

# High Pressure Studies Of Uranium Hydride – UH<sub>3</sub>

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A large variety of the lanthanide and actinide elements have the RH<sub>2</sub> and RH<sub>3</sub> compounds. Usually the RH<sub>2</sub> have the CaF<sub>2</sub> structure. The Uranium has an exemption and exhibits only the UH<sub>3</sub> compound. The UH<sub>3</sub> has two different unit cell parameters, one of which is unstable. The exposure of uranium metal to gaseous hydrogen leads to a direct reaction in which the metal is converted into the hydride UH<sub>3</sub>. This work was done to gain a better understanding of why the uranium system is different.

The uranium hydride sample was prepared by a direct reaction in which the metal is converted into the hydride UH<sub>3</sub>. The sample was kept under argon atmosphere. SEM examined the microstructure and the phases analyzed by EDS to determine the chemical compositions.

X-Ray powder diffraction measurements were taken at ambient and high pressure (up to 27GPa), using a Merrill-Bassett type diamond anvil cell (DAC). The experiments were carried out in a "Tel-Aviv"-type diamond-Anvil-Cell.

The UH<sub>3</sub>( $\beta$ ) -type structure of the UH<sub>3</sub> actinide compound is described by the cubic space group Pm3n (223). The unit cell contains eight formula units in the cell, H 24k (0.000 0.156 0.313), U1 2a (0 0 0) and U2 in 6c (0.250 0.000 0.500) (Fig 1.).

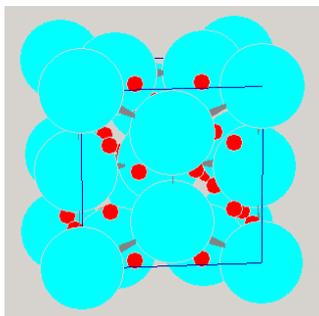


Fig 1. The crystallographic structures of UH<sub>3</sub>

**X-ray at ambient pressure:** The powder diffraction spectra of the sample taken by Cu K $\alpha$  radiation and SiO<sub>2</sub> monochromator is shown in Fig 2.

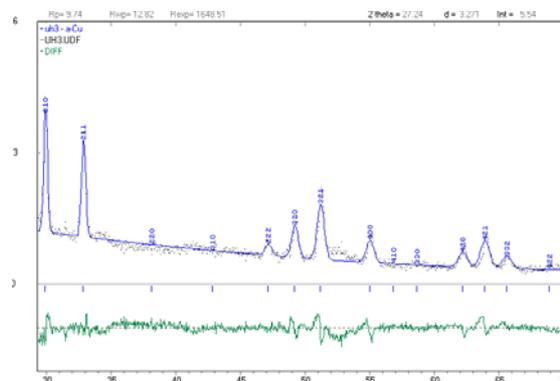


Fig 2. The powder diffraction spectra of UH<sub>3</sub> at ambient pressure.

UH<sub>3</sub> compound has a UH<sub>3</sub>( $\beta$ )-type, Pm3n symmetry cubic structure, with the lattice parameter  $a=6.6452(5)\text{\AA}$  close to the unit cell that was reported by Rundle[1], Si was used as an external standard.

**X-ray under pressure:** The UH<sub>3</sub> maintains the cubic symmetry as a function of pressure, as was determined by the Rietveld[2] analysis. The volume-pressure curve calculated from the X-ray analysis is given in Fig 3.  $V/V_0=0.87$  at 26.0GPa.

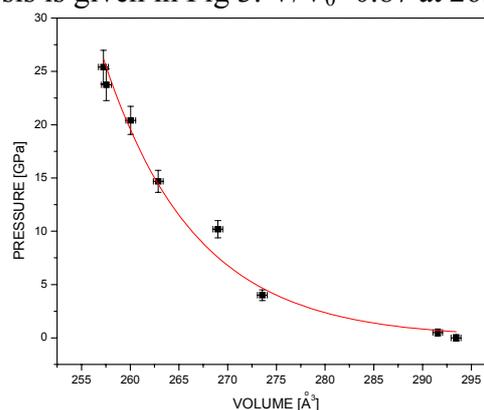


Fig. 3. Volume-Pressure curve calculated from the X-ray analysis.

## References

- [1] Rundle, R. , J. Am. Chem. Soc., 69 1719 (1947)  
Common Met., 149 (1989) 237.
- [2] R.A. Young, A. Sakthivel, T.S. Moss, C.O. Paiva-Santos, " Rietveld Analysis of X-Ray and Neutron Diffraction Patterns, Georgia Institute of Technology Atlanta, GA 30332, 19 Dec. 1994.