

# Phase Diagram Studies on the Mixed Rare Earth Oxide System ( $DY_2O_3$ & $ZRO_2$ )

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## Abstract

The ICDD is a non-profit organization established by a number of international scientific organizations. This organization provides the determination of powder patterns. Most of the patterns published in PDF obtained from useful sources of information for crystal structure. The PDF is a collection of single phase x-ray diffraction patterns in the form of tables of inter planar spacing and corresponding relative intensities. A typical data “card” obtained from the computer version of the PDF.

**Keywords:** Non-Profit Organisation, Patterns, Diffraction Systems, Inter Planar Spacing

## Introduction

Materials are used in every activity of life, living things and non-living things are made of materials. Advances in technology depend increasingly on the development of better materials of greater strength, lightness, safety, electrical conductivity, resistance to corrosion and to heat etc. .

Material Science involves investigating the relationships that exist between the structures and properties of materials [1]. Material Science is a field of science that emphasizes studies of relationships between the microstructures, synthesis, processing and the properties of materials . The materials making up the surrounding world consist of discrete particles having sub-microscopic size .Their behaviour is explained by atomic theories .

## Powder Diffraction

The ICDD is a non-profit organization established by a number of international scientific organizations .This organization provide the determination of powder patterns. Most of the patterns published in PDF obtained from useful

sources of information for crystal structure. The PDF is a collection of single phase x-ray diffraction patterns in the form of tables of inter planar spacing and corresponding relative intensities. A typical data “card” is obtained from the computer version of the PDF.

Generally, materials are broadly classified as

1. Simple materials.
2. Composite materials.

Composites are materials that occur due to combination of materials in definite proportions. Based on these properties a material falls into one of the following categories.

- i. Mechanical properties (strength, ductility etc)
- ii. Thermal properties (specific heat, melting point etc)
- iii. Magnetic properties (permeability, hysteresis etc)
- iv. Chemical properties (chemical composition, acidity etc) Optical properties (reflectivity, refractive index etc)
- v. Physical properties.

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According to these properties a material is classified as,

A pure metal consist entirely of atoms of only one element. It has its own unique physical properties such as melting point, boiling point, and thermal or heat conductivity. Example: Fork

An alloy is the intimate combination of two or more pure metals that have been dissolved while molten, one into the other. An alloy could be harder, softer, stronger, less easily corroded and so forth, than the individual metals composing it.

They are chemical compounds of the type metals and non-metals. They are generally good thermal and electrical insulators. They are hard and brittle. Example: Scissors, a glass vase, floor tile, a china tea-cup [1].

They are organic compounds, generally in the form of long chain. They have low density. Example: Spoon, fork, knife [1]

They have electrical properties that are intermediate between conductors (eg: metals and alloys) and insulators (Example: Ceramics and polymers). They are composed of more than one type of materials. Most common and familiar composites is fiberglass, in which small glass are embedded within a polymeric material [4].

### Powder Data Analysis

Subsequently, the high purity samples like  $BaF_2$ ,  $NaCl$ ,  $LaB_6$ ,  $URh_3$ ,  $UGa_3$ ,  $Dy_2O_3$ ,  $CsCl$  and  $Tm_2O_3$ ,  $ZrO_2$ ,  $KCl$ ,  $KBr$  and  $Se$  theoretical pattern was generated by using the PCW software [5]. Input parameters like the crystal structure, atom positions were taken from Pearson's Intermetallic Database [6-10]. The computed theta values along with respective hkl values and intensities are used to index the experimentally observed Bragg peaks. Care was taken to compare the intensities of every peak and assign the right hkl value. A graph comparing the experimentally observed theta, intensity and theoretical intensity, calculated difference in both values ( $2\theta_{EXP} - 2\theta_{CAL}$ ) was prepared for this purpose. The indexed experimental data was fed into NBS\*AIDS83 SOFTWARE, in order to get the lattice parameter, cell, volume, figure of merit etc [11]. The output giving the lattice parameter found using the least square fitting method (NBS\*AIDS83), volume and the figure of merit.

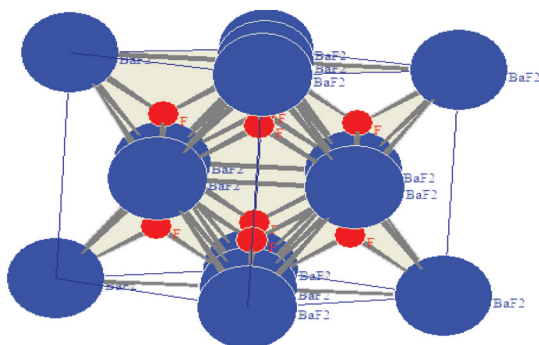


Figure (1) Crystal Structure of  $BaF_2$

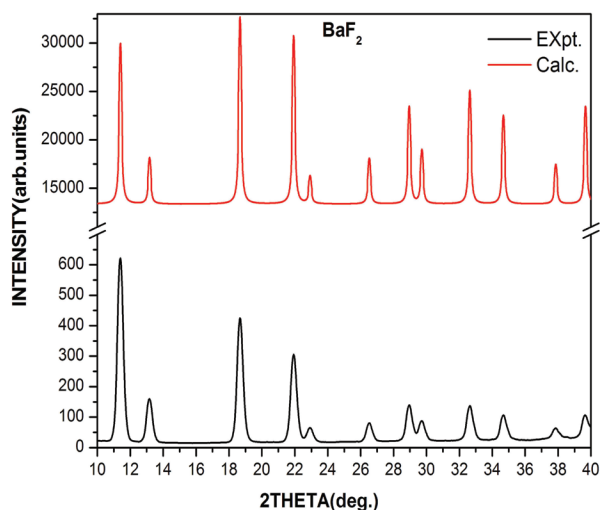


Figure (2) XRD Pattern of  $BaF_2$

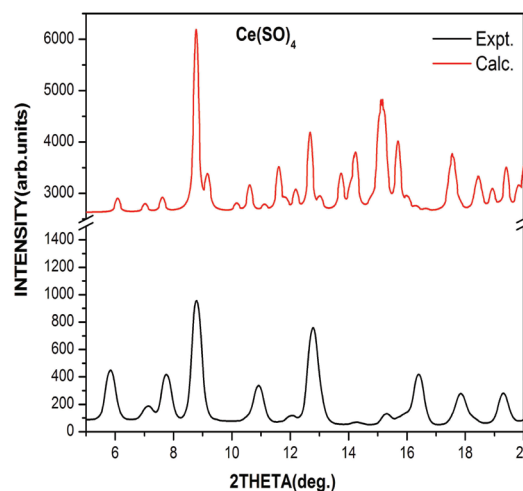


Figure (3.) XRD Pattern Of  $Ce(SO_4)_4$

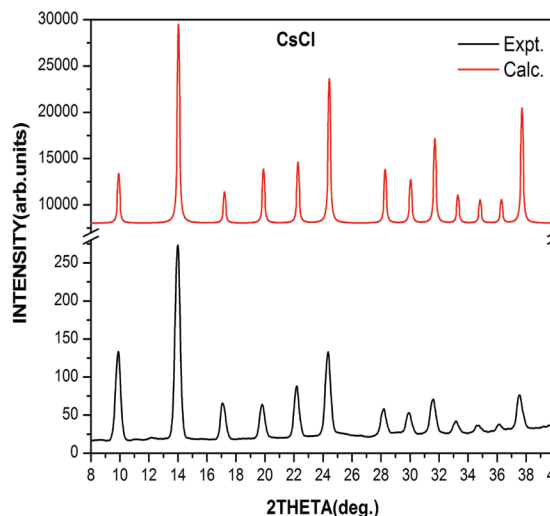


Figure (4) XRD Pattern Of  $CsCl$ .

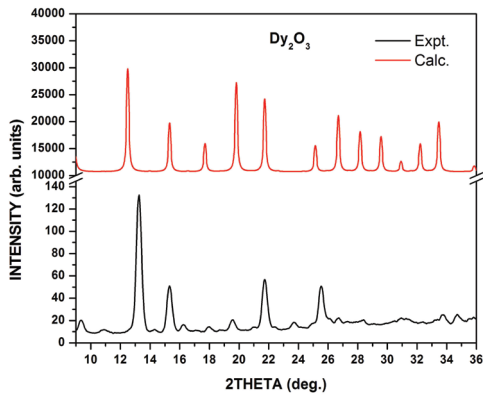
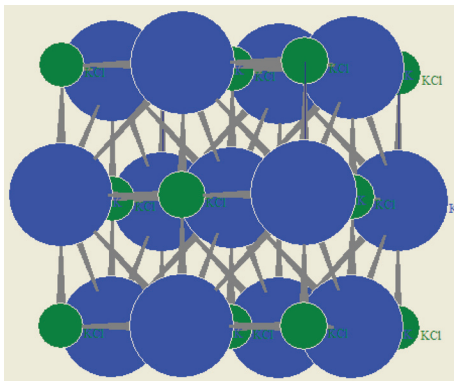


Figure (5) XRD Pattern Of  $Dy_2O_3$



Figure(6) Crystal Structure of KCl

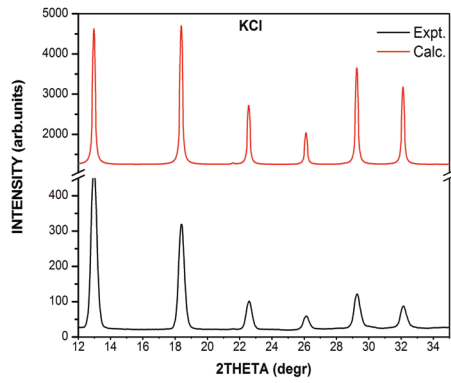


Figure (7) XRD Pattern of KCl

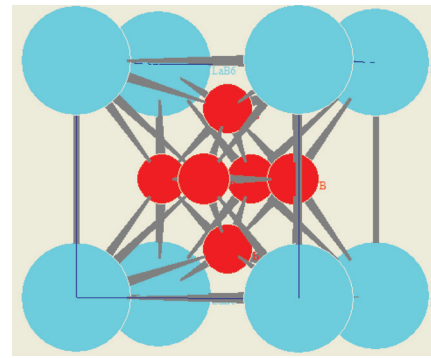


Figure (8) Crystal Sturcture of  $LaB_6$

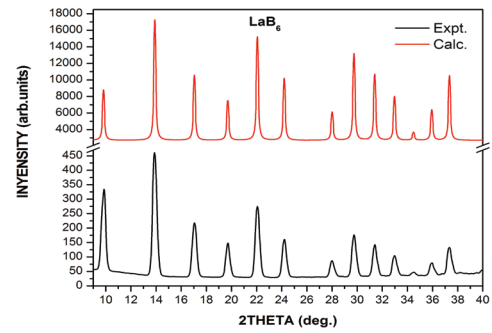


Figure (9) XRD Pattern of  $LaB_6$

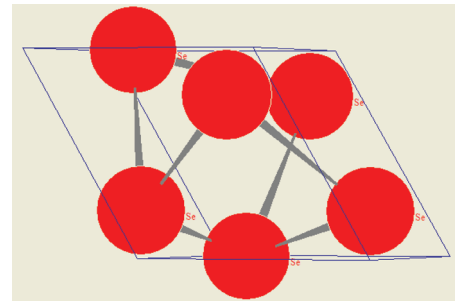


Figure (10) Crystal Sturcture of Se

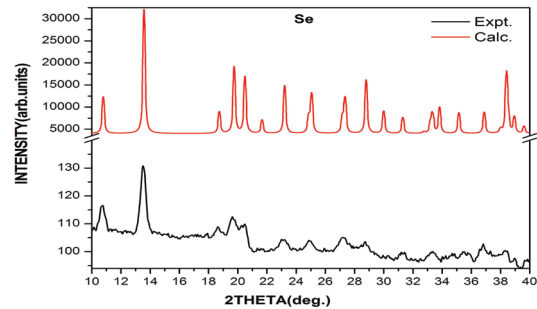


Figure (11) XRD Pattern of Se

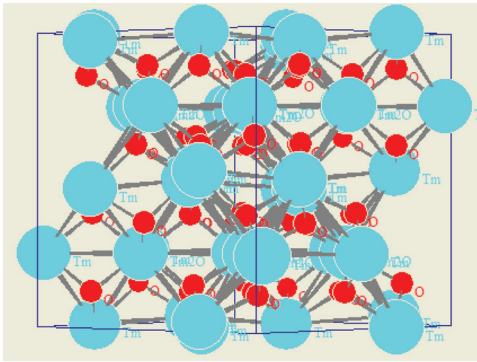


Figure (12) Crystal Structure of Tm<sub>2</sub>O<sub>3</sub>

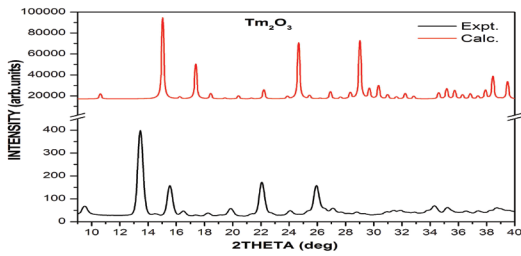


Figure (13) XRD Pattern of Tm<sub>2</sub>O<sub>3</sub>

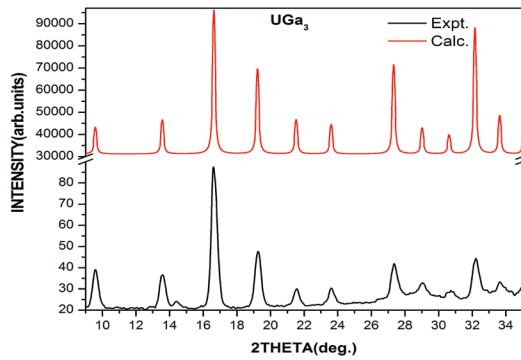


Figure (14) XRD Pattern of UGa<sub>3</sub>

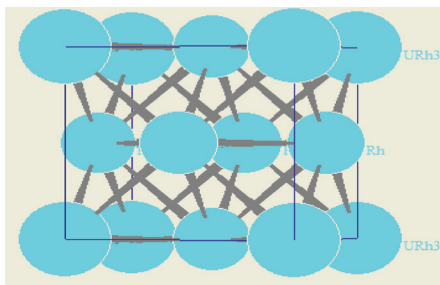


Figure (15) Crystal Structure of URh<sub>3</sub>

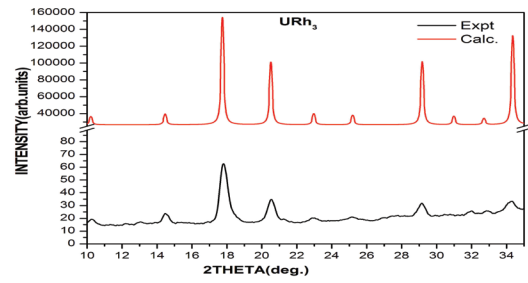


Figure (16) XRD Pattern of URh<sub>3</sub>

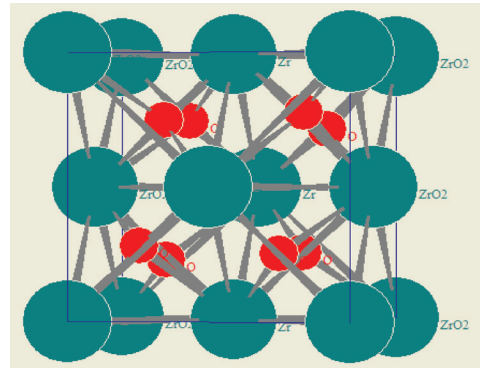


Figure (17) Crystal Structure Of ZrO<sub>2</sub>

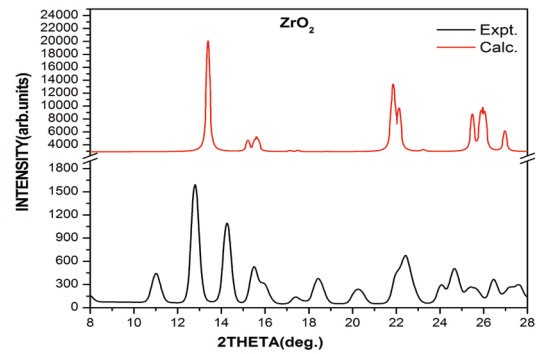


Figure (18) XRD Pattern of ZrO<sub>2</sub>

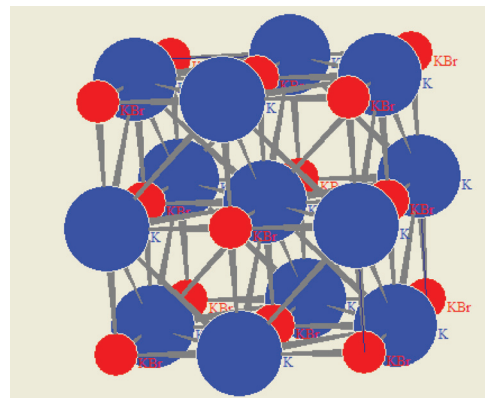


Figure (19) Crystal Structure of KBr

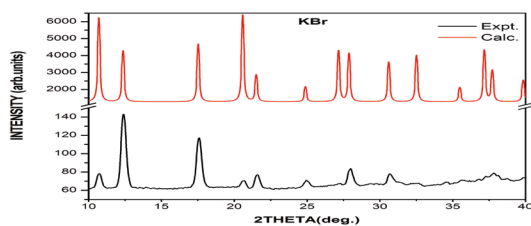


Figure (20) XRD Pattern of KBr

## Conclusion

The XRD pattern of mixed oxides of  $\text{Dy}_2\text{O}_3$  &  $\text{ZrO}_2$  were prepared in the pestle and mortar with few drops of hexane. An annealed in the double zone furnace for 12 hours at  $1000^\circ\text{C}$ . The compositions vary from 0.9% to 52.9%. From the composition of 0.9% to 10.9% and 42.7% to 50.3% the 100% peak of  $\text{Dy}_2\text{O}_3$  is visible due to its high intensity. And its peak is very sharp. But from 0.9% to 45.4% the 100% peak of  $\text{ZrO}_2$  is not visible, and it is hidden in the shoulder of 100% peak of  $\text{Dy}_2\text{O}_3$ . This is due to low intensity in  $\text{ZrO}_2$ . In these compositions 21.6%, 32.2% and 52.9%, the 100% peak of  $\text{ZrO}_2$  is visible independently due to its high intensity. In these compositions the 100% peak of  $\text{Dy}_2\text{O}_3$  is not visible, but it is hidden in the shoulder of  $\text{ZrO}_2$ . This is because of low intensity compared to  $\text{ZrO}_2$ .

In summary, a high temperature phase diagram studies on the mixed rare earth oxide system ( $\text{Dy}_2\text{O}_3$  &  $\text{ZrO}_2$ ) was performed. This study reveals their C-type Cubic crystal structure and the B-type Monoclinic at higher temperature. This helps to identify the phase stability. Rare earth oxides  $\text{Dy}_2\text{O}_3$  &  $\text{ZrO}_2$  corrected data were analyzed using NBS\*AIDS83, FIT2D&PCW software.

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