

UNCLASSIFIED

102924

Armed Services Technical Information Agency

Reproduced by

DOCUMENT SERVICE CENTER

KNOTT BUILDING, DAYTON, 2, OHIO

This document is the property of the United States Government. It is furnished for the duration of the contract and shall be returned when no longer required, or upon recall by ASTIA to the following address: Armed Services Technical Information Agency, Document Service Center, Knott Building, Dayton 2, Ohio.

NOTICE: WHEN GOVERNMENT OR OTHER DRAWINGS, SPECIFICATIONS OR OTHER DATA ARE USED FOR ANY PURPOSE OTHER THAN IN CONNECTION WITH A DEFINITELY RELATED GOVERNMENT PROCUREMENT OPERATION, THE U. S. GOVERNMENT THEREBY INCURS NO RESPONSIBILITY, NOR ANY OBLIGATION WHATSOEVER; AND THE FACT THAT THE GOVERNMENT MAY HAVE FORMULATED, FURNISHED, OR IN ANY WAY SUPPLIED THE SAID DRAWINGS, SPECIFICATIONS, OR OTHER DATA IS NOT TO BE REGARDED BY IMPLICATION OR OTHERWISE AS IN ANY MANNER LICENSING THE HOLDER OR ANY OTHER PERSON OR CORPORATION, OR CONVEYING ANY RIGHTS OR PERMISSION TO MANUFACTURE, USE OR SELL ANY PATENTED INVENTION THAT MAY IN ANY WAY BE RELATED THERETO.

UNCLASSIFIED

BEST

AVAILABLE

COPY

AD No. 102934

ASTIA CUE COPY

FC
BAC

Studies of the Rare-Earth Hydrides
Technical Report IX

LANTHANUM MONOXIDE

Office of Naval Research
Physical Sciences Division

Project No. NR 356-290

Contract No. Nonr 228(03)

By

James C. Warf, Project Supervisor
and
William L. Korst, Research Assistant

The Department of Chemistry
University of Southern California
Los Angeles 7, California

June 30, 1956

Lanthanum Monoxide

It was reported in Technical Report VII (p.96) that a sample of lanthanum metal filings which had been annealed at 475° for 52 hours gave a powder diffraction pattern corresponding to two f.c.c. lattices, with cell constants of $5.307 \pm 0.005\text{\AA}$. and $5.249 \pm 0.005\text{\AA}$. The data are given in Table 20 at the end of this report. The first number corresponds to the lattice constant of f.c.c. lanthanum, which was reported as $5.296 \pm 0.005\text{\AA}$. by Ziegler et al. (1953).

An attempt was made to identify the second f.c.c. phase, having a_0 equal to 5.249 Å. It seemed most likely that this would be either an oxide or a nitride phase. No hitherto reported lanthanum oxide has a f.c.c. structure,¹ but Ellinger & Zachariassen (1953) have reported a f.c.c. samarium monoxide, SmO, with the NaCl-type structure, and a slightly variable lattice constant. The Sm-Sm distance in this monoxide is approximately 0.04 Å. smaller than in the metal. The authors have identified this phase with the grey coating formed on metal pieces on heat treatment.

Lanthanum nitride, LaN, has been described by Young & Miegler (1952) as having the NaCl-type structure, with a_0 equal to 5.295 ± 0.004 Å. (5.284 ± 0.004 kX.)² on the basis of calculations of intensity ratios for adjacent lines for this and other structures considered as possible. A comparison of measured intensity ratios from the two sets of lines found in this investigation with the intensity ratios calculated by Young & Miegler for f.c.c. lanthanum and for LaN indicates that the lattice to which the cell constant of 5.249 Å. corresponds does have the NaCl-type structure.

¹The strong lines of the b.c.c. modification of La_2O_3 form an apparent f.c.c. lattice, but with a much larger a_0 , namely 5.7 Å. This is discussed further below.

²Landelli & Eotti (1937) had previously reported 5.286 Å. (5.275 kX.).

This comparison is made in Appendix III, Table 21. The principal evidence lies in the relative weakness of the lines with $\sum h_i^2$ equal to 11 and 19. Since this lattice constant is so different from that of the nitride, it is concluded that it corresponds to that of a monoxide, LaO , similar to SmO . It may be noted that the La-La distance in the presumed monoxide is also about 0.04 Å. less than in the metal.

Ziegler, Young & Lloyd (1953), in an investigation of the crystal structure of lanthanum metal, reported that the principal f.c.c. pattern, ascribed to f.c.c. (β -) lanthanum, was usually accompanied by some lines from h.c.p. (α -) lanthanum, as well as by a number of weak lines which were assignable to two different f.c.c. lattices, one having a lattice constant 0.5-1% less than that of f.c.c. lanthanum, and the other having a_0 between 5.63 and 5.66 Å. The former of these, called the "γ" structure by the authors, and for which no explanation was suggested, corresponds to the lanthanum monoxide lattice described above.

The second extra f.c.c. lattice appearing was ascribed by the authors to lanthanum hydride. However, it seems unlikely that there should have been any hydride present in the metal sample. If there had been some originally present, the hydrogen should have been effectively

removed by the 13 hours' heating at 700° C. under high vacuum to which one sample was subjected, but for which the "hydride" structure persisted. It is suggested here that the lines of the "hydride" pattern are actually the stronger lines of the b.c.c. form of La_2O_3 (often called the C-modification) (Löfberg, 1955; Bommer, 1939), which appear to belong to a f.c.c. lattice half as large, because the metal atoms lie in a slightly distorted face-centered cubic array (Strukturbericht II, 38-40). If this were so, the b.c.c. lattice to which the lines of the "hydride" structure would correspond would have a cell constant of twice 5.66 Å., or 11.32 Å., to make it as large as possible. The actual b.c.c. parameter has not been very accurately determined, being reported as 11.4 kX. units by Löfberg on the basis of only four blurred measured lines, with considerable deviation. This appears to be a reasonable explanation, if not a conclusive one, of what Ziegler, Young & Floyd considered to be an appearance of the hydride.

TABLE 20.--X-ray diffraction data for lanthanum metal sample, annealed. Cu radiation, Ni filter. See p. 96.

<u>sin θ obs.</u>	<u>f.c.c. La</u>			<u>LaO</u>		
	<u>$\sum h_1^2$</u>	<u>a_0</u>	<u>I_{obs}</u>	<u>$\sum h_1^2$</u>	<u>a_0</u>	<u>I_{obs}</u>
0.2547	3	5.243		3	5.243	
.2932	4	5.258				
.2955				4	5.217	
.4129	8	5.281	40	8	5.221	40
.4176						
.4836	11	5.286	70	11	5.226	50
.4892						
.5050	12	5.289	20	12	5.234	20
.5102						
.5822	16	5.296	10	16	5.236	10
.5890						
.6345	19	5.296	40	19	5.239	35
.6414						
.6712	20	5.294	40	20	5.241	40
.6578						
.7120	24	5.304	30	24	5.243	35
.7203						
.7551	27	5.305*	40	27	5.248*	40
.7634						
.8222	32	5.304*	10	32	5.247*	15
.8311						
.8589	35 α_1	5.306*	70	35 α_1	5.248*	
.8684						
.8705	36 α_1	5.309*	40	36 α_1	5.248*	50
.8807						
.9178	40 α_1	5.308*		40 α_1	5.252*	60
.9276						
.9520	43 α_1	5.305*		43 α_1	5.249*	
.9623	44 α_1	5.310*		44 α_1	5.253*	
.9727						

Average a_0 values computed using starred values are 5.307 $\pm$ 0.005 Å. (f.c.c. La) and 5.249 $\pm$ 0.005 Å. (LaO).

See Technical Report I for references