IV CONCLUSION

Previously, Simon¹⁶, Saxena¹⁷ et al. and Brown¹⁸ have calculated certain elastic properties of sodium. Saxena¹⁷ et al. calculated the binding energy and compressibility both for bcc and hcp phases; however, the Born repulsive term was not included. Simon¹⁶ has done similar calculations on the elastic modulus and its first pressure derivative using an improved theory of Gombas.¹⁹ The theoretical results were compared with experimental values obtained by Diederichs and Trivisonno²⁰ and Daniels²¹. Quite a discrepancy was found for the bulk modulus and its pressure. derivative. The present calculation shows much better agreement between the theoretical values and experimental values for the isothermal bulk modulus and its first pressure derivative. There is a slight discrepancy in the second pressure derivative. However, if we consider the simplicity of the theory and the fact that there is an estimated 10% standard error in the experimental value for the second derivative the result is actually fairly good.

ACKNOWLEDGMENT

We acknowledge the support of the United States Atomic Energy Commission and use of the Central Facilities of the Cornell Materials Science Center.

-9-