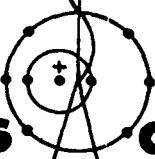


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Hugoniot Equations of State
of Li⁶H, Li⁶D, LiⁿH, and LiⁿD (U)


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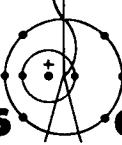
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Hugoniot Equations of State of Li⁶H, Li⁶D, LiⁿH, and LiⁿD (U)

by

S. P. Marsh

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HUGONIOT EQUATIONS OF STATE OF Li⁶H, Li⁶D, Li⁷H, AND Li⁷D

by

S. P. Marsh

Energy per gram

ABSTRACT

An experimental value of $1.08 \text{ cm}^3/\text{g}$ for $(\partial E/\partial P)_v$ was determined from the Hugoniot data on pressed Li⁶D of various porosities which agrees well with the thermodynamic value of $(\partial E/\partial P)_v$. For each of the isotopic combinations Li⁶H, Li⁷H, Li⁶D, and Li⁷D crystal-density Hugoniots and equations of state were calculated from Hugoniot data on pressed material having porosity of 5% or less and from the thermodynamic values of $(\partial E/\partial P)_v$. No evidence of a polymorphic transition was observed.

Ultrasonic measurements of longitudinal and shear-wave velocities in these materials pressed to near crystal density were obtained, and the isotropic elastic moduli were calculated. In addition, elastic constants were determined for crystals of Li⁷H, Li⁷D, and Li⁷D.

INTRODUCTION

Hugoniot data for Li⁷H and Li⁷D have been reported by a number of investigators. In 1954, Walsh¹ studied pressed Li⁷H to pressures of 160 kbar. He also reported data² on single crystals of Li⁷(D_{0.80}H_{0.20}) shocked to pressures above 500 kbar. In 1954, Burton and Landeen³ obtained two Hugoniot data points on pressed Li⁷H at pressures above 700 kbar using a spherically convergent driver system. In 1965, Kusubov⁴ obtained data on pressed Li⁶H, Li⁶D, Li⁷H, and Li⁷D at pressures up to 500 kbar, as well as longitudinal and shear-wave velocities for each material. In 1969, May⁵ and Guess⁶ reported Hugoniot data below 200 kbar, obtained using quartz-gage detectors on pressed Li⁷H. Guess also reported longitudinal and shear-wave velocities in this same material.

In this endeavor, we have determined the Hugoniots of Li⁶D pressed to near crystal density and at porosities of approximately 5, 8, 17, 28, 36, and 44%. The wide range of initial densities enabled us to determine the average values of the Grüneisen gamma, $\gamma = V(\partial P/\partial E)_v$. These values are compared with the Grüneisen ratio determined from thermodynamic measurements at standard temperature and pressure. The Hugoniots of the isotopic combinations, Li⁶H, Li⁷H and Li⁷D, and Li⁷D for pressed samples with an initial 5% porosity have also been determined. Further, we have obtained Hugoniot data on a few crystals of Li⁷H and Li⁷D, and have determined elastic constants for crystals of Li⁷H, Li⁷H, and Li⁷D.

MATERIAL

The crystals used in these experiments were prepared by Pretzel and Thrasher of LASL

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Group CMB-3. All pressed materials and their chemical analysis and densities were obtained from Y-12 at Oak Ridge. The Li⁶ had a composition of 95.5 at. % Li⁶ and 4.5 at. % Li⁷; the normal Liⁿ was 7.5 at. % Li⁶ and 92.5 at. % Li⁷; and the Li⁷ was 100 at. % Li⁷. The isotopic purity of the hydrogen and deuterium was above 99 at. %. The densities of the four pure isotopic combinations were calculated from the lattice parameters reported by Anderson et al.⁷ We used these densities to compute the value of the crystal densities of the particular isotopic compositions used here. The Oak Ridge analysis provided an estimate of the oxygen in our samples, and we calculated their ideal or crystal density by mixing 2 wt % of the hydroxide or deutoxide using the densities that Vier⁸ reported and the density for oxygen-free materials. These densities were used to determine the porosity and are listed in Table I at the end of this report.

ELASTIC CONSTANTS IN LITHIUM HYDRIDE AND LITHIUM DEUTERIDE

Elastic-wave velocities were determined ultrasonically using a pulse-echo method⁹ on pressed isotropic samples at near crystal density for all isotopic compositions. The longitudinal and shear-wave velocities (V_L and V_s) were measured, and the bulk sound velocity (C_b) was calculated from the expression

$$C_b^2 = V_L^2 - 4/3 V_s^2.$$

Table II lists the results of sound-velocity measurements on the various isotopic combinations of lithium hydride at near crystal density and at approximately 5% porosity. The bulk modulus (B_b), shear modulus (μ), Young's modulus (Y), longitudinal modulus (B_L), and Poisson's ratio (σ) are also listed.

Using the same apparatus used for making pulse-echo sound-velocity measurements in pressed specimens, we determined the sound velocities in LiⁿH, Li⁷H, and Li⁷D crystals. Because of the difficulty in seeing the arrival time

of the echo signals, we used a different method of measurement in which we observed the arrival times through several different thicknesses of sample and obtained the velocity from the derivative of the resulting x-t data. Table III shows the results of these measurements for both the longitudinal and shear-wave velocities in the [100] direction and the longitudinal and two shear-wave velocities in the [110] direction. The errors in the longitudinal and shear-wave velocities are 0.5 and 1.0%, respectively. We obtained the elastic constants using the following set of equations which involve only the velocities in the [100] direction and the bulk sound velocities (C_b) deduced from sound-velocity measurements on pressed isotropic samples.

$$c_{11} = \rho V_L^2 [100],$$

$$c_{44} = \rho V_s^2 [100],$$

$$c_{12} = \frac{3\rho C_b^2 - \rho V_L^2 [100]}{2}.$$

We compared the measured velocities in the [110] direction with the calculated velocities in that direction to check the derived elastic constants. The expressions used to calculate these velocities were

$$V_L[110] = \sqrt{\frac{c_{11} + c_{12} + 2c_{44}}{2\rho}},$$

$$V_{1s}[110] = \sqrt{\frac{c_{44}}{\rho}}$$

(for particle motion in the [001] direction),

$$V_{2s}[110] = \sqrt{\frac{c_{11} - c_{12}}{2\rho}}$$

(for particle motion in the [110] direction).

The measured and calculated velocities in this direction differ by several percent, a greater amount than would be anticipated from the error in the elastic constants, which is 1, 2, and 20% in c_{11} , c_{44} , and c_{12} , respectively. The reason for this discrepancy is not understood. Another disturbing feature observed in this table is that the c_{11} and c_{44} values for Li⁷D are lower than the corresponding values for LiⁿH and Li⁷H. We

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feel that because the lattice spacing is smaller for Li⁷D, the zero-point energy is also lower and, consequently, the elastic constants should be higher. We can give no reason for this apparent reversal. The differences in the elastic constants for Li⁶H and Li⁷H are not statistically significant.

The elastic moduli in Table II show a distinct relationship to isotopic composition. The bulk modulus, shear modulus, Young's modulus, and longitudinal modulus agree for Li⁶H and Li⁷H and are 3% smaller than the elastic moduli of the deuterides, which are also in agreement. In this case, the elastic moduli are consistent with the zero-point energy assumption mentioned above.

SHOCK-WAVE EXPERIMENTS

We used Walsh and Christian's shock-impedance match technique¹⁰ with 2024 Al as a standard, and the optical flash-gap method to determine the Hugoniots. The shock velocities in both the standard and lithium hydride samples were measured with a sweeping-image camera. Having the equation of state of the standard and the shock velocities in both the standard and lithium hydride samples permits determination of the pressure and particle velocity. A linear shock-particle velocity Hugoniot,

$U_s = 5.328 + 1.338 U_p$ (in units of km/sec), and a constant value of

$$\rho\gamma = 5.57 \text{ g/cm}^3$$

(where ρ and γ are the density and Grüneisen parameter) specify the equation of state of the standard required for the impedance-match solution.

We determined the density and internal energy of the shocked materials using the Rankine-Hugoniot relations. Tables IV - XIII show the experimental data and derived Hugoniot parameters.

ANALYSIS AND INTERPRETATION

Performing shock-wave experiments on materials of different porosities allows examination of different pressure-energy regions because

the greater the initial volume, the greater the pressure and internal energy at the same final density. This can be seen in Fig. 1 where the pressure-density shock-wave data for the Li⁶D are plotted. The curve in this figure is a linear fit of the U_s - U_p data for the high-density samples, presented in Fig. 2.

An average $(\partial E / \partial P)_v$ value can be determined for each experimental point of the low-initial-density material by differencing its pressure and energy with the pressure and energy calculated at the same density from the fit of the high-density data. These $(\partial E / \partial P)_v$ values for the five most porous materials are shown in Fig. 3. The $(\partial E / \partial P)_v$ values calculated from porous-material data having small energy and pressure offsets from the high-density Hugoniot (which includes all the low-pressure data) have large errors; consequently, we show only $(\partial E / \partial P)_v$ values for which ΔP was larger than 15 kbar. The average $(\partial E / \partial P)_v$ value determined from these experiments is $1.08 \text{ cm}^3/\text{g}$, and within experimental error it does not appear to be a function of density or pressure.

We calculated the Grüneisen γ 's at standard conditions, along with the corresponding $(\partial E / \partial P)_v$ values, using the expressions

$$\gamma = \frac{\beta C_v^2}{c_p}, \quad \gamma = V_0 / (\partial E / \partial P)_v.$$

We obtained the values of β (volume coefficient of thermal expansion at 25°C) for all isotopic compositions by interpolating the thermal-expansion data of Anderson et al.⁷ (Fig. 4) and used them for our isotopic compositions. The resulting Li⁷H value ($31.0 \times 10^{-6}/^\circ\text{K}$) is somewhat lower than that obtained by a number of other investigators,¹¹⁻¹⁴ but it agrees with the compilation by Goldsmith et al.¹⁵ for Li⁶H at 25°C. The 6.682-cal/mol°K value of c_p (specific heat at a constant pressure) for Li⁶H was obtained from the JANAF Tables.¹⁶ This value was used for all isotopic compositions. The specific volume (V_0)

and bulk sound velocities are those given in Table I (for chemically pure compounds) and Table II. A summary of these values and the resulting values of γ and $(\partial E/\partial P)_v$ is given in Table XIV. The experimental ($1.08 \text{ cm}^3/\text{g}$) value of $(\partial E/\partial P)_v$ for Li^6D agrees well with the thermodynamic value of $1.13 \text{ cm}^3/\text{g}$ shown in this table.

The values of $(\partial E/\partial P)_v$ in Table XIV were used to transform the Hugoniot data obtained on the porous samples to pressures and energies corresponding to samples at crystal density.⁹ These transformations are made at constant volume. The recentered Hugoniot data and a linear least-squares fit of the U_s-U_p points are shown in the U_s-U_p plane and the $P-\rho$ plane in Figs. 5-12 along with the bulk sound velocities with which the intercepts should agree if there are no transitions. The least-squares fits summarized in Table XV were determined only from points having $U_p > 0.9 \text{ km/sec}$. We chose this lower particle-velocity limit to avoid the extreme sensitivity of the transformed U_s-U_p points to small errors in the original data at low pressure. The bulk sound velocities and the Hugoniot intercepts agree well.

From these Hugoniots and the thermodynamic data in Table XIV, one can determine a complete equation of state for each material.⁹ Tables XVI-XIX list the thermodynamic parameters on the Hugoniot, the foot of the release isentrope, and the isotherms at room temperature and at 0°K . The zero-pressure parameters used in calculating these loci are summarized at the head of each table.

In Figs. 13-16 our Hugoniot results are compared with the results of other investigators studying similar isotopic compositions. We converted isothermal compression data to Hugoniot shock velocity-particle velocity points using the equations of state shown in Tables XVI-XIX. Stevens and Lilley's results¹⁷ generally show less compressibility than ours, and Kusubay's results⁴

for Li^6H and Li^6D show considerably greater compressibility. Burton and Landeen's high-pressure data³ for Li^6H are also more compressible than our extrapolated Hugoniot.

TRANSITION IN LITHIUM HYDRIDE AND LITHIUM DEUTERIDE

Several investigators have postulated a transition from the NaCl to the CsCl structure in lithium hydride, but no transformation has been observed. Schumacher,¹⁸ using two different Born-Mayer potentials for lithium hydride, calculates a transition at 3 to 4 kbar involving a volume change of about 0.4%. Voronov et al.¹⁹ showed experimentally that there is a linear relationship of compressibility to pressure up to 20 kbar which did not fit the simple Born model. Using Schumacher's potential function, he predicted that the transition should occur at 140 kbar with a volume change of 14%. Other experimenters^{20,21} have also examined lithium hydride and have found no evidence for a transformation. Stephens and Lilley¹⁷ determined the isothermal compressibility of Li^6H , Li^6D , Li^7H , and Li^7D up to 40 kbar and found that their data fit the simple Born-Mayer model, but observed no transition.

There is no evidence of a transition for any of the isotopic combinations in the pressure region (60 to 450 kbar) that we investigated. The data for the low-porosity Li^6D are particularly important, and although the low-pressure data have considerable scatter there is no evidence of a kink indicative of a transition in the U_s-U_p curve. It is somewhat more difficult to detect transitions in the more porous samples, however. Although we see no evidence for a transition in our data, it may possibly be present but have too small a volume change to be observed as a discontinuity in the shock-wave data.

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TABLE I

CRYSTAL DENSITIES OF Li⁶H, Li⁶D, Li⁷H, AND Li⁷D

Pure Isotopes		Isotopic Purity for this Study		Corrected for Assumed 2 wt% Hydroxide or Deuterioxide	
Material	Density (g/cm ³)	Material	Density (g/cm ³)	Density (g/cm ³)	
Li ⁶ H	0.6842	(Li ⁶ _{0.955} Li ⁷ _{0.045})H	0.689	0.696	
Li ⁶ D	.7905	(Li ⁶ _{0.955} Li ⁷ _{0.045})D	.795	.802	
Li ⁷ H	.7830	(Li ⁶ _{0.075} Li ⁷ _{0.925})H	.776	.783	
Li ⁷ D	.8900	(Li ⁶ _{0.075} Li ⁷ _{0.925})D	.883	.890	

TABLE II

SOUND VELOCITIES AND ELASTIC MODULI OF Li⁶H, Li⁶D, Li⁷H, AND Li⁷D

Material	Density (g/cm ³)	Sound Velocities (km/sec)			Elastic Moduli (kbar)				
		V _L	V _S	C _b	B _s	μ	Y	B _L	σ
Li ⁶ H	0.698	10.67	7.18	6.72	314	359	781	794	.086
Li ⁶ D	.799	10.10	6.80	6.35	322	369	802	815	.085
Li ⁷ H	.782	10.05	6.75	6.34	314	356	776	789	.089
Li ⁷ D	.894	9.56	6.43	6.02	324	369	803	817	.087
Li ⁶ H	.666	10.42	6.86						
Li ⁶ D	.764	9.72	6.53						
Li ⁷ H	.743	9.84	6.61						
Li ⁷ D	.840	9.36	6.31						

TABLE III
ELASTIC CONSTANTS OF Li⁷H, Li⁷H, AND Li⁷D

Mat	ρ_0 (g/cm ³)	V _L [100] (km/sec)	V _S [100] (km/sec)	C _b (km/sec)	c ₁₁ (kbar)	c ₄₄ (kbar)	c ₁₂ (kbar)	V _L [110] (km/sec)	V _{1s} [110] V _{2s} [110]	
									ξ along [001]	ξ along [110]
Li ⁷ H	0.776	9.06	7.40	6.34	636	424	150	10.55	7.62	5.77
								10.26*	7.40*	5.60*
Li ⁷ H	.783	9.02	7.39	6.34	637	428	154	10.50	7.70	5.80
								10.25*	7.39*	5.56*
Li ⁷ D	.890	8.38	6.81	6.02	626	419	171	-	-	-
								9.58*	6.81*	5.05*

*Calculated

TABLE VIII

HUGONIOT DATA FOR Li⁶D ($\rho_0 = 0.579 \text{ g/cm}^3$)

ρ_0 (g/cm ³)	U_s (km/sec)	U_r (km/sec)	P (Mbar)	ρ (g/cm ³)	U_{state} (km/sec)	ρ_0 (g/cm ³)	U_s (km/sec)	U_r (km/sec)	P (Mbar)	ρ (g/cm ³)	U_{state} (km/sec)
0.592	2.87	0.85	0.014	0.840	5.98	0.456	2.60	1.18	0.014	0.825	6.17
.571	3.39	1.12	.022	.851	6.17	.456	3.45	1.84	.026	.889	6.53
.572	4.40	1.56	.039	.885	6.53	.450	4.34	2.12	.041	.879	6.91
.587	5.34	2.00	.061	.908	6.91	.457	5.80	2.83	.078	.924	7.58
.571	6.81	2.77	.108	.962	7.59	.482	6.25	3.19	.092	.942	7.81
.574	7.29	3.01	.128	.977	7.81	.452	6.58	3.35	.100	.920	7.95
.572	7.49	3.17	.136	.991	7.95	.448	7.05	3.73	.118	.953	8.28
.570	7.83	3.54	.158	1.039	8.26	.425	7.48	3.93	.125	.897	8.41
.571	8.29	3.69	.175	1.029	8.41	.458	7.83	4.09	.147	.959	8.59
.575	8.59	3.88	.191	1.048	8.59	.447	8.37	4.48	.167	.958	8.89
.587	9.09	4.19	.224	1.089	8.89	.441	8.42	4.48	.167	.944	8.91
.583	9.23	4.21	.227	1.072	8.91	.448	8.75	4.73	.185	.974	9.12
.579	9.35	4.47	.242	1.109	9.12	.451	9.09	4.75	.194	.943	9.18
.570	9.54	4.52	.257	1.083	9.16	.444	9.39	4.92	.205	.933	9.31
.578	9.91	4.65	.266	1.089	9.31	.446	9.71	5.09	.220	.937	9.46
.596	10.23	4.79	.292	1.121	9.47	.444	9.61	5.11	.218	.948	9.47
.588	9.93	4.83	.282	1.145	9.48	.450	10.04	5.36	.242	.965	9.70
.594	10.48	5.06	.315	1.148	9.70	.434	10.20	5.49	.243	.940	9.78
.594	10.62	5.14	.325	1.152	9.78	.442	10.02	5.49	.243	.977	9.78
.578	10.52	5.18	.315	1.138	9.78						

TABLE X

HUGONIOT DATA FOR Li⁶D ($\rho_0 = 0.448 \text{ g/cm}^3$)

TABLE IX

HUGONIOT DATA FOR Li⁶D ($\rho_0 = 0.514 \text{ g/cm}^3$)

ρ_0 (g/cm ³)	U_s (km/sec)	U_r (km/sec)	P (Mbar)	ρ (g/cm ³)	U_{state} (km/sec)	ρ_0 (g/cm ³)	U_s (km/sec)	U_r (km/sec)	P (Mbar)	ρ (g/cm ³)	U_{state} (km/sec)
0.522	3.01	1.14	0.018	0.840	8.17	0.669	6.42	0.72	0.031	0.753	5.94
.526	4.08	1.59	.034	.861	6.53	.661	6.38	.73	.031	.747	5.95
.530	4.88	2.05	.053	.913	6.91	.664	6.31	.73	.031	.751	5.95
.532	6.35	2.83	.096	.961	7.59	.666	6.86	1.00	.045	.779	8.18
.514	6.76	3.10	.108	.950	7.81	.671	6.81	1.02	.046	.789	8.20
.527	7.12	3.24	.122	.968	7.95	.881	7.49	1.42	.070	.816	6.57
.519	7.54	3.62	.142	.998	8.26	.661	8.13	1.79	.096	.847	8.90
.521	7.98	3.78	.157	.988	8.41	.662	9.02	2.43	.145	.907	7.49
.539	8.35	3.94	.177	1.021	8.59	.660	8.91	3.19	.208	.974	8.19
.506	9.04	4.34	.198	.972	8.91	.666	10.71	3.74	.268	1.023	8.71
.499	8.70	4.35	.189	.999	8.89	.671	10.97	3.99	.294	1.054	8.95
.511	9.08	4.60	.213	1.035	9.12	.664	11.71	4.63	.380	1.098	9.54
.502	9.38	4.64	.218	.993	9.16	.664	12.31	4.95	.404	1.110	9.86
.508	9.72	4.78	.235	.996	9.31	.664	12.41	4.99	.411	1.111	9.90
.508	10.05	4.96	.253	1.003	9.47	.664	12.73	5.40	.457	1.154	10.28
.499	9.86	4.98	.245	1.009	9.46	.664	12.93	5.48	.471	1.153	10.36
.510	10.27	5.23	.274	1.038	9.70						
.497	10.51	5.33	.279	1.010	9.78						
.500	10.33	5.35	.276	1.035	9.78						

TABLE XI

HUGONIOT DATA FOR Li⁶H ($\rho_0 = 0.666 \text{ g/cm}^3$)

TABLE XII
HUGONIOT DATA FOR LiⁿH AND Li⁷H
(ρ₀ = 0.739 g/cm³)

ρ ₀ (g/cm ³)	U _s (km/sec)	U _s (km/sec)	P (Mbar)	ρ (g/cm ³)	U _{s+4} (km/sec)	ρ ₀ (g/cm ³)	U _s (km/sec)	U _s (km/sec)	P (Mbar)	ρ (g/cm ³)	U _{s+4} (km/sec)
0.743	6.11	0.71	0.032	0.841	5.91	0.840	5.88	0.70	0.034	0.953	5.94
.740	5.97	.73	.032	.842	5.95	.839	5.79	.71	.035	.956	5.95
.742	6.52	.98	.048	.874	6.18	.840	6.30	.96	.051	.992	6.18
.735	6.44	1.01	.048	.872	6.20	.839	6.33	.98	.052	.993	6.20
.736*	7.15	1.37	.072	.911	6.54	.840	7.50	1.72	.108	1.090	6.90
.729*	7.17	1.39	.072	.904	6.55	.839	8.29	2.35	.163	1.170	7.49
.741*	7.12	1.40	.074	.923	6.56	.844	9.11	3.07	.236	1.274	8.19
.741	7.15	1.40	.074	.922	8.57	.838	9.86	3.61	.298	1.321	8.71
.742*	6.90	1.41	.072	.933	6.56	.840	10.21	3.84	.330	1.347	8.95
.736*	6.92	1.41	.072	.825	6.56	.842	10.97	4.48	.414	1.423	9.57
.739*	7.57	1.73	.097	.957	6.86	.841	11.03	4.59	.426	1.441	9.68
.732*	7.63	1.73	.097	.946	6.86	.840	11.24	4.78	.451	1.461	9.86
.741	7.73	1.76	.101	.959	6.90	.841	11.56	4.80	.466	1.437	9.90
.741*	7.83	1.80	.104	.962	6.93	.840	11.85	5.20	.517	1.496	10.28
.749*	7.84	1.80	.106	.972	6.94	.840	11.85	5.29	.527	1.518	10.36
.742	8.66	2.39	.154	1.025	7.49	crystal data (ρ ₀ = 0.890 g/cm ³)					
.733*	8.67	2.42	.154	1.017	7.51	0.890*	7.13	0.94	0.060	1.025	6.20
.739*	8.62	2.45	.158	1.032	7.54	.890*	10.01	3.54	.315	1.376	8.71
.737*	8.83	2.64	.171	1.050	7.71	.890*	10.96	4.42	.431	1.491	9.57
.735*	8.76	2.64	.170	1.052	7.71	*Li ⁷ D					
.742	9.46	3.14	.221	1.111	8.19	*Li ⁷ D					
.740*	9.58	3.17	.225	1.105	8.22	*Li ⁷ D					
.737*	9.58	3.18	.225	1.104	8.23	*Li ⁷ D					
.741	10.25	3.68	.280	1.156	8.71	*Li ⁷ D					
.740*	10.27	3.72	.283	1.160	8.75	*Li ⁷ D					
.737*	10.24	3.73	.282	1.160	8.75	*Li ⁷ D					
.740*	10.22	3.74	.283	1.167	8.75	*Li ⁷ D					
.735*	10.19	3.75	.280	1.162	8.75	*Li ⁷ D					
.735*	10.34	3.77	.287	1.157	8.78	*Li ⁷ D					
.743	10.56	3.93	.307	1.183	8.95	*Li ⁷ D					
.740*	11.22	4.52	.375	1.238	9.50	*Li ⁷ D					
.741	11.30	4.55	.381	1.240	9.54	*Li ⁷ D					
.741	11.45	4.57	.388	1.233	9.57	*Li ⁷ D					
.734*	11.63	4.86	.414	1.260	9.82	*Li ⁷ D					
.741	11.91	4.86	.429	1.251	9.86	*Li ⁷ D					
.741	12.07	4.90	.438	1.247	9.90	*Li ⁷ D					
.738*	11.80	4.90	.427	1.262	9.88	*Li ⁷ D					
.741	12.16	5.32	.480	1.318	10.28	*Li ⁷ D					
.741	12.15	5.42	.488	1.338	10.36	*Li ⁷ D					
crystal data (ρ ₀ = 0.777 g/cm ³)											
0.783*	7.59	0.96	0.057	0.896	6.20	*Li ⁷ D					
.776*	7.67	1.20	.071	.920	6.42	*Li ⁷ D					
.763*	7.50	1.30	.075	.924	6.50	*Li ⁷ D					
.777*	9.13	2.58	.183	1.083	7.71	*Li ⁷ D					
.783*	10.54	3.61	.298	1.190	8.71	*Li ⁷ D					
.783*	11.71	4.49	.411	1.269	9.57	*Li ⁷ D					
.778*	12.04	4.82	.450	1.293	9.88	*Li ⁷ D					

*Li⁷H

TABLE XIII
HUGONIOT DATA FOR LiⁿD AND Li⁷D
(ρ₀ = 0.840 g/cm³)

TABLE XIV

THERMODYNAMIC GRÜNEISEN γ AND $(\partial E / \partial P)_V$ FOR Li⁶H, Li⁶D, Li⁸H, AND Li⁸D

<u>Mat</u>	<u>ρ_0 (g/cm³)</u>	<u>V_0 (cm³/g)</u>	<u>β (/°K)</u>	<u>C_0 (cm/sec)</u>	<u>c_p (cal/mol °K)</u>	<u>c_p (erg/g °K)</u>	<u>γ</u>	<u>$(\partial E / \partial P)_V$ (cm³/g)</u>
Li ⁶ H	0.689	1.451	84.9×10^{-6}	6.72×10^6	6.682	3.953×10^7	0.970	1.496
Li ⁶ D	.795	1.258	95.1	6.35	6.682	3.461	1.108	1.135
Li ⁸ H	.776	1.289	93.0	6.34	6.682	3.518	1.063	1.213
Li ⁸ D	.883	1.133	107.7	6.02	6.682	3.122	1.250	0.906

TABLE XV

SUMMARY OF CRYSTAL-DENSITY HUGONIOT
COEFFICIENTS OF THE EQUATION $U_s = C_0 + S U_p$
FOR Li⁶H, Li⁶D, Li⁸H, AND Li⁸D FOR DATA
HAVING $U_p > 0.9$ km/sec

<u>Material</u>	<u>ρ_0 (g/cm³)</u>	<u>C_0 (km/sec)</u>	<u>S</u>
Li ⁶ H	0.696	6.740	1.189
Li ⁶ D	.802	6.464	1.130
Li ⁸ H	.783	6.426	1.167
Li ⁸ D	.890	6.138	1.151

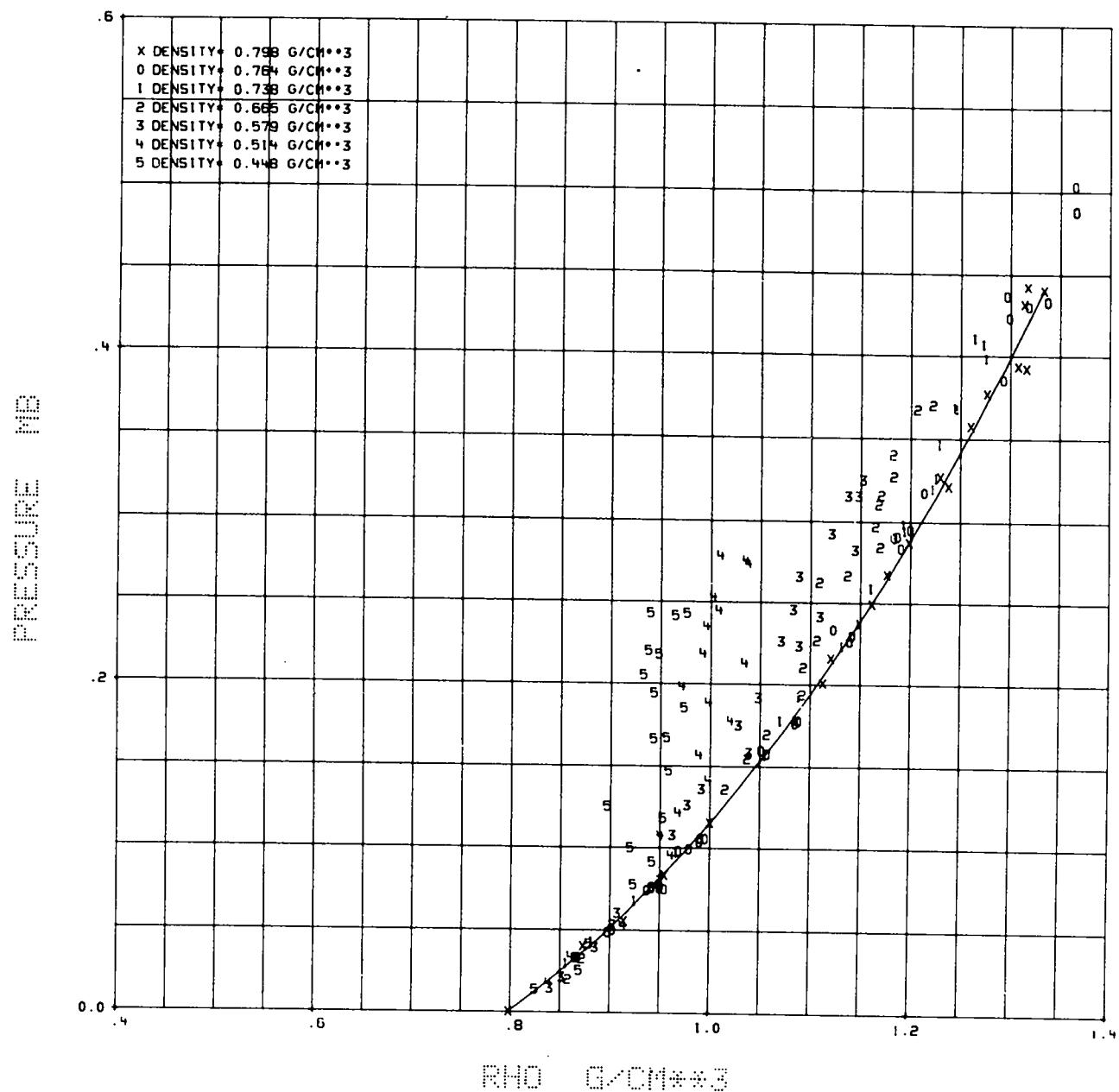


Fig. 1. Experimental pressure-density Hugoniot data for Li^6D at seven different porosities. The curve is the fit shown in Fig. 2.

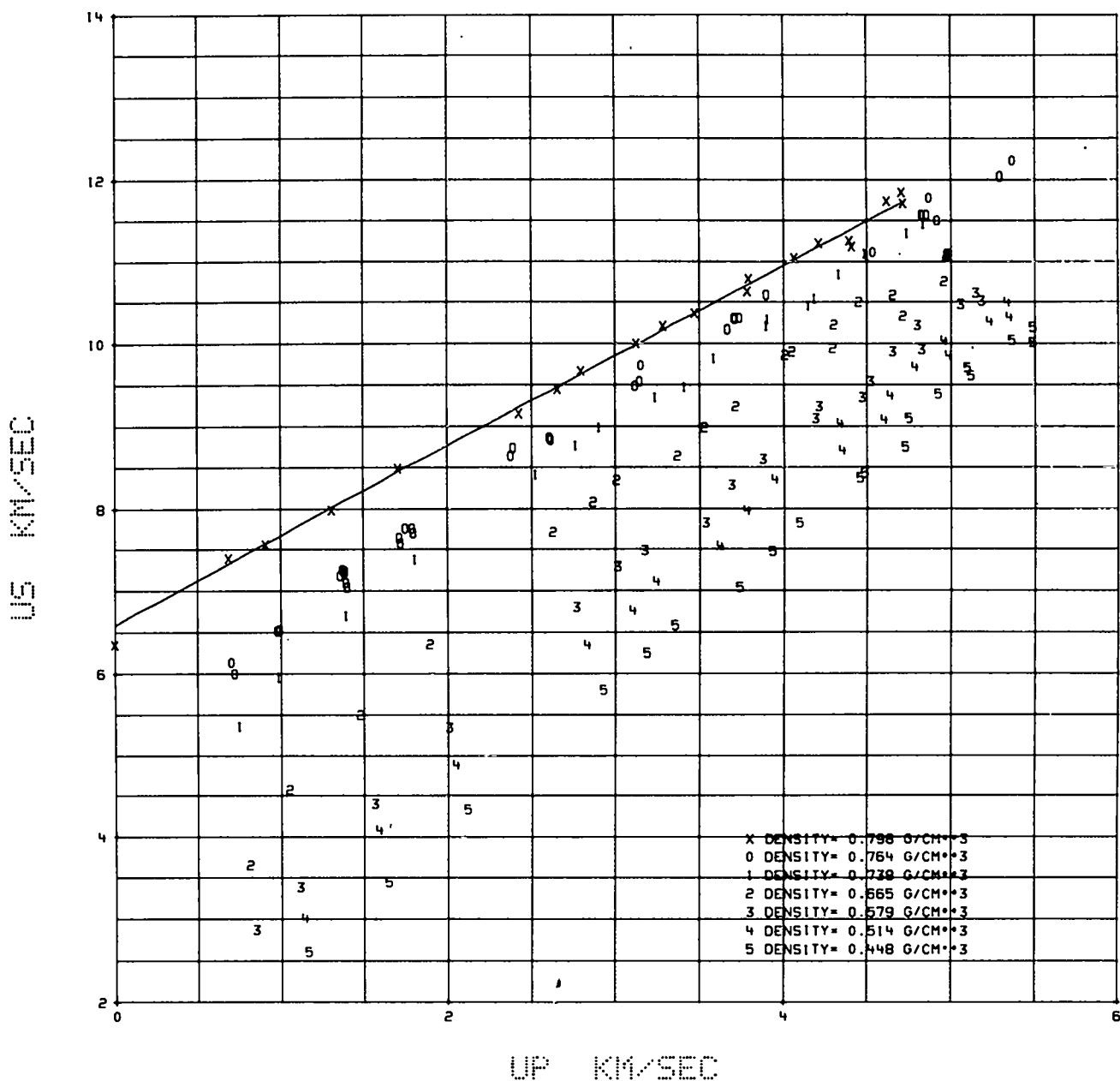


Fig. 2. Experimental shock velocity - particle velocity Hugoniot for Li^6D at seven different porosities. The linear least-squares fit of the data (excluding the bulk sound velocity) of the lowest porosity material is shown. The average density at each porosity is indicated in the legend. The Hugoniots of the porous samples curve downward in a manner predictable by calculations.

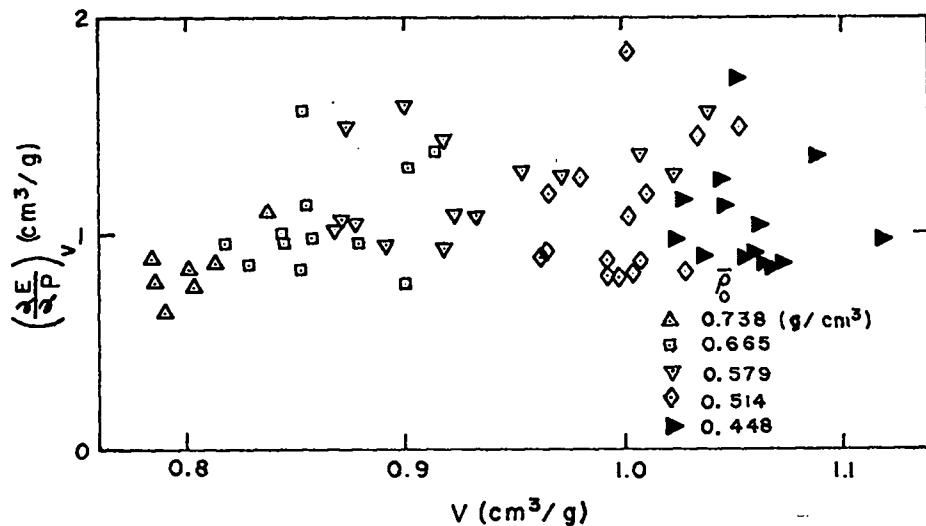


Fig. 3. $(\partial E / \partial P)_V$ vs specific volume for Li^6D for all data having $\Delta P > 15$ kbar.

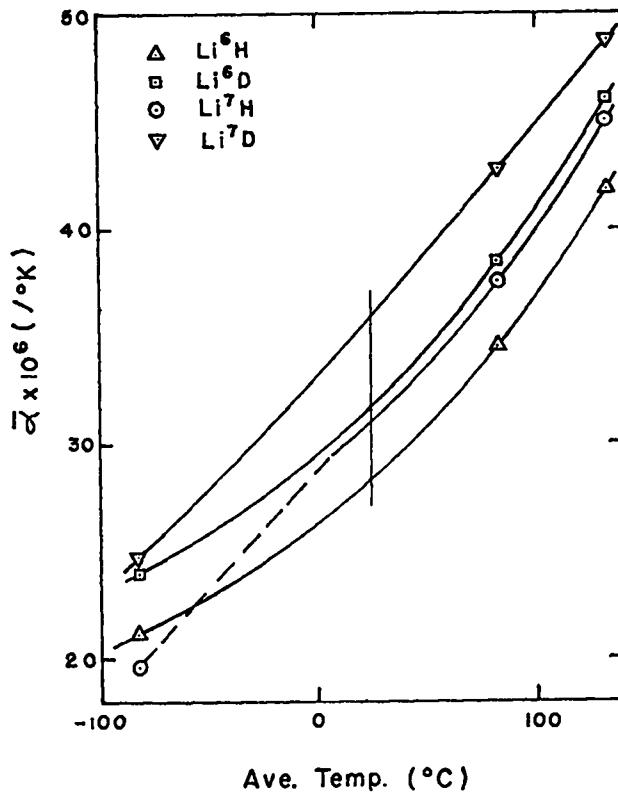


Fig. 4. Average coefficients of linear thermal expansion for Li^6H , Li^6D , Li^7H , and Li^7D over the interval T to 25°C using the data of Anderson et al.⁷ The dashed line connects the unexplained low value of Li^7H at low temperature. The $\bar{\alpha}$ values at 25°C are identical to the instantaneous α values at that temperature.

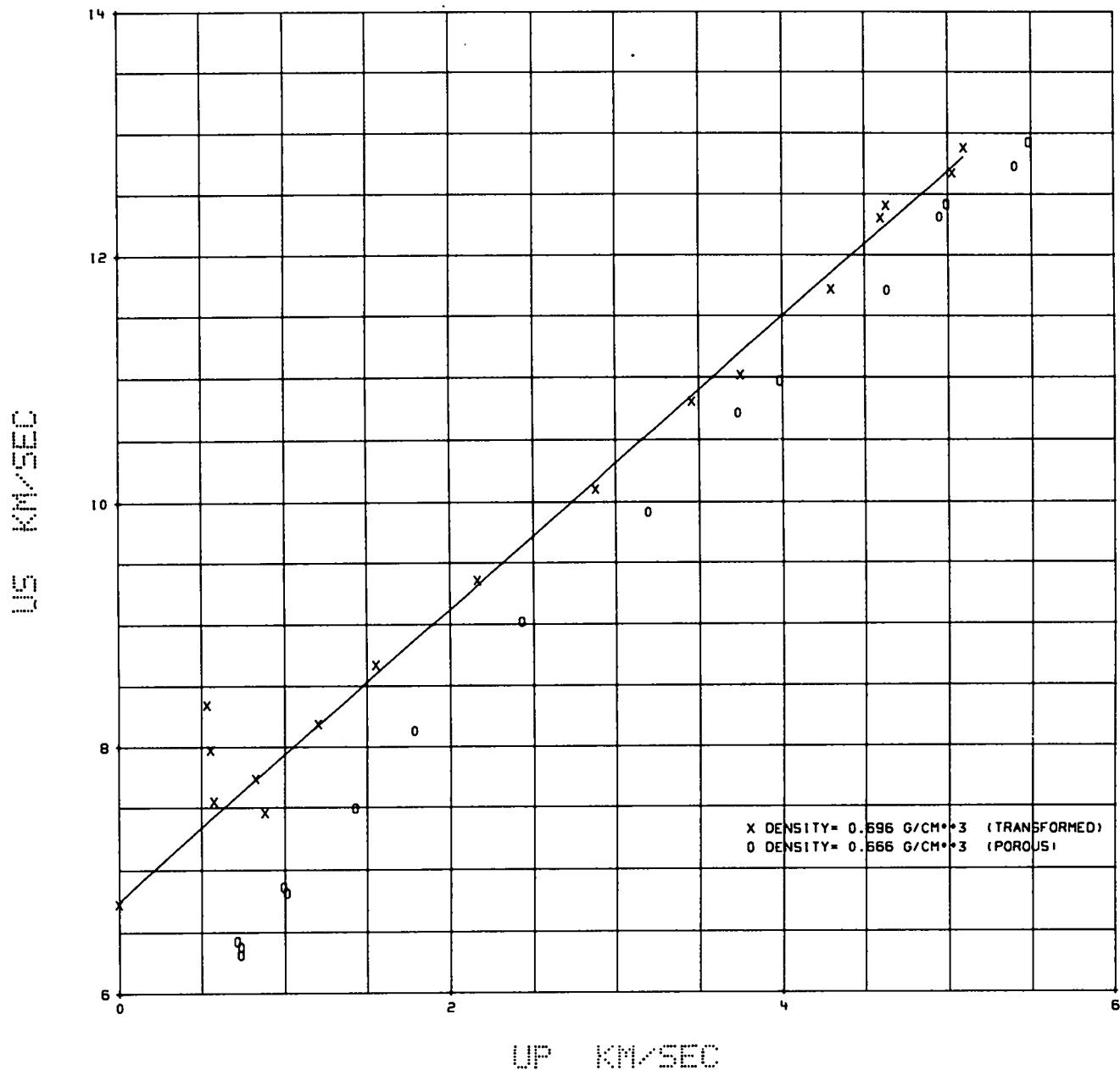


Fig. 5. Shock velocity - particle velocity Hugoniot data for Li⁶H. Experimental porous Hugoniot data and transformed crystal-density Hugoniot points are shown. A linear least-squares fit of the transformed crystal-density points is shown for points having $U_p > 0.9$ km/sec.

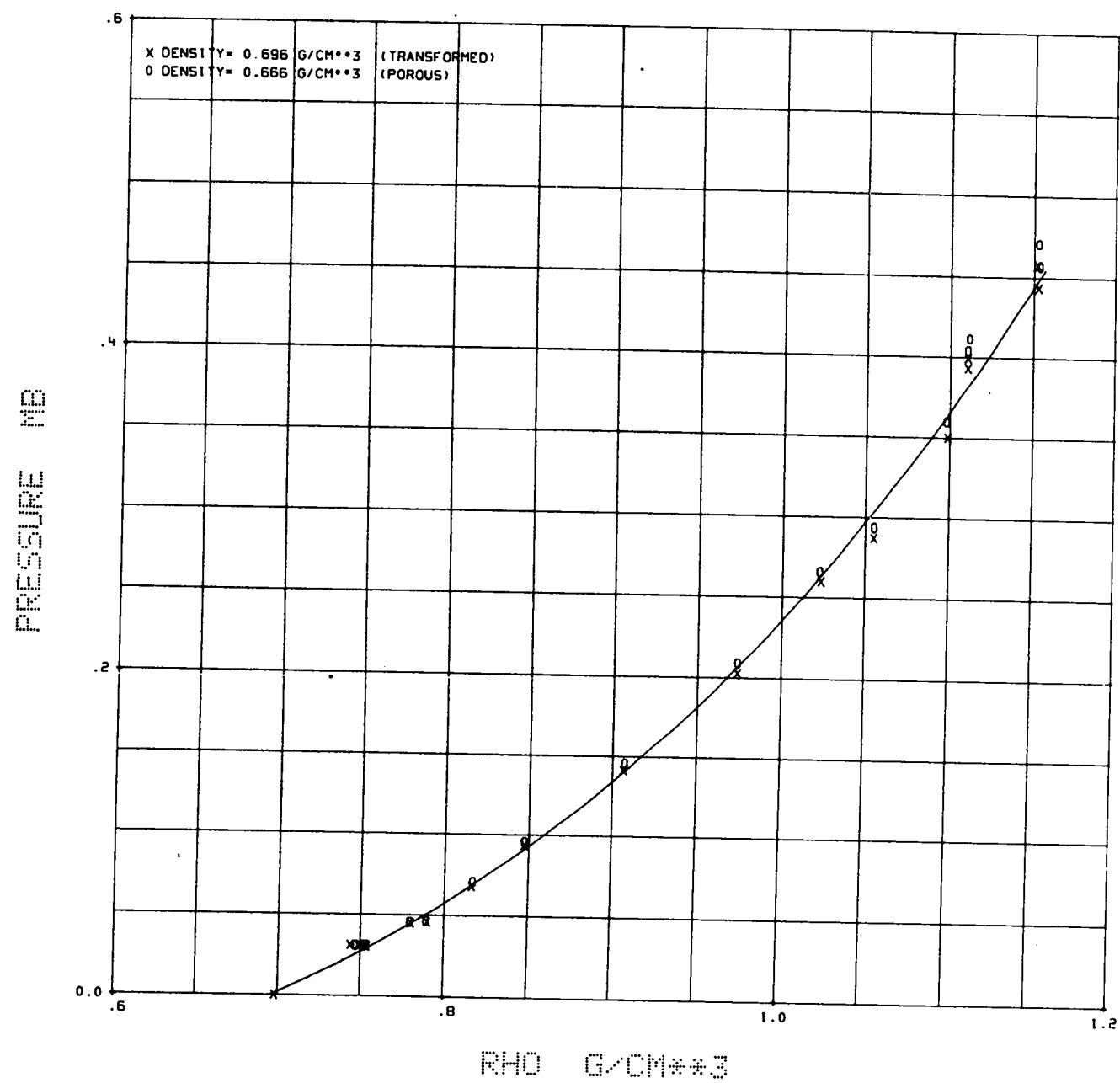


Fig. 6. Pressure-density Hugoniot data for Li^6H . Experimental porous Hugoniot data and transformed crystal-density Hugoniot points are shown. The curve is the fit shown in Fig. 5.

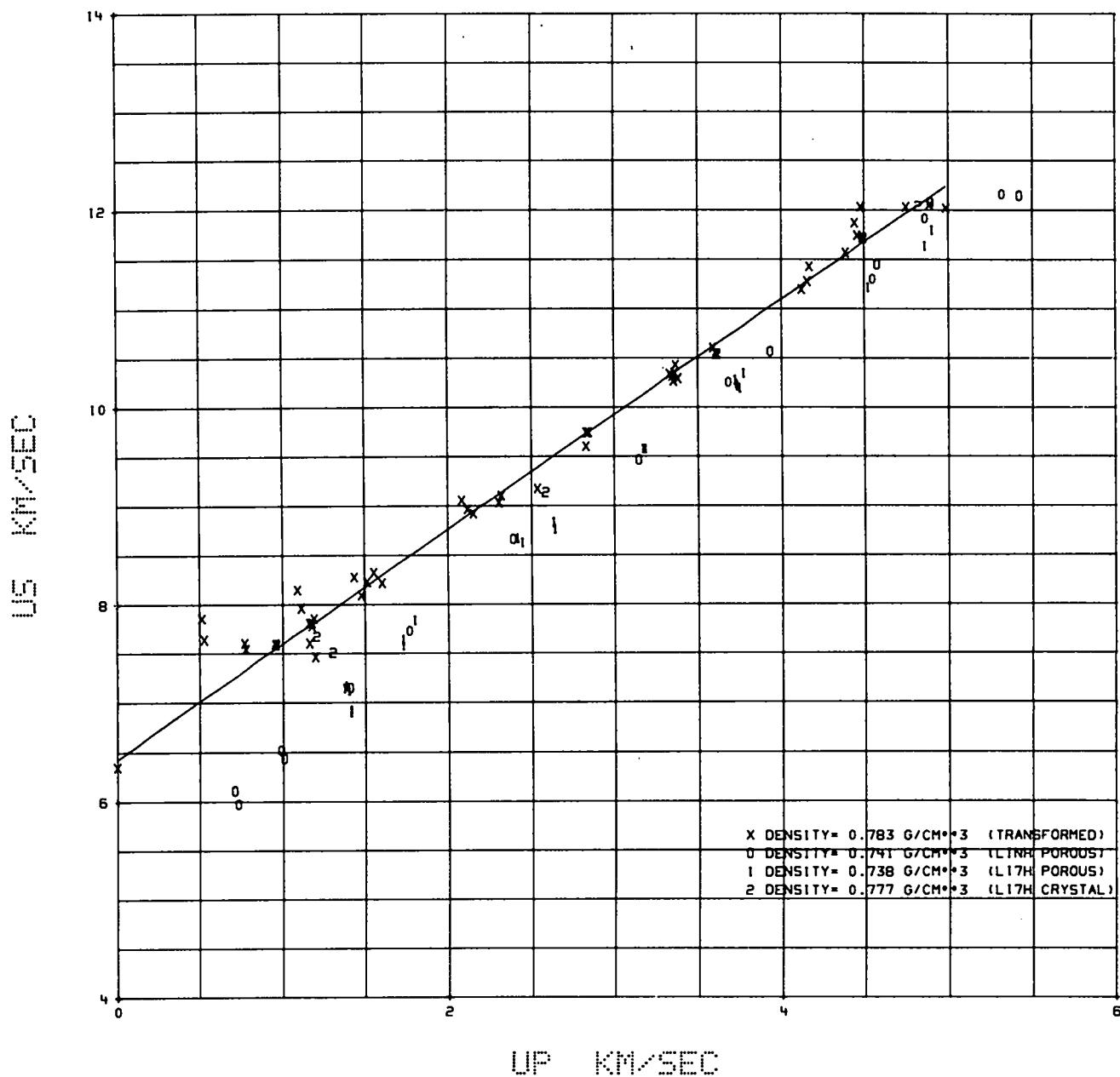


Fig. 7. Shock velocity - particle velocity Hugoniot data for Li^6H and Li^7H . Experimental porous Hugoniot data for Li^6H and Li^7H are shown along with single-crystal Li^7H data and transformed crystal-density points of all the data. A linear least-squares fit of the transformed crystal-density points is shown for points having $U_p > 0.9$ km/sec. No difference between the Li^6H and Li^7H Hugoniot data is resolved.

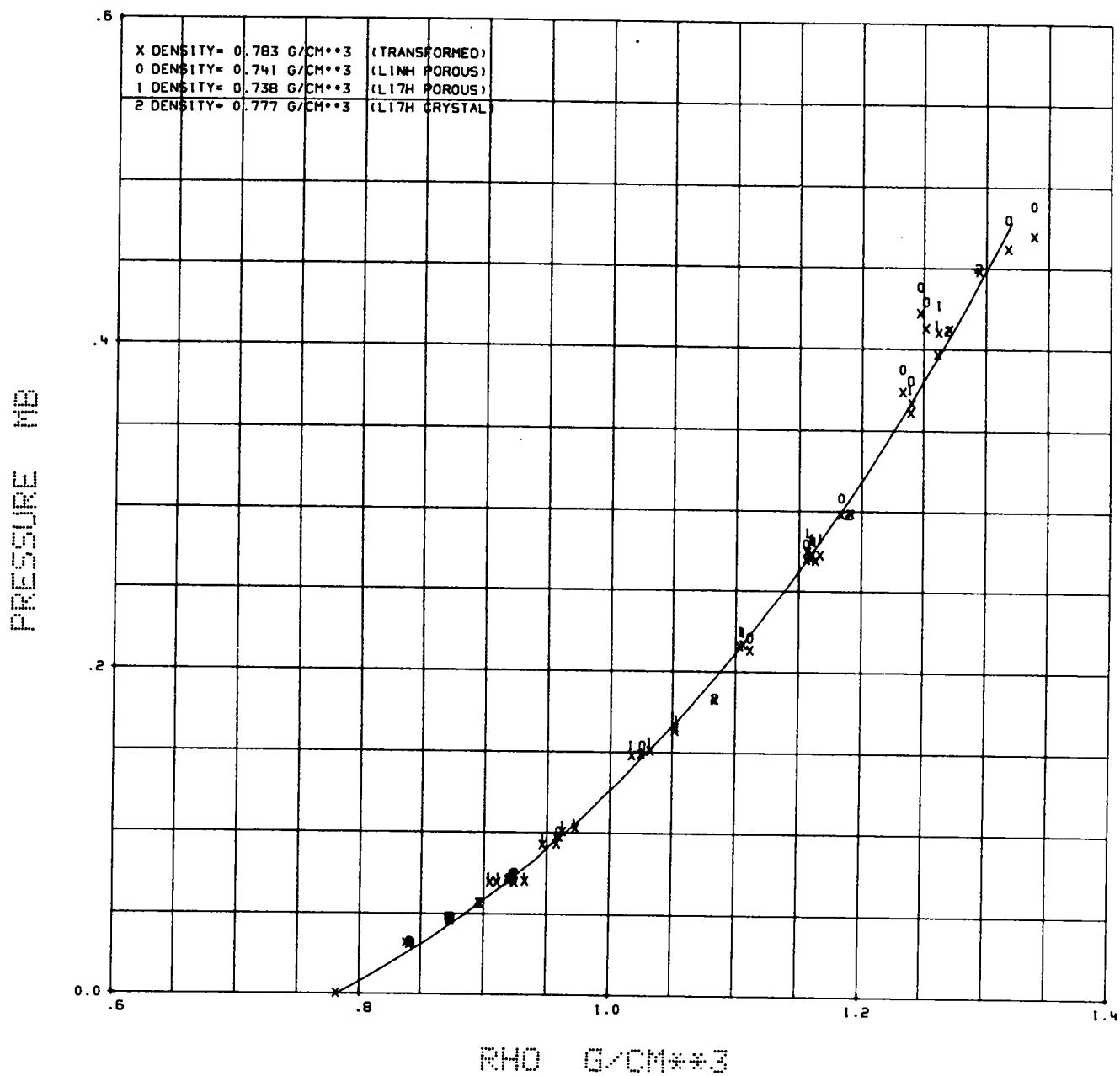


Fig. 8. Pressure-density Hugoniot data for Li⁶H and Li⁷H. Experimental porous Hugoniot data for Li⁶H and Li⁷H are shown along with experimental single-crystal Li⁷H data and transformed crystal-density points for all the data. The curve is the fit shown in Fig. 7.

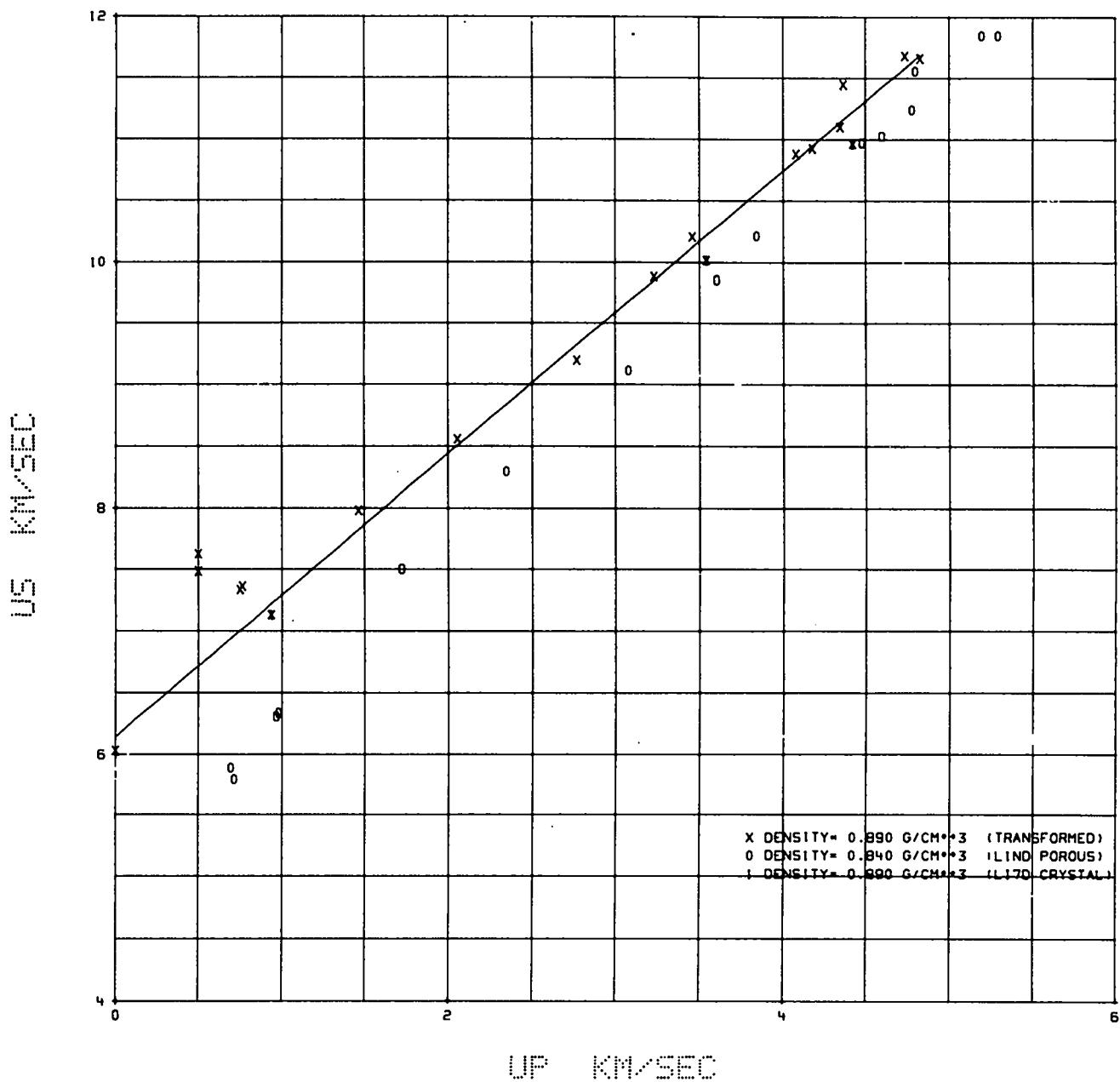


Fig. 9. Shock velocity - particle velocity Hugoniot data for Li^6D and Li^7D . Porous Hugoniot data for Li^6D and experimental single-crystal data for Li^7D are shown along with transformed crystal-density points. A linear least-squares fit of the experimental and transformed crystal-density points is shown for points having $U_p > 0.9$ km/sec. The three single-crystal data points lie below the fit. The reason for this difference is not known.

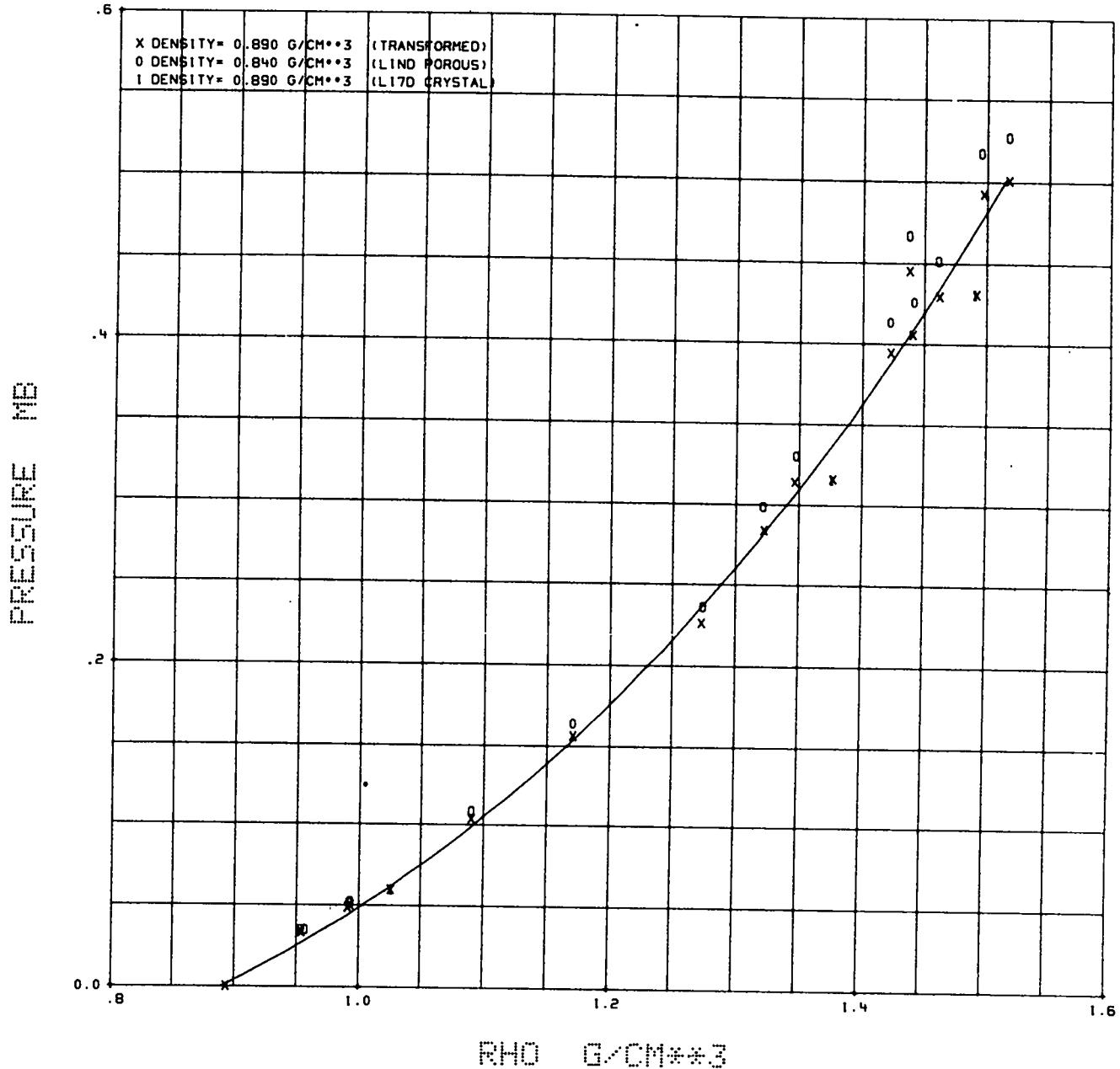


Fig. 10. Pressure-density Hugoniot data for Li^6D and Li^7D . Porous Hugoniot data for Li^6D and experimental single-crystal data for Li^7D are shown along with transformed crystal-density points. The curve is the fit shown in Fig. 9.

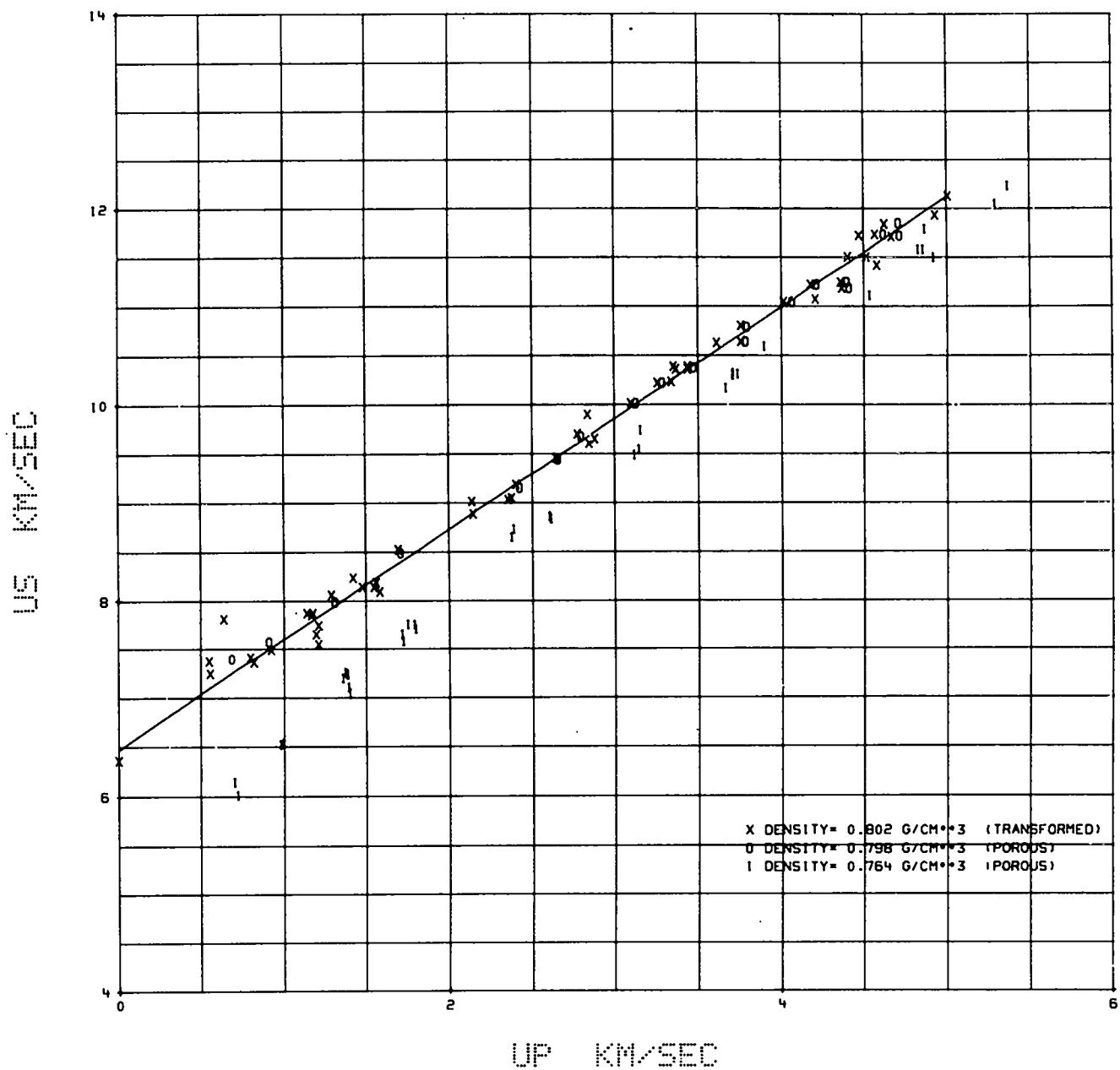


Fig. 11. Shock velocity - particle velocity Hugoniot data for Li^6D . Experimental porous Hugoniot data ($\rho_0 = 0.798$ and 0.764 g/cm^3) and transformed crystal-density Hugoniot points are shown. A linear least-squares fit of the transformed crystal-density points is shown for points having $U_p > 0.9 \text{ km/sec}$.

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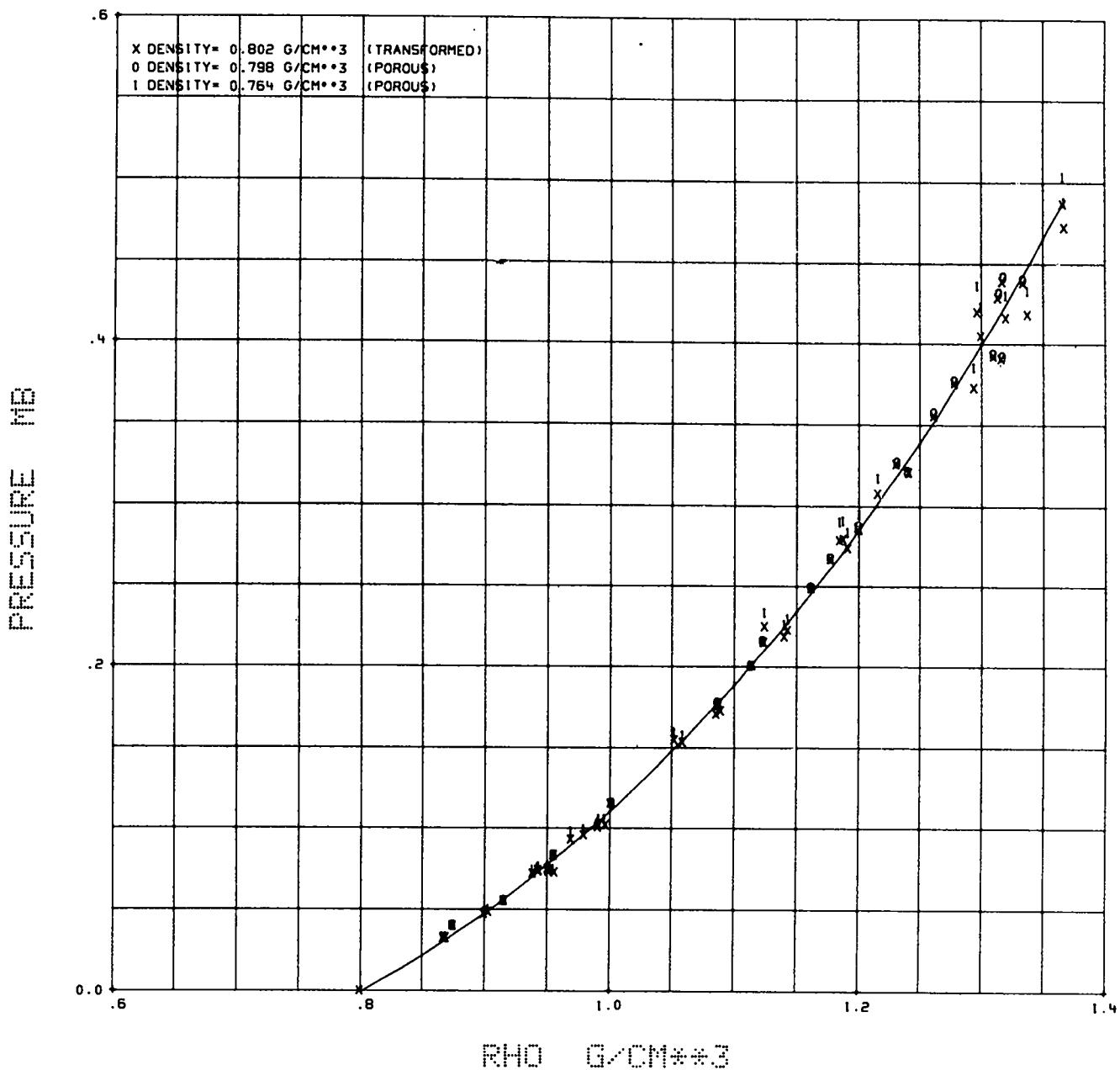


Fig. 12. Pressure-density Hugoniot data for Li⁶D. Experimental porous Hugoniot data ($\bar{\rho}_0 = 0.798$ and 0.764 g/cm³) and transformed crystal-density Hugoniot points are shown. The curve is the fit shown in Fig. 11.

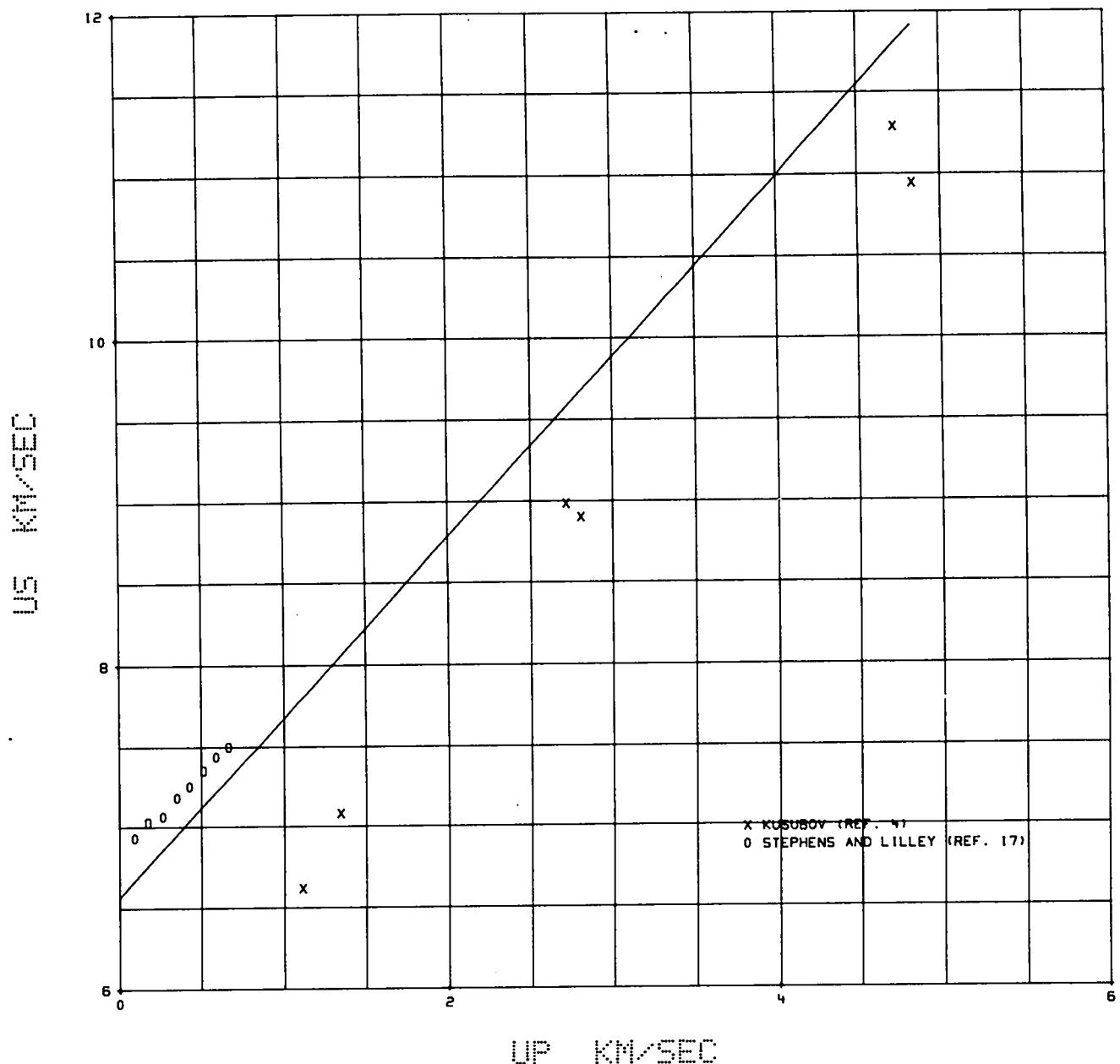


Fig. 13. Summary of other investigators' Li⁶H Hugoniot data. The straight line shows the fit reported in Table XV.

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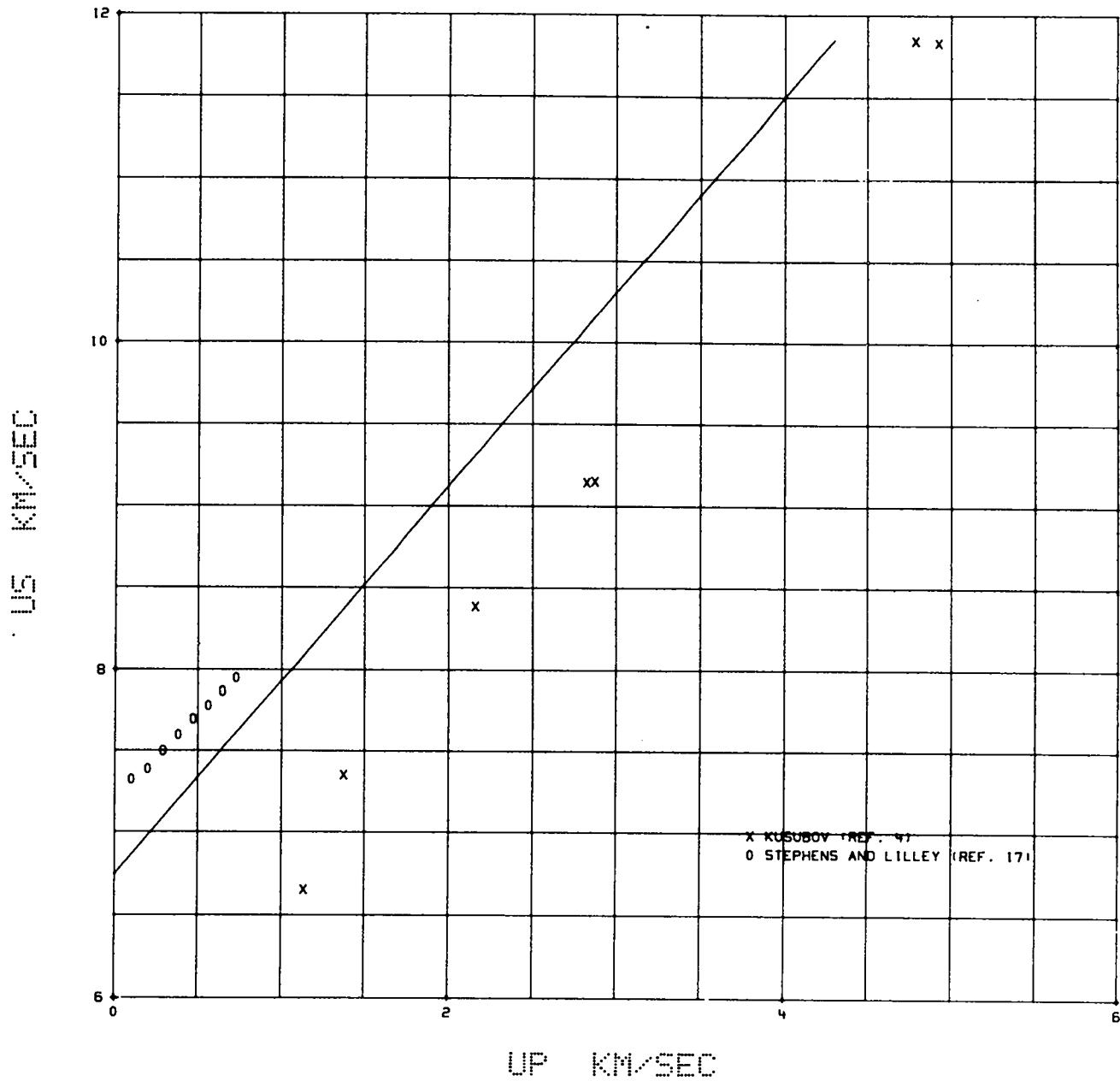


Fig. 14. Summary of other investigators' Li⁶D Hugoniot data. The straight line shows the fit reported in Table XV.

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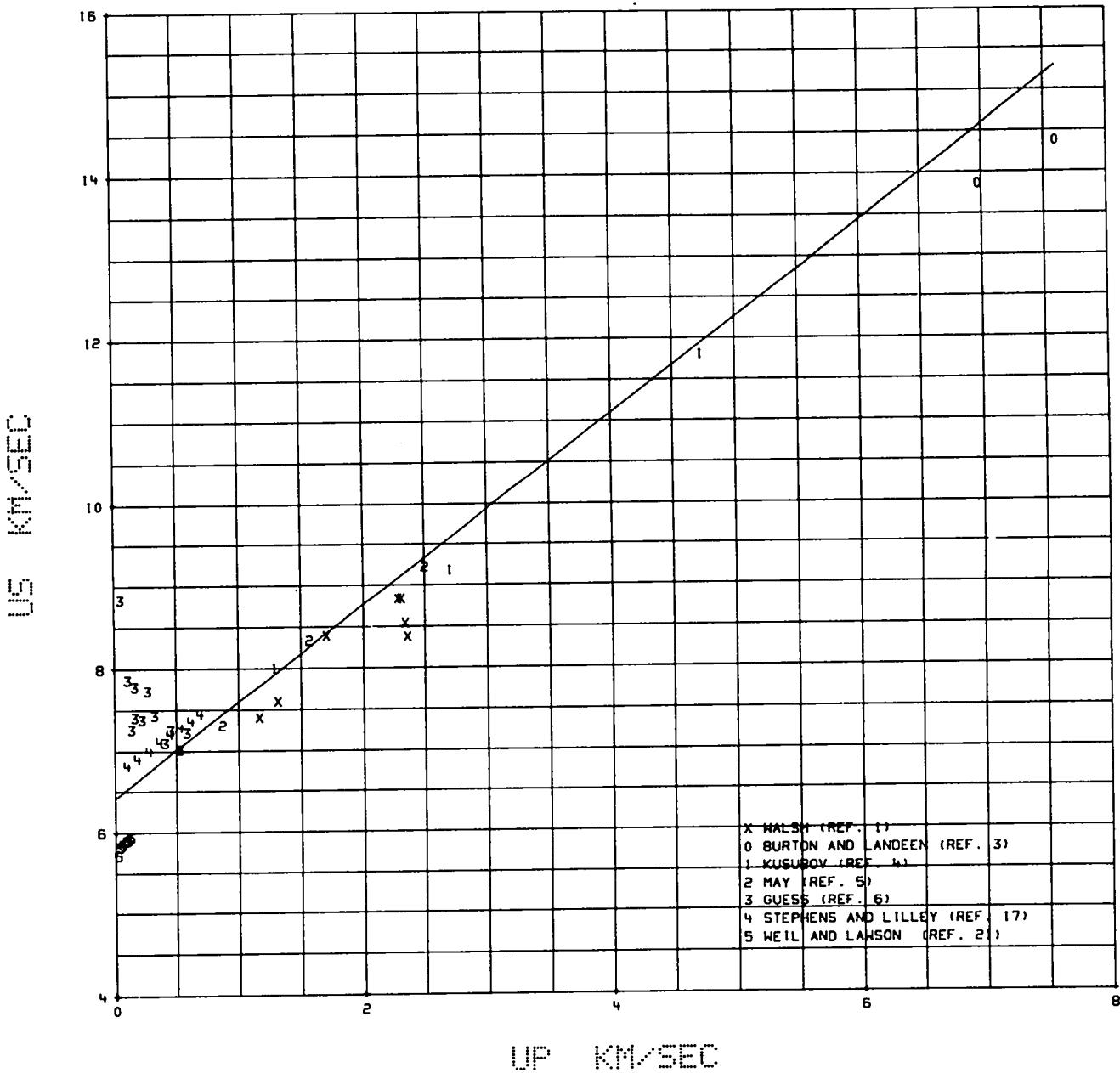


Fig. 15. Summary of other investigators' LiH Hugoniot data. The straight line shows the fit reported in Table XV.

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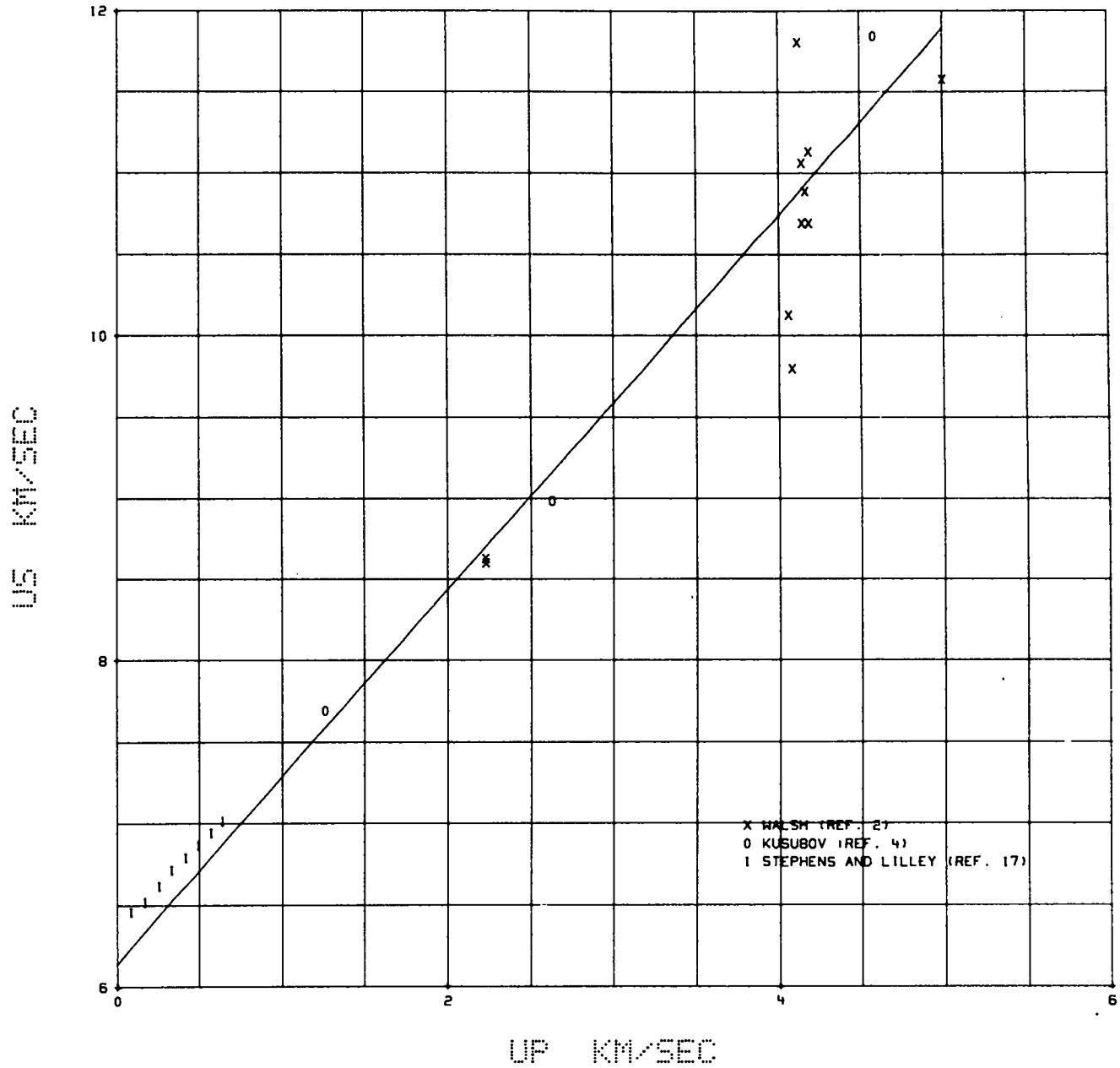
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Fig. 16. Summary of other investigators' LiD Hugoniot data. The straight line shows the fit reported in Table XV.

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