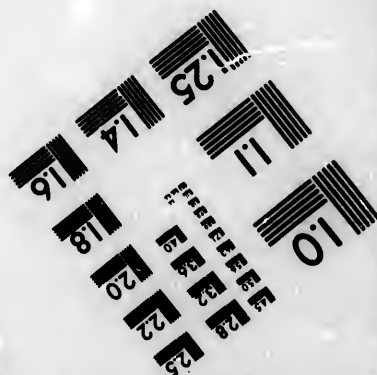
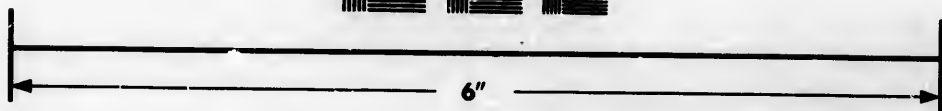
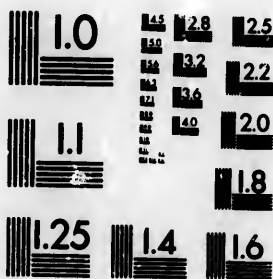


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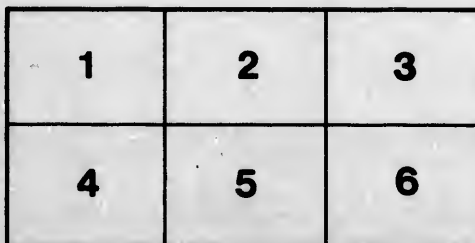
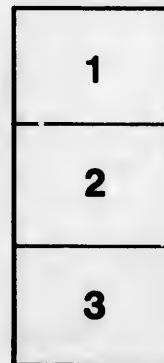
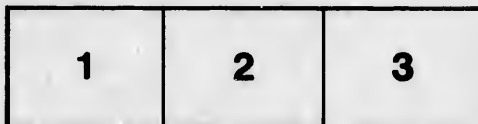
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I.—A Table of the Cubical Expansions of Solids.

By Prof. J. G. MACGREGOR, D.Sc.

(Read May 23, 1888.)

It has been found by experiment that in general the volume of any body at temperature t° may be expressed in terms of its volume at any other temperature τ° , and of the difference of temperatures, by means of the expression,

$$V_t = V_\tau (1 + A(t - \tau) + B(t - \tau)^2), \quad (1)$$

where V_t and V_τ are the volumes at t° and τ° respectively and A and B are constants for the substance of which the body consists. In some cases an additional term $C(t - \tau)^3$ is necessary. But in general the value of C is so small that it may be neglected.

As the constants A, B, C have different values for different values of τ , it is convenient to choose some one temperature as a temperature of reference, and for this purpose the temperature $0^\circ C$ is now universally chosen, the above expression becoming therefore,

$$V_t = V_0 (1 + at + bt^2). \quad (2)$$

The constants a and b (and c also, the coefficient of t^3 , in cases in which a term in t^3 is found necessary) having been determined for any substance, the volume of any body of that substance, whose volume at $0^\circ C$ is known, may be determined at any other temperature within the temperature limits of the experiments by which the values of a, b and c were found.

Density may be substituted for volume in the above formula, provided the signs of a and b be changed. For if ρ_t and ρ_0 are the densities of a substance at t° and 0° respectively, we have, in general,

$$\rho_t / \rho_0 = V_0 / V_t = 1 / (1 + at + bt^2) = 1 - at - bt^2,$$

since a and b are, in general, small quantities.

In the case of isotropic solids, length may be substituted for volume in the above formula, provided the constants a and b be divided by 3, change of length in any direction in such cases being numerically equal to one third of the corresponding change of volume. The formula in the case of the linear expansion of such solids becomes therefore:—

$$L_t = L_0 (1 + \frac{a}{3}t + \frac{b}{3}t^2). \quad (3)$$

In the case of ælotropic solids, the linear expansion is different for different directions, and must therefore be specially determined.

To determine the volume at t'' when that at t' is given, we have:—

$$\begin{aligned} V_t &= V_o(1 + at + bt^2), \\ \text{and} \quad V_o &= V_o(1 + at' + bt'^2). \\ \text{Hence} \quad V_o &= V_t \frac{1 + at' + bt'^2}{1 + at + bt^2}, \\ &= V_t(1 + a(t' - t) + b(t'^2 - t^2)), \end{aligned} \quad (4)$$

since a and b are small quantities.

The mean coefficient of thermal expansion between two temperatures is by some writers defined as the change of volume per degree and per unit volume at the lower temperature, and by others as the change of volume per degree and per unit volume at the temperature of reference (0°C). In terms of the symbols used above, it is in the former case the value of $\frac{V_o - V_t}{V_t(t' - t)}$, and in the latter case the value of $\frac{V_o - V_t}{V_o(t' - t)}$. From (4) it follows at once that:—

$$\frac{V_o - V_t}{V_t(t' - t)} = a + b(t' + t) \quad (5)$$

provided a and b are so small that their powers and product may be neglected. From (2) and a similar equation with t' substituted for t , it follows that:—

$$V_o - V_t = V_o(a(t' - t) + b(t'^2 - t^2))$$

exactly, and therefore that

$$\frac{V_o - V_t}{V_o(t' - t)} = a + b(t' + t), \quad (6)$$

whatever the magnitudes of a and b may be. Provided a and b are sufficiently small, therefore, the mean coefficient has the same value between given temperature limits according to both modes of definition; and if a and b are known, this value may be determined for any temperature range.

The "true" coefficient of thermal expansion at any temperature is the rate at which volume varies with temperature at that temperature per unit volume at zero. It is thus $\frac{1}{V_o} \frac{dV_t}{dt}$, and by differentiation of (2) is seen to have the value $a + 2bt$. The true coefficient at a given temperature is clearly the mean coefficient (per unit volume at 0°C) between two temperatures indefinitely near one another and including the given temperature; and the above value is also obtained from (6) by noting that ultimately $t' + t = 2t$.

Sometimes, but rarely, the true coefficient at any temperature is defined as the rate at which volume varies with temperature at that temperature per unit volume at that temperature—in symbols, $\frac{1}{V_t} \frac{dV_t}{dt}$; and by (5), its approximate value, provided a and b be small, is $a + 2bt$, the same as the value of the true coefficient according to the former definition.

The following table contains the values of a and b (and of c also in cases in which c is found to have an appreciable value) in the case of the more important and interesting solids. A table containing an exhaustive list of the determinations of these constants would be so long that in most cases I have thought it well to give only the most recent and most accurate determinations, though in some important cases, as in that of glass, a

large number of such determinations have been given. Many observers of thermal expansion have determined, not the values of the constants a and b , but the values of the mean coefficient throughout given ranges of temperature. In many cases, especially in cases in which the values of a and b are not known, these results have been given. They are contained in the fourth column, and in the fifth column are stated the temperature ranges to which they apply. In general, also, these temperatures are the ranges throughout which observations were made. A more complete list of the older determinations of mean coefficients of expansion will be found in Part III of the "Constants of Nature," by Prof. F. W. Clarke, published in Vol. XIV of the Smithsonian Miscellaneous Collections, but Prof. Clarke's tables do not give the values of the constants a and b .

SUBSTANCE.	$a \times 10^6$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Coefficient $\times 10^6$	Temperature Range. (°C.)	OBSERVER.
Aluminium (cast).	6664	3435	7008	0—100	Fizeau
Do. (commercial).	6660	0—100	Calvert, Johnson & Lowe
Do.	7439.4	0—100	L. Pfaff
Do.	7062	0—100	Glatzel
Alums (in powder)—						
Ammonium-Aluminium Alum	2840	0—20	Spring
.....	2585	0—60	"
.....	5343	0—100	"
Potassium-Aluminium Alum	2635	0—20	"
.....	4198	0—60	"
.....	31118	0—90	"
Rubidium-Aluminium Alum	2070	0—20	"
.....	2635	0—60	"
.....	5045	0—100	"
Cesium-Aluminium Alum	2610	0—20	"
.....	2621	0—60	"
.....	7699	0—100	"
Potassium-Chromium Alum	2330	0—20	"
.....	4968	0—60	"
.....	59503	0—80	"
Ammonium Sulphate.	8335	0—20	"
.....	9051	0—60	"
.....	11191	0—100	"
Antimony	2770	3970	3167	0—100	Matthiessen
Do. (pure).	2940	0—100	Calvert, Johnson & Lowe
Do. (crystal).	3387	870	3379	0—100	Fizeau

SUBSTANCE.	$a \times 10^3$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Coefficient $\times 10^6$	Temperature Range. ($^{\circ}\text{C}$.)	OBSERVER.
Aragonite.....	6105	39.5—40.5	Fizeau
Arsenic (sublimed).....	1158.6	6480	1800	0—100	"
Augite	2305	1800	2485	0—100	"
Beryl.....	105	0—100	F. Pfaff
Bismuth.....	3502	4400	3948	0—100	Matthiessen
Do.	3900	0—100	Calvert, Johnson & Lowe
Do. (crystal).....	3706	4155	4121	0—100	Fizeau
Brass (71 % by mass, 66.15 % by volume, of Cu.).....	5101	5580	5719	0—100	Matthiessen
Do. (71.5 % of Cu, 27.7 % of Zn, 0.3 of Sn, 0.5 of Pb).....	5341	2940	5637	0—100	Fizeau
Bromide of Silver (HlgBr).....	10100	—60— 0	Rodwell
	10429	0—100	"
	11033	0—400	"
Bronze (86.3 % of Cu, 9.7 % of Sn, 4 of Zn.).....	5101	3060	5406	0—100	Fizeau
Do. (25 % of Sn.).....	5532	16.6—100	Daniell
Cadmium.....	8078	14000	9478	0—100	Matthiessen
Do.	9363	0—100	Glatzel
Do. (compressed powder).....	8816	4800	9306	0—100	Fizeau
Calc spar.....	1541	39.5—40.5	"
Do.	2010	0—100	F. Pfaff
Caoutchouc (commercial grey).....	67000	0— 30	Russnor
Do. (vulcanized, unstretched).....	95784	14— 37	Lebedeff
Do. do. do.....	67646	11—295	"
Do. (do., stretched).....	56158	14.5—37.5	"
Do. do. do.	66906	12.5— 30	"
Carbon (graphite).....	877	16.6—100	Daniell
Do. do.	632	10.6—350	"
Do. do.	2237	1515	2388	0—100	Fizeau
Do. (gas-coke).....	1488	1650	1653	0—100	"
Do. (diamond).....	181.2	2160	390	0—100	"
Cassiterite.....	925.6	1355	1061	0—100	"

CUBICAL EXPANSION OF SOLIDS.

7

SUBSTANCE	$a \times 10^4$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Co-efficient $\times 10^6$	Temperature Range. ($^{\circ}\text{C}$)	OBSERVER.
Chloride of Silver (AgCl).....	9245	-60— 0	Rodwell
	9492	0—100	"
	9865	0—400	"
Cobalt (compressed powder).....	3612	1200	3732	0—100	Fizeau
Copper.....	5150	0—100	Dulong & Petit
Do.	5650	0—300	"
Do.	4443	5550	4098	0—100	Matthiessen
Do.	4788	3075	5094	0—100	Fizeau
Do. (native, Lake Superior).....	4850	2740	5124	0—100	"
Do.	5265	0—100	L. Pfaff
Do.	5233	0—100	Rodwell
Do.	5115	0—100	Glatzel
Copper-Gold Alloy (60% % by mass, 48.06% by volume, of Gold)....	4015	6420	4657	0—100	Matthiessen
Copper Oxide (CuO)....	27	3150	342	0—100	Fizeau
Corundum.....	1443	3275	1771	0—100	"
Darcet's fusible metal (Bi ₁₅ Sn ₁₀ Pb ₅).....	52143	0— 35	Spring
	-36026	35— 50	"
	102534	50— 90	"
Diopside.....	2330	0—100	F. Pfaff.
Ebonite (See Vulcanite)..
*Emerald.....	16	1900	206	0—100	Fizeau
Epidote.....	2627	3830	2410	0—100	"
Felspar (Adular).....	1399.5	1900	1590	0—100	"
Fluorspar.....	5851	0—100	F. Pfaff
Galena.....	5578	0—100	"
Garnet.....	2543	0—100	"
German Silver.....	5509	0—100	L. Pfaff
Glass (crystal of Choisy- le-Roy).....	2231	50	100—280	Bosscha †
Do. do.	2523	70	"	"
Do. do.	2439	300	"	"
Do. (ordinary)....	2578	1330	0.37	"	"

* See also Beryl.

† Bosscha's constants were obtained by a recalculation of Regnault's experimental results.

SUBSTANCE	$a \times 10^9$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Co-efficient $\times 10^6$	Temperature Range (°C.)	OBSERVER.	
Glass (crystal of Choisy-le-Roy) (ordinary).....	2507	1020	0.51	100—280	Bosscha *	
Do. do.	2281	1540	0.46	"	"	
Do. (green).....	2143	890	0.37	"	"	
Do. (Swedish) ...	2218	3570	1.15	"	"	
Do. (crystal, Choisy-le-Roy) {	2270	0—10	Regnault	
	2270	0—50	"	
	2280	0—100	"	
	2300	0—150	"	
	2310	0—200	"	
	2320	0—250	"	
	2330	0—300	"	
	2340	0—350	"	
	Do. (ordinary) {	2628	0—10	"
		2687	0—50	"
.....		2761	0—100	"	
.....		2835	0—150	"	
.....		2908	0—200	"	
.....		2982	0—250	"	
.....		3056	0—300	"	
Do. (rod) {	2410	1700	3131	0—350	"	
				2580	0—100	Dulong & Petit	
				2750	0—200	"	
.....	3040	0—300	"		
Do. (tube)	2435	0—100	Lavoisier & LaPlace	
Do. (plate).....	2673	0—100	"	
Do. (crown plate).....	2627	0—100	"	
Do. do.	2693	0—100	"	
Do. do.	2753	0—100	"	
Do. (thin tube).....	588	31500	0—30	Hällström	
Do. (white, French)...	2553	0—100	Kopp	
Do. (white, tube).....	2648	0—100	Regnault	
Do. (do. globulo)....	2592	0—100	"	
Do. (do. do.)....	2514	0—100	"	

* Bosscha's constants were obtained by a recalculation of Regnault's experimental results.

CUBICAL EXPANSION OF SOLIDS.

SUBSTANCE.	$a \times 10^6$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Co-efficient $\times 10^6$	Temperature Range. ("C.)	OBSERVER.
Glass (green, tubo).....	2200	0—100	Regnault.
Do. (do. globule)....	2132	0—100	"
Do. (Swedish, tube)....	2363	0—100	"
Do. (do. globule)...	2441	0—100	"
Do. (hard, French, tube)	2142	0—100	"
Do. (do. globule)	2242	0—100	"
Do. (ordinary), from	2431	0—100	"
to...	2758	0—100	"
Do. (ord. crystal, tube).	2101	0—100	"
Do. (do. globule)	2330	0—100	"
Do. (Cholsy-le-Roy crystal) from.....	2144	0—100	"
to..	2442	0—100	"
Do. (St. Gobain).....	2141	2370	2378 ⁴	0—100	Fizeau
Do. (3 pts. of sand, 2 of Pb, and 1 of alkali) .	2187	0—100	Matthiessen
Do. (Thuringian soft, tube).....	3585	0—100	Weinhold
Do. (do. vessel)....	3590	0—100	"
Do. (before heating)...	2784	0—100	Crafts
Do. (after having been heated for 100 hours in boiling sulphur)...	2741	0—100	"
Gold.....	4075	3361	4411	0—100	Matthiessen
Do. (cast).....	4220	1245	4353	0—100	Fizeau
Do. (pure).....	4140	0—100	Calvert, Johnson & Lowe
Do.	4401	0—100	L. Pfaff
Gutta-percha (purified)...	49606	496000	69500	0— 40	Russner
Gypsum	6700	6940	7304	0—100	Fizeau
Do.	7500	0—100	F. Pfaff
Hornblende.....	2355	3025	2658	0—100	Fizeau
Ice.....	11250	-20— 0	Brunner
Do	15850	-20— 0	Flicker & Goisler
Do	3500	-12— 0	Marchand
Indium (cast).....	7424	63570	13782	0—100	Fizeau
Iodine.....	{ 23500 (?) 31231	40·3—107	Billet

SUBSTANCE.	$a \times 10^6$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Coefficient $\times 10^6$	Temperature Range. ($^{\circ}$ C.)	OBSERVER.
Iodides—						
AgI (crystal)	—200.0	—754.5	—282	0—100	Fizeau
Do. (compressed precipitate)	—218	—2400	—458 *	0—100	"
Do. (cast)	—240	—2100	—240	0—100	"
Do.	—1740	70—142	Rodwell
Do.	—450000	142—145.5	"
Do.	2844	145.5—300	"
Do.	—8010	—60—142	"
Do.	2843	142—400	"
PbI ₂ (cast)	9378.6	8700	10255	0—100	Fizeau
Do.	7614	0—205	Rodwell
Do.	8317	0—205	"
Do.	63780	205—253	"
Do.	18000	253—m.p.	"
HgI ₂ (cast)	9523	0—200	Rodwell & Elder
Iridium (cast)	2005	1185	2124	0—100	Fizeau
Iron (forged)	3061	0—100	Lavoisier & Laplace
Do. (wire drawn)	3705	0—100	"
Do. (soft)	3408	2775	3684	0—100	Fizeau
Do. (cast, grey)	3019	2055	3235	0—100	"
Do. (cast)	3360	0—100	Calvert, Johnson & Lowe
Do. (wrought)	3570	0—100	"
Do.	3742	0—100	L. Pfall
Do.	3307	0—100	Rodwell
Do.	4161	0—100	Glatzel
Iron-glance (oxide, rhombohedral)	2241	3215	2565	0—100	Fizeau
Iron-stone (magnetite)	2862	0—100	F. Pfall
Iron Pyrites	3025	0—100	"
Do.	2721	39.5—40.5	Fizeau
Lipowitz's Fusible Metal (Bi ₁₁ Pb ₆ Cd, Sn ₂)	15400	0—25	Spring
Do.	—39514	25—40	"
Do.	38381	40—60	"
Lead	8177	2220	8399	0—100	Matthiessen

* The negative coefficient shewn qualitatively to extend to -10° C.

CUBICAL EXPANSION OF SOLIDS.

SUBSTANCE.	$a \times 10^6$	$b \times 10^{11}$	$c \times 10^{18}$	Mean Coefficient $\times 10^6$	Temperature Ranges (°C.)	OBSERVER.
Lead (pure).....	9030	0-100	Calvert, Johnson & Lowe
Do. (cast).....	8485	3585	8844	0-100	Fizeau
Do.	8725	0-100	L. Pfaff
Do.	9003	0-100	Rodwell
Do.	8808	0-100	Glatzel
Lead - Antimony alloy } (Sn, Pb, 22.28% by vol. } of Pb).....	0200	0880	7188	0-100	Matthiesson
Do. (Sn Pb, 82.00% by } vol. of Pb).....	8087	3320	8419	0-100	"
Magnesium (cast).....	7201	10200	8280	0-100	Fizeau
Do.	7042	0-100	L. Pfaff
Nickel (comp. powder)...	3752	1005	3858	0-100	Fizeau
Osmium (semi-fused)....	1700	3270	2037	0-100	"
Palladium.....	3032	2800	3312	0-100	Matthiesson
Do. (forged, annealed)	3370	1980	3567	0-100	Fizeau
Paraffin.....	58400	90200	0-33	Russner
	260000	33.5-37.7	"
	666000	37.7-41	"
Do. (high boiling } point).....	115000 *	690000 *	41-52	"
	31985	0-15.55	Rodwell
	36090	15.55-37.6	"
Do. (Rangoon, melt- } ing at 56°C)	143118	37.6-48.85	"
	71651	148890	244358	48.85-61.11	"
Pewter.....	70989	0-56	Fizeau
Do.	6099	16.6-100	Daniell
Do.	5982	16.6-206	"
Phosphorus.....	37988	0-37.5	Erman
Do.	38300	0-44	Kopp
Do.	20000	11500	20506	0-44	Pisati & de Franchis
Platinum.....	2644	16.6-100	Daniell
Do.	2650	0-100	Dulong and Petit
Do.	2750	0-300	"
Do.	2554	1040	2658	0-100	Matthiesson
Do. (cast).....	2603	1170	2721	0-100	Fizeau

* These constants are for the formula: $-V_t = V_0 (1 + at + bt^2)$.

SUBSTANCE.	$a \times 10^8$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Co-efficient $\times 10^7$	Temperature Range. ($^{\circ}$ C.)	OBSERVER.
Platinum.....				2583	0—100	J. Pfaff
Do. (commercial)...				2040	0—100	Culvert, Johnson & Lowe
Platinum-iridium alloy (10% of Ir.) (cast).....	2561	1140		2675	0—100	Fizeau
Porcelain (Bayeux).....				4950	1000—1400	Deville & Troost
Do. do.				6000	about and above 1500	Weinhold
Do.				866	0—99	"
Potassium.....	24070	17980		24990	0—50	Hagen
	23800	23870				
				10570	0—20	Spring
Potassium Chromate.....				10731	0—60	"
				11344	0—100	"
				9520	0—20	"
Potassium Sulphate.....				10037	0—60	"
				12645	0—100	"
Quartz.....	3358	3265		3685	0—100	Fizeau
Do.				3840	0—100	F. Pfaff
Rhodium(semi-fused)...	2453	1215		2574	0—100	Fizeau
Rock Salt.....	11578	6735		12252	0—100	"
Rose's Fusible Metal † ..	6784.7	-181580	553.07†		0—90	Kopp
				33575	0—40	Spring
Do. (Bi, Sn, Pb).....				-129922	40—55	"
				54546	55—90	"
				9760	0—20	"
Rubidium Sulphate.....				10020	0—60	"
				11148	0—100	"
Ruthenium (semi-fused).....	2552	4215		2973	0—100	Fizeau
Rutile.....	2196	2225		2419	0—100	"
Selenium (cast).....	9702	16725		11376	0—100	"
				14780	0—20	Spring
Do. (crystallized).....				17430	0—60	"
				19810	0—100	"

† Kopp uses the formula: $-V_t = V_0(1 + at + bt^2 + ct^3 + dt^4)$; a , b , and c are given in the table; $d = 10^{-13} \times 5250$.

‡ Erman gives the formula: $-V_t = V_0(1 + 10^{-8}.21864t - 10^{-8}.93485 \sqrt{(t - 34^{\circ}.9)(78^{\circ}.5 - t)})$, the temperatures being expressed in Reaumur degrees and the imaginary values of the irrational factor being put equal to 0.

SUBSTANCE.	$a \times 10^6$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Coefficient $\times 10^6$	Temperature Range. ($^{\circ}C$)	OBSERVER.
Selenium (com. powder) {	13070	0—20	Spring
	16440	0—60	"
	17510	0—100	"
Silicium (cast).....	2086	2535	2340	0—100	Fizeau
Silver.....	5426	4050	5831	0—100	Matthiessen
Do. (cast).....	5587	2205	5808	0—100	Fizeau
Do.	5516	5—26	Plantamour & Hirsch
Do.	5573	0—100	L. Pfaff
Do.	5792	0—100	Rodwell
Silver-platinum alloy (33.4 % by mass, and 19.65 % by vol. of Pt.).....	4246	3220	4568	0—100	Matthiessen
Sodium.....	14178	52000	26	16674	0—40	de Lucchi.
Do.	23670	0—90	"
Do. {	20510	22730	21630	0—50	Hagon
	20280	25730			
Solder (2 parts Pb, 1 Sn.)	7524	0—100	Sneaton
Speculum Metal.....	5790	0—100	"
Steel (not tempered).....	3238	0—100	Lavoisier & LaPlace
Do. (tempered yellow)	4158	0—100	"
Do. (French cast, tempered).....	4086	0—100	Fizeau
Do. (French cast, annealed).....	3339	0—100	"
Do. (English cast, annealed).....	3103	2280	3331	0—100	"
Do. (wire drawn).....	3768	0—100	Glatzei
Do. (soft).....	3060	0—100	Calvert, Johnson & Lowe
Do. (bar at max. softness).....	3270	0—100	"
Do. (same bar at max. hardness).....	4260	0—100	"
Sulphur.....	10458	265880	-146.73	0—78	Kopp
Do.	-83804	7878840 §	78—115	"
Do. (Sicilian crystal) .	15222	50440	20244	0—100	Fizeau
Do. (crystallized from solution in Carbon Bisulphide) {	21220	0—20	Spring
	24380	0—60	"
	35410	0—100	"

§ These constants are for the formula: $-V_t = V_{78} (1 + a(t-78) + b(t-78)^2)$

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SUBSTANCE	$a \times 10^6$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Coefficient $\times 10^6$	Temperature Range. ($^{\circ}C$)	OBSERVER
Sulphur (Sicilian).....	24300	0—20	Spring
Do.	25000	0—60	"
Do.	26000	0—100	"
Do. (crystallized from solution in Carbon Bi-sulphide)	12800	186000	153	29468	0—60	Russner
Do. (crystal).....	17596	20.5—21.5	Schrauf
Do. (natural crystal)	19991	0—60	Sciuchione
Do. (after heating to $140^{\circ}C$).....	—4410	62—70	"
Do. (after heating to $140^{\circ}C$).....	28483	70—110	"
Do. (after heating to $140^{\circ}C$).....	20958	0—82	"
Do. (after heating to $140^{\circ}C$).....	—13731	82—85	"
Do. (after heating to $240^{\circ}C$).....	201700	85—100	"
Do. (after heating to $240^{\circ}C$).....	18246	0—90	"
Do. (after heating to $240^{\circ}C$).....	—29385	90—92	"
Do. (after heating to $240^{\circ}C$).....	195180	92—100	"
Tellurium (crystallized)	10320	0—20	Spring
Do. (comp. powder)	11210	0—60	"
Do. (comp. powder)	11060	0—100	"
Do. (cast).....	4335	8625	10410	0—20	"
Do. (cast).....	4335	8625	10110	0—60	"
Do. (cast).....	4335	8625	10630	0—100	"
Thallium (cast).....	7694	17115	5196	0—100	Fizeau
Tin.....	6100	7890	9405	0—100	"
Do. (Malacca, compressed powder).....	6281	5265	6889	0—100	Matthiessen
Do. (pure).....	6807	0—100	Fizeau
Topaz.....	6567	0—100	L. Pfaff
Tourmaline.....	2137	0—100	F. Pfaff
Type Metal.....	2181	0—100	"
Do.	6099	16.6—100	Daniell
Do.	5856	16.6—264	"
Vulcanite.....	18300	114000	24239	16.7—35.4	Kohlrausch
Do.	24600	18—19	Fuess
Do. (22.2 per cent. of Sulphur).....	40000	20—60	Russner

SUBSTANCE.	$a \times 10^6$	$b \times 10^{11}$	$c \times 10^{10}$	Mean Co-efficient $\times 10^6$	Temperature Range. (°C.)	OBSERVER.
Vulcanite (27 per cent. of Sulphur)	35000	20— 60	Russner
Do. (30.7 per cent. of Sulphur)	23000	20— 60	"
Do. (31 per cent. of Sulphur)	41000	20— 60	"
Do. (36.4 per cent. of Sulphur)	38000	20— 60	"
Wood's Fusible Metal { (Bi, Pb Cd, Sn ₂)... }	-16600	0— 25	Spring
	27827	25— 70	"
Zinc	8919	16.6—100	Daniell
Do.	8222	7000	8928	0—100	Matthiessen
Do. (pure, rolled).....	6600	0—100	Calvert and Lowe
Do. (compres'd powder)	8906	1905	8715	0—100	Fizeau
Do.	8663	0—100	Rodwell

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