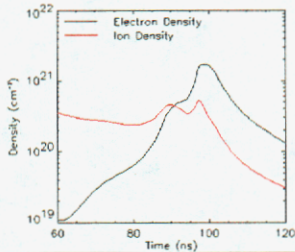
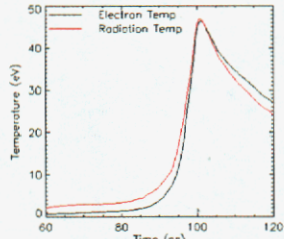


Incident Radiation Temperatures

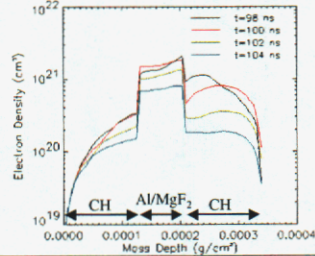
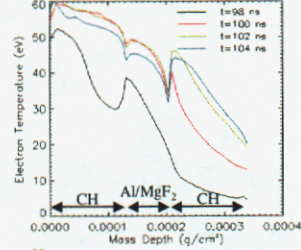


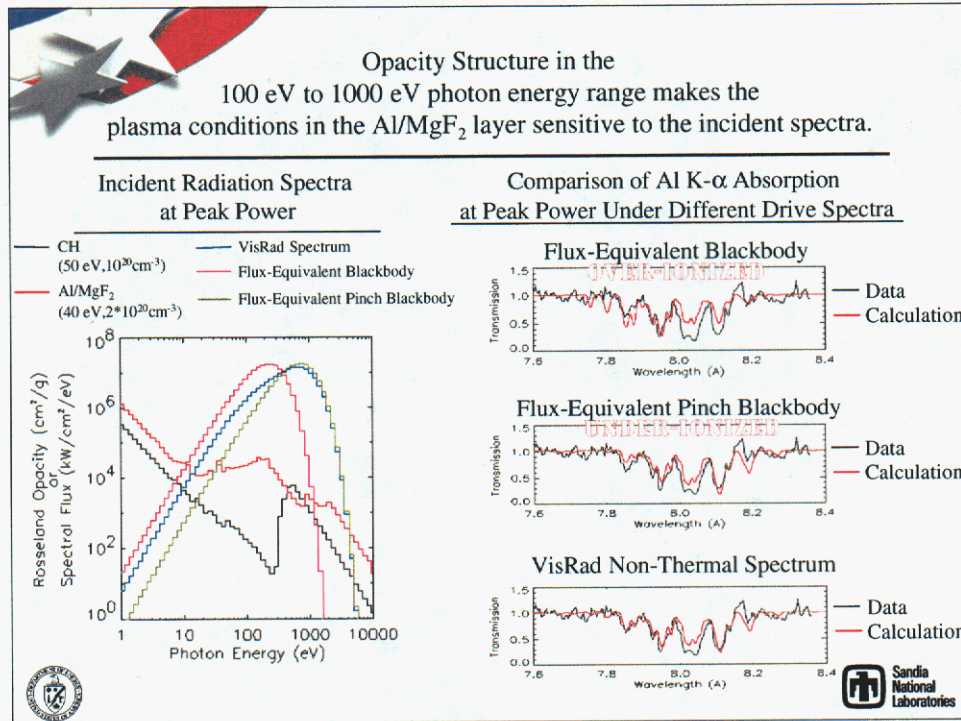
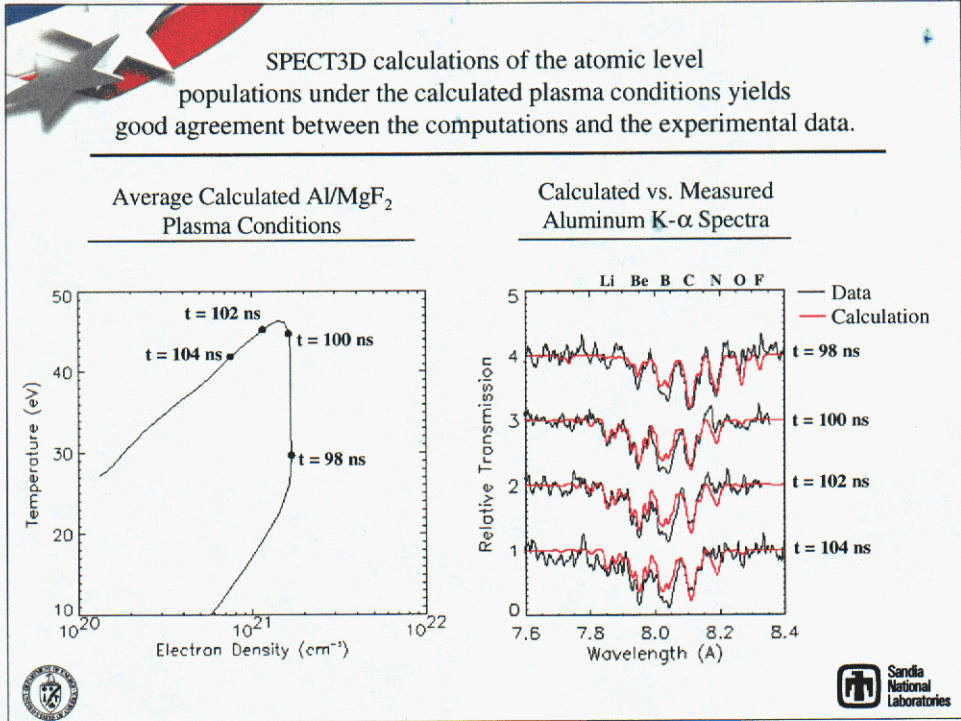
1-D simulations using the BUCKY radiation-hydrodynamics code are utilized to calculate the plasma temperature and density evolution.

Average Conditions In the Al/MgF₂ Plasma



Spatial Profiles of the Calculated Plasma Conditions



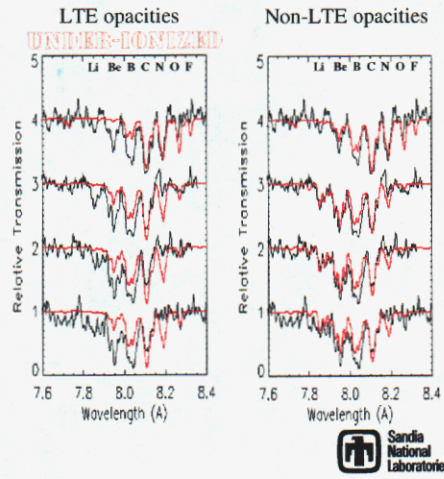
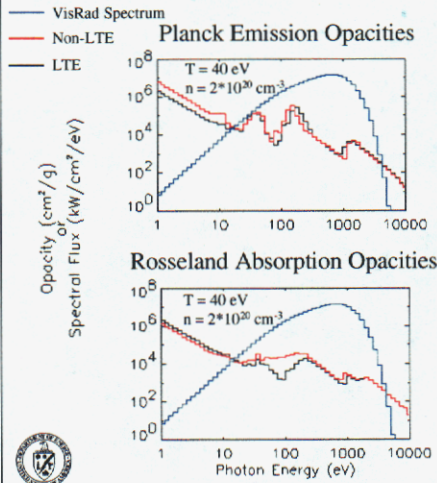




The calculated plasma conditions are also sensitive to non-LTE opacity effects requiring a detailed calculation of atomic transition rate coefficients.

Comparison between LTE and non-LTE absorption and emission opacities

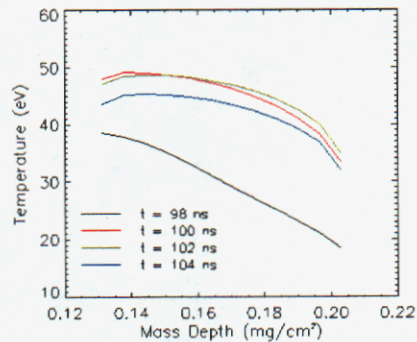
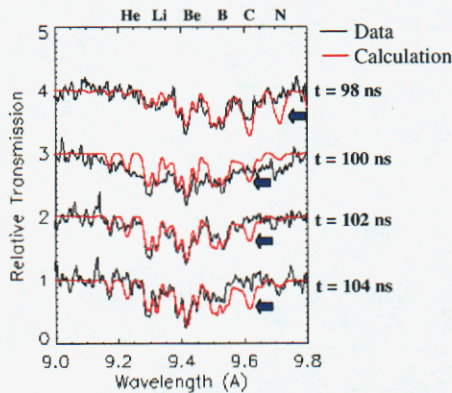
Comparison of Time-Dependent Al K- α Absorption



Calculated Mg K- α spectra are also in fair agreement with experimental data, but indicate that the calculations may contain too large of a temperature gradient.

Calculated vs. Measured Magnesium K- α Spectra

Calculated Temperature Profiles in the Al/MgF₂ Layer



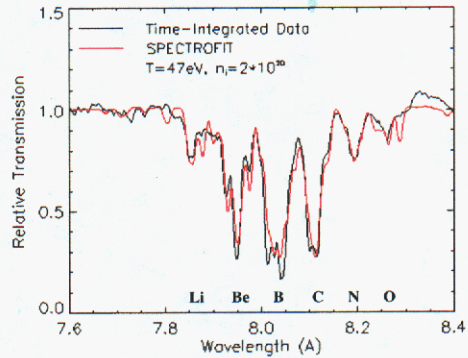


In addition to the time-dependent analysis, the time-integrated data can be analyzed by an automatic chi-squared fitting program called SPECTROFIT.

SPECTROFIT

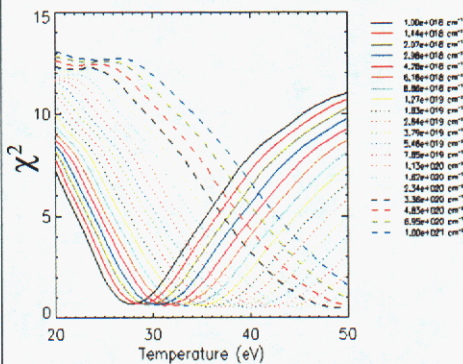
- Reads detailed line-opacity tables generated by the JATBASE code
 - LTE and Non-LTE opacities
 - Arbitrary plasma mixtures
- Compares two relative transmission absorption spectra (i.e. experiment vs. calculation) at discrete temperatures and densities.
- Computes chi-squared fit parameter for a user-defined grid resolution over a specified temperature and density range.
- Reports fit-variance, χ^2 , and $+1\sigma$

Comparison between time-integrated Al K- α data and the SPECTROFIT best fit for $n_i = 2 \times 10^{20} \text{ cm}^{-3}$.

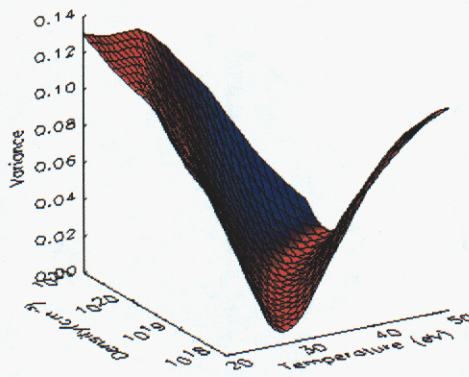


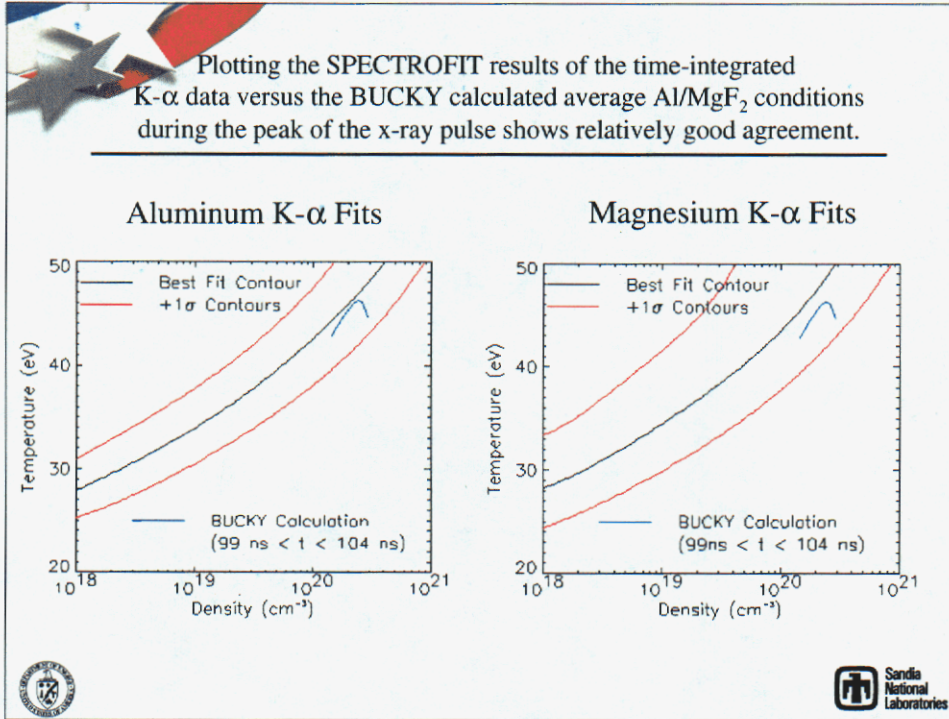
The SPECTROFIT χ^2 and fit-variance curves show the contour in temperature-density phase space where the calculated spectra best match the data.

χ^2 plots at discrete densities for the comparison between time-integrated data and the calculated Al K- α spectra



Fit variance over the searched T- ρ phase space





Summary

- Time-resolved and time-integrated data of Aluminum and Magnesium K- α absorption lines were measured in an Al/MgF₂ foil irradiated by z-pinch radiation on Sandia's Z machine.
- The data has been successfully analyzed and computationally reproduced by coupling together simulations of the radiation drive spectra, the sample radiation hydrodynamics, and detailed configuration analysis of atomic energy level transitions.
- Calculating the proper distribution of ionization states (and therefore the plasma temperature and density conditions) requires detailed consideration of the time-dependent non-thermal drive spectra and a thorough calculation of the atomic transition rate coefficients in non-LTE multi-group opacities.
- A new χ^2 spectral fitting code called SPECTROFIT has been applied to the time-integrated K- α absorption data and shows the aluminum and magnesium to be in near-equilibrium, and in good agreement with peak plasma conditions in 1-D BUCKY radiation-hydrodynamics calculations.

Sandia National Laboratories