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Comparison of Absorption, Fluorescence, and Polarization Spectroscopy of Atomic Rubidium

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Spectroscopy of Atomic Rubidium Joaquin Romero, Cayla Stifler, Seth Ashman

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The D1 transitions of rubidium are recorded using the Doppler-free saturated absorption technique (red line). This overlays the purple line which is a simple absorption spectrum. The vertical lines mark the positions of the positions of the transitions between the hyperfine levels of the upper and lower state. Transitions are labeled using the notation $5S_{1/2}(F') \rightarrow 5P_{1/2}(F)$ where just the hyperfine numbers are displayed ($F' \rightarrow F$). An "extra" peak is seen between pairs of labeled hyperfine transitions. This phenomenon is explained below.



As the laser frequency is scanned, the narrow peaks in the saturated absorption spectrum occur when the probe and pump beam interact with the same atoms. The labeled hyperfine transitions occur when both beams interact with the v=0 velocity group. The "extra" line in between each pair of labeled hyperfine transitions occurs when the laser is at a frequency halfway between two hyperfine transitions. For a particular velocity group, the pump and probe beams are each Doppler shifted by an amount equal to half the separation of the hyperfine transitions, and each induces a transition originating from the ground state. So both beams are competing for the same ground state atoms, leaving less available atoms to absorb the probe beam.

Future Work

Next we plan to perform polarization spectroscopy of our rubidium vapor. In our two-beam experimental setup, note we have included a Fresnel rhomb, capable of making our pump beam circularly polarized. This, along with the two linear polarizers in the path of the probe beam, will allow us to investigate using this additional spectroscopic technique.