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BASIC ENERGY-LEVEL
AND EQUILIBRIUM DATA FOR
ATMOSPHERIC ATOMS AND MOLECULES

Forrest R. Gilmore

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SUMMARY

This memorandum presents tables of the formation energies, dissociation energies, ionization energies, electronic energy levels, and vibrational level spacings for most atomic, diatomic, and triatomic molecules involving hydrogen, carbon, nitrogen, oxygen, and argon. Many positively and negatively charged ions are included. Tables of equilibrium fractional electronic-state populations, and graphs of the equilibrium constants for dissociation, ionization, and detachment for most of the atomic and diatomic species are appended. A brief discussion of the significance of such data precedes the tables and graphs.

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I. INTRODUCTION

An important property of any reaction is its reaction energy, i.e., the amount of energy taken up or given off when unit amounts of the reactants are converted to the products. Heats of formation, dissociation energies, ionization energies, etc., represent specific types of reaction energies. When a reaction is exothermic, the reaction energy represents the energy available for excitation or heating of the products of the reaction. When a reaction is endothermic, the reaction energy must be supplied by the thermal or excitational energy of the reactants. Consequently, endothermic reactions will be very slow when the mean thermal energy is much less than the reaction energy, unless the reactants have above-thermal excitational energy. (The converse, however, is not true; reactions are not necessarily fast just because sufficient energy is available.)

An understanding of reactions involving excited states also requires a knowledge of the excited energy levels of the reacting or product species. Moreover, for the case of thermal equilibrium (i.e., for a Boltzmann distribution of excited-state populations), such energy levels can be used to calculate "equilibrium constants" which relate the forward to the backward rates of reactions (see Section 4). Finally, since most reaction energies and energy levels can be determined (by spectroscopic, calorimetric, or other measurements) with an accuracy far surpassing that attainable in reaction-rate measurements, such data form a firm foundation on which to build the often crude and speculative edifices of reaction mechanisms and rates. The present memorandum presents such data for atoms and simple molecules involving hydrogen, carbon, nitrogen, oxygen, and argon. These data have been obtained by a critical study of the recent literature, combined with new thermodynamic calculations for some of the species.

II. REACTION ENERGIES

The energy of a reaction may be defined as the amount of energy taken up when one molecule or one mole of reactants reacts to form products. Unless otherwise stated, all reactants and products are assumed to be in their ground rotational, vibrational, and electronic levels (i.e., no excitation). Such a reaction energy is equivalent to what chemists call "the heat of reaction at 0°K", since at absolute zero (in thermal equilibrium) all particles are in their lowest levels. Moreover, for gases at absolute zero no distinction needs to be made between the reaction enthalpy (or heat) and the reaction energy, since $\Delta H_0^{\circ} - \Delta E_0^{\circ} = \Delta(pV)_{T=0} = 0$, where by international convention the subscript "0" designates 0°K and the superscript "o" the ideal gas state. (A similar relation holds for the free energy: $\Delta F_0^{\circ} = \Delta A_0^{\circ} = \Delta E_0^{\circ}$.)

Chemists usually consider the reaction energy to vary with temperature, because of the varying thermal energy of the reactants and products. However, only the zero-temperature reaction energy will be tabulated here, for two reasons. In the first place, the variation with temperature is usually relatively small except at high temperatures, and can be readily estimated if necessary. Secondly, and more important in the present context, it is highly unlikely that the reaction products will be formed with an initial energy distribution corresponding to the ambient gas temperature. Hence, the temperature-dependent reaction energy is not really pertinent to the reaction-rate problem, but only to the question of the net heating after the products are thermalized by subsequent collisions.

The sign of the reaction energy depends upon whether it is defined as the energy absorbed or the energy released on reaction. For the cases considered here it is conventional to define it as the energy absorbed during formation, dissociation, or ionization, or equivalently the excess of the internal energy of the products over that of the reactants.

Before standard heats of formation can be determined, it is also necessary to establish "reference states" which are the standard substances from which all the other species are formed. For the species of present interest the conventional reference states are H₂, N₂, O₂, and Ar in the ideal-gas (or isolated-molecule) state, and C in the form of graphite.

Table 1 gives the molecular weight, energy of formation, dissociation energy, and ionization energy for the (gaseous) atoms and small molecules of present interest, together with the relevant references. All energies are given in three different units: physical (electron volts per particle), spectroscopic (reciprocal of the wavelength of photons with that energy, in 10³ cm⁻¹, also known as kilokaysers), and thermochemical (kilocalories per mole). Conversion factors are taken from the recent NAS-NRC list.⁽¹⁾ The accuracy of each value is indicated roughly by the number of decimal places shown. The energy of any other reaction involving these species may be readily calculated by adding the formation energies of the products and subtracting those of the reactants.

III. ENERGY LEVELS

The energy levels of atoms and atomic ions depend upon the arrangement of their orbital electrons. The lower energy levels of the atoms and ions of present interest are listed in Tables 2 to 11, together with their electronic state designation and statistical weight. (For an explanation of the latter terms see Herzberg⁽²⁾ or Moore⁽³⁾.) The ions H⁻, H⁺, and C⁺ have only one bound state, so they are not tabulated. All of the listed species except O⁻ actually have an infinite number of highly-excited states. The present tables, however, list only those states where all the bound electrons have principal quantum numbers less than 4. In most physical situations higher states will not play an important role, but if needed they can be obtained from the more extensive tabulations of Moore,⁽³⁾ or they may be calculated from the Rydberg formula.⁽²⁾

Also included in Tables 2 to 11 are the equilibrium fractional populations of the different electronic states, for various temperatures up to 10,000°K. These results are often useful in problems concerning equilibrium gases. The reader is cautioned, however, that in many situations involving low-density gases, or transient processes even at high densities, equilibrium will not obtain and these tabulated populations will not be applicable.

Molecules, because of their additional rotational and vibrational degrees of freedom, have so many individual energy levels that it is rather impractical to tabulate them. Fortunately, however, for each degree of electronic excitation the rotational and vibrational levels are usually quite regular and can be represented by simple formulas, only the coefficients of which need to be tabulated (see Herzberg⁽⁴⁾). For present purposes it is probably sufficient to note that the vibrational levels of each electronic state are fairly evenly spaced, only slowly converging near the dissociation limit. The rotational levels are not evenly spaced, but vary approximately quadratically with the rotational quantum number; however, the spacing is generally so close (10^{-3} to 10^{-2} eV) that for most reaction-rate purposes the rotational energy levels can be treated as if they formed a continuum.

Tables 12 to 18 present the electronic energy, lowest vibrational interval, and fractional population for the lower electronic states of several diatomic molecules of present interest. Similar values for other diatomic molecules, but without the fractional population numbers, are given in Table 19. (Again, the reader is cautioned against use of the equilibrium population values in nonequilibrium situations.)

For triatomic molecules, existing knowledge of the lower excited electronic states is quite incomplete. Consequently, only the ground state and its lowest vibrational intervals, for the three normal vibrational modes, are listed in Table 20.

IV. EQUILIBRIUM CONSTANTS

In any ideal-gas mixture in complete thermal and chemical equilibrium that contains three or more species which can be related by a possible reaction, such as $XY \rightleftharpoons X + Y$ or $W + X \rightleftharpoons Y + Z$, the concentration ratios $(X)(Y)/(XY)$, $(Y)(Z)/(W)(X)$, etc., can be shown to depend only on the temperature.⁽⁵⁾ Since these ratios are independent of the individual species concentrations, they are called equilibrium constants.

A reacting gas mixture that is initially out of chemical equilibrium will tend to approach equilibrium, and the ratios defined above will tend to approach their equilibrium values. As equilibrium is approached the various reaction rates do not actually become small, but instead each reaction becomes balanced by its reverse reaction. Consequently, it can be shown that in equilibrium the ratio of the forward to backward rate coefficients for each reaction is equal to its equilibrium constant. Unfortunately, a rate coefficient is measurable and has practical significance only in nonequilibrium situations. Among each reactant species there will always be particles with a range of velocities, and usually with a range of rotational, vibrational, or electronic levels. Generally some of these levels or velocities will be more reactive than others. In nonequilibrium situations the more reactive ones will be removed (by reaction) more rapidly than the others, resulting in an internal distribution of levels or velocities in each reacting species which makes it less reactive than if it had an equilibrium distribution.⁽⁶⁾

Molecular dissociation calculations based on a simple model show that when the mean thermal energy is much less than the dissociation energy the dissociation and association coefficients are only slightly smaller than their equilibrium values. Moreover, both coefficients are decreased by the same fraction, so that their ratio still equals the equilibrium constant.⁽⁷⁾ However, other types of reactions may not have such a convenient behavior.

In the present work, equilibrium constants have been calculated for several pertinent dissociation and ionization reactions. Results up to 10,000°K are presented graphically in Figs. 1 to 4.

Table 1. MOLECULAR WEIGHTS AND ENERGIES OF FORMATION, DISSOCIATION,
AND IONIZATION FOR SELECTED ATOMS AND MOLECULES

Species	Molec. Weight*	Energy (or Heat) of Formation**	Dissociation Energy**	Ionization Energy**	Reference ***
H^-	1.00852	1.~85 ev	-	0.754	8
		$11.974 \times 10^3 \text{ cm}^{-1}$		6.083	
		34.235 kcal/mole		17.392	
H	1.00797	2.239	-	13.598	3
		18.057		109.679	
		51.627		313.585	
H^+	1.00742	15.837	-	-	-
		127.736			
		365.213			
C^-	12.01170	6.24	-	1.13	9,10
		50.4		9.1	
		144.0		26.0	
C	12.01115	7.371	-	11.259	3
		59.452		90.814	
		169.979		259.648	
C^+	12.01060	18.630	-	24.382	3
		150.265		196.659	
		429.627		562.272	
N	14.0067	4.880	-	14.532	3
		39.359		117.214	
		112.532		335.129	
N^+	14.0062	19.412	-	29.601	3
		156.573		238.751	
		447.661		682.618	
O^-	15.9999	1.079	-	1.478	11
		8.705		11.925	
		24.89		34.10	
O	15.9994	2.558	-	13.18	3
		20.630		109.837	
		58.984		314.037	

Table 1 (Cont.)

Species	Molec. Weight*	Energy (or Heat) of Formation**	Dissociation Energy**	Ionization Energy**	References***
O ⁺	15.9989	16.175 ev 130.467 x10 ³ cm ⁻¹ 373.021 kcol/mole	-	35.117 283.244 809.829	3
Ar	39.948	0 0 0	-	15.759 127.110 363.423	3
Ar ⁺	39.947	15.759 127.110 363.423	-	27.629 222.848 637.149	3
H ₂	2.01594	0 0 0	4.477 36.114 103.254	15.425 124.414 355.715	13, 14
H ₂ ⁺	2.01539	15.425 124.414 355.715	2.651 21.379 61.125	-	14
CO	28.0106	-1.179 -9.513 -27.200	11.108 89.595 256.163	14.013 113.029 323.163	15 to 18
CO ⁺	28.0100	12.834 103.516 295.963	8.354 67.380 192.648	27.8 224 640	19
N ₂	28.0134	0 0 0	9.759 78.717 225.061	15.580 125.667 359.297	17, 20, 21
N ₂ ⁺	28.0129	15.580 125.667 359.297	8.711 70.264 200.893	27.1 219 626	22
NO ⁻	30.0066	0.6 5 14	5.3 43 123	0.3 2 7	†

Table 1 (Cont.)

Species	Molec. Weight*	Energy (or Heat) of Formation**	Dissociation Energy**	Ionization Energy**	References***
NO	30.0061	0.931 ev $7.506 \cdot 10^3 \text{ cm}^{-1}$ 21.46 kcal/mole	6.507 52.483 150.055	9.267 74.747 213.711	15, 23, 24
NO ⁺	30.0056	10.198 82.253 235.17	10.857 (N - O ⁺) 87.573 250.382	30.5 246 703	19
O ₂ ⁻	31.9993	-0.43 -3.5 -10.0	4.08 32.9 94.0	0.43 3.5 10.0	25
O ₂	31.9988	0 0 0	5.115 41.260 117.967	12.063 97.295 278.178	26, 27
O ₂ ⁺	31.9983	12.063 97.295 278.178	6.670 53.802 153.826	24.2 195 558	22
OH ⁻	17.0079	-1.43 -11.51 -32.9	4.75 (O ⁻ - H) 38.27 109.4	1.83 14.75 42.2	28
OH	17.0074	0.401 3.24 9.26	4.395 35.45 101.33	13.34 107.6 307.6	29 through 31
OH ⁺	17.0068	13.74 116.8 316.8	4.65 (O - H ⁺) 37.5 107.2	-	-
H ₂ O	18.0153	-2.476 -19.972 -57.103	5.116 (H - OH) 41.27 117.98	12.619 101.78 291.0	15, 32
H ₂ O ⁺	18.0148	10.143 81.81 233.9	5.84 (H - OH ⁺) 47.1 134.7	-	-
CO ₂	44.0100	-4.075 -32.865 -93.965	5.453 (CO - O) 43.982 125.750	13.769 111.06 317.5	15, 32

Table 1 (Cont.)

Species	Molec. Weight*	Energy (or Heat) of Formation**	Dissociation Energy**	Ionization Energy**	References***
CO_2^+	44.0094	9.694 ev $78.19 \times 10^3 \text{ cm}^{-1}$ 223.5 kcal/mole	5.179 ($\text{CO} - \text{O}^+$) 41.77 119.4	22.6 182 521	19
NO_2^-	46.0060	-3.6 -29 -83	5.6 ($\text{NO} - \text{O}^-$) 45 128	4.0 32 92	33
NO_2	46.0055	0.372 3.00 8.59	3.116 ($\text{NO} - \text{O}$) 25.13 71.86	9.78 78.9 225.6	15, 34
NO_2^+	46.0050	10.15 81.9 234.2	2.60 ($\text{NO}^+ - \text{O}$) 21.0 60.0	-	-
N_2O	44.0128	0.881 7.107 20.32	1.677 ($\text{N}_2 - \text{O}$) 13.523 38.66	12.894 104.00 297.35	15, 32
N_2O^+	44.0123	13.775 111.11 317.67	1.302 ($\text{N} - \text{NO}^+$) 10.50 30.02	-	-
O_3^-	47.9987	-0.4 -3 -9	1.5 ($\text{O}_2 - \text{O}^-$) 12 34	1.9 15 44	35
O_3	47.9982	1.506 12.15 34.74	1.051 ($\text{O}_2 - \text{O}$) 8.48 24.25	12.80 103.2 295.2	15, 36
O_3^+	47.9977	14.31 115.4 329.9	0.32 ($\text{O}_2^+ - \text{O}$) 2.5 7.3	-	-

* Molecular weights are for the normal isotopic mixture, based on $\text{C}^{12} = 12.00000$.

** All reaction energies are for isolated particles in their lowest rotational, vibrational, and electronic state (microscopic description), or for ideal gases at 0°K (equivalent macroscopic description). All energies are given in three units: physical (electron volts), spectroscopic (10^3 cm^{-1} or kilokaysers), and thermochemical (kilocalories per mole).

*** To avoid unnecessary duplication, references are indicated only where they give directly a formation, dissociation, or ionization energy. Where a dissociation energy is calculated from the formation energies of the molecule and its dissociation products, references to the latter are given only opposite the products. Similarly, no direct references are given for formation energies calculated from measured dissociation or ionization energies.

† Electron affinity estimated, since observation of negative charge transfer (37) to O_2^- shows that the magnetron measurement (33) is inaccurate.

Table 2. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF H

STATE (CP-1)	LEVEL (EV)	STAT- WT.	TEMPERATURE (DEG K)						
			200	250	300	400	500	600	
1s	C	C.CCCC	2	1.00E 00					
2s	82259	10.1986	2	C.	0.	0.	0.	0.	0.
2p	82259	10.1986	6	0.	0.	0.	0.	0.	0.
3s	97492	12.7482	2	0.	0.	0.	0.	0.	0.
3p	97492	12.7482	6	0.	0.	0.	0.	0.	0.
3d	97492	12.0872	1C	C.	0.	0.	0.	0.	0.
4s	1C2824	12.7482	2	0.	0.	0.	0.	0.	0.
4p	1C2824	12.7482	6	C.	0.	0.	0.	0.	0.
4d	102324	12.7482	1C	0.	0.	0.	0.	0.	0.
4f	102324	12.7482	14	0.	0.	0.	0.	0.	0.

ENERGY LEVELS FROM MOORE (3).
NOTE: STATES INVOLVING ELECTRONS WITH PRINCIPAL QUANTUM NUMBERS ABOVE N = 4 ARE NOT INCLUDED.

LEVEL (CP-1)	2000	2500	3000	3500	TEMPERATURE (DEG K)							
					4000	4500	5000	6000	7000	8000	9000	10000
C	1.CCE CC	1.COE CO	1.00E 00	1.00E 00	1.00E 00	1.00E 00	1.00E 00	1.00E 00	1.00E 00	1.00E 00	1.00E 00	1.00E 00
82259	1.99E-26	2.-5E-21	7.36E-18	2.06E-15	1.41E-13	3.78E-12	5.25E-11	2.71E-09	4.54E-08	3.76E-07	1.94E-06	7.24E-06
82259	5.98E-26	8.26E-21	2.21E-17	6.18E-15	4.24E-13	1.13E-11	1.57E-10	8.14E-09	1.36E-08	1.13E-06	5.83E-06	2.17E-05
97492	3.47E-31	4.29E-25	4.94E-21	3.93E-18	5.89E-16	2.9CE-14	6.55E-13	7.03E-11	1.98E-09	2.43E-08	1.70E-07	8.09E-07
97492	1.04E-30	1.29E-24	1.48E-20	1.18E-17	1.77E-15	8.7CE-14	1.97E-12	2.11E-10	5.95E-09	7.28E-08	5.11E-07	2.43E-06
97492	1.74E-3C	2.15E-24	2.47E-20	1.97E-17	2.95E-15	1.45E-13	3.28E-12	3.51E-10	9.92E-09	1.21E-07	8.52E-07	4.05E-06
1C2824	7.5CE-33	1.99E-26	3.83E-22	4.39E-19	8.66E-17	5.27E-15	1.41E-13	1.96E-11	6.63E-10	9.30E-09	7.26E-08	3.76E-07
1C2824	2.25E-32	5.98E-26	1.15E-21	1.32E-18	2.6CE-16	1.58E-14	4.24E-13	5.87E-11	1.59E-09	2.79E-08	2.18E-07	1.13E-06
1C2824	3.75E-32	9.97E-26	1.91E-21	2.20E-18	4.33E-16	2.64E-14	7.C6E-13	9.78E-11	3.31E-09	4.65E-08	3.63E-07	1.88E-06
1C2824	5.25E-32	1.40E-25	2.68E-21	3.07E-18	6.06E-16	3.69E-14	9.89E-13	1.37E-10	4.64E-09	6.51E-08	5.08E-07	2.63E-06

Table 3. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF C

STATE	(CP-1)	LEVEL (EV)	STAT.	WT.	2CC	250	300	400	5CC	600	8CC	1CCC	1500
2s ² 2p ² 3p ⁰	0	0.CC00	1	1.37E-01	1.28E-01	1.23E-01	1.19E-01	1.17E-01	1.16E-01	1.14E-01	1.14E-01	1.14E-01	1.14E-01
3p ¹	16	0.CC20	3	3.64E-01	3.58E-01	3.54E-01	3.49E-01	3.46E-01	3.44E-01	3.41E-01	3.40E-01	3.38E-01	3.38E-01
3p ¹	43	C.CC54	5	4.99E-01	5.11E-01	5.18E-01	5.28E-01	5.33E-01	5.37E-01	5.42E-01	5.44E-01	5.48E-01	5.48E-01
3D ⁰	10194	1.26E-3	5	9.67E-33	2.18E-26	5.74E-22	7.34E-17	1.10E-11	1.44E-11	2.47E-07	2.47E-07	2.47E-07	2.47E-07
3s ²	21648	2.6839	1	0.	0.	0.	0.	0.	0.	0.	0.	0.	1.10E-10
2s 2p ³ 3s ⁰	33735	4.1625	5	0.	0.	0.	0.	0.	0.	0.	0.	0.	5.06E-15
3p ⁰	64091	7.9461	15	0.	0.	0.	0.	0.	0.	0.	0.	0.	3.43E-27
3p ⁰	75256	9.3203	5	0.	0.	0.	0.	0.	0.	0.	0.	0.	4.60E-32
3p ⁰	105801	13.1173	3	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
3p ⁰	97878*	12.1350	5	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
2p ⁴	119878*	14.8626	3	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
15COC0*	16.5972	6	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
15ACCO*	19.5890	5	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
181CCC*	22.4406	1	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
2s ² 2p ³ (P ^o) 3s	6C776	7.5351	12	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
3p	39	69722	8.6442	36	0.	0.	0.	0.	0.	0.	0.	0.	3.72E-29
3p	78426	9.7233	66	0.	0.	0.	0.	0.	0.	0.	0.	0.	1.47E-32
4s	78184	9.6933	12	0.	0.	0.	0.	0.	0.	0.	0.	0.	3.70E-33
4p	80866	10.0258	36	0.	0.	0.	0.	0.	0.	0.	0.	0.	8.47E-34
4d	83850*	10.3958	6C	0.	0.	0.	0.	0.	0.	0.	0.	0.	8.06E-35
2s 2p ³ (P ^o) 3s	84CCCC*	1C.4144	84	0.	0.	0.	0.	0.	0.	0.	0.	0.	9.78E-35
3p	1160C0*	1C.3816	24	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
3p	125CC0*	15.4976	72	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
3p	132CC0*	16.3655	12C	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
4	136169*	16.8824	24	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
2s 2p ³ (D ^o) 3s	138CCCC*	17.1044	2L	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
3p	145CCC*	17.9772	60	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
3d	154CCC*	15.0931	1CC	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
4	158C00*	19.5890	32C	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
(P ^o) 3	184CCC*	22.8125	108	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
4	184CCC*	24.0523	192	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
(S ^o) 3	17COC0*	21.0768	3t	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
4	179CCC*	22.1926	6t	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
2p ³ (S ^o) 3	216CCC*	26.7799	72	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
4	225CC0*	27.8957	128	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
(D ^o) 3	224CCC*	27.7717	18C	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
4	213CC0*	28.8876	32C	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
(P ^o) 3	243CC0*	30.1274	1Ct	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.
4	252CCC*	31.2432	192	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.

ESTIMATE.

♦ INFLUENCE ESTIMATED SUBLVELS.

ACNSTARRED ENERGY LEVELS FROM MCCRT (38).

NOTE: STATES INVOLVING ELECTRONS WITH PRINCIPAL QUANTUM NUMBERS ABITIVE N = 4 ARE NOT INCLUDED.
ALL LEVELS AND CP-1 ARE SUBJECT TO AUTOMIZATION.

Table 3 (Cont.). ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF C

LEVEL (CM-1)	TEMPERATURE (K)						9CCO			10000	
	2CCO	25CO	35CO	35CC	4000	4500	5000	6CCO	7000	80CO	9CCO
0. 1.13E-01	1.13E-01	1.12E-01	1.12E-01	1.11E-01	1.10E-01	1.09E-01	1.07E-01	1.04E-01	1.02E-01	1.00E-01	9.80E-02
16 3.36E-01	3.35E-01	3.34E-01	3.32E-01	3.30E-01	3.28E-01	3.25E-01	3.19E-01	3.12E-01	3.06E-01	2.59E-01	2.43E-01
43 5.50E-01	5.50E-01	5.50E-01	5.48E-01	5.45E-01	5.41E-01	5.37E-01	5.28E-01	5.17E-01	5.07E-01	4.97E-01	4.87E-01
10194 3.71E-04	1.60E-03	4.22E-03	8.44E-03	1.41E-02	2.11E-02	2.89E-02	4.63E-02	6.42E-02	8.17E-02	9.81E-02	1.13E-01
21648 1.96E-08	4.38E-07	3.48E-06	1.52E-05	4.60E-05	1.08E-04	2.14E-04	5.94E-04	1.22E-03	2.08E-03	3.14E-03	4.35E-03
33735 1.64E-11	2.09E-09	5.28E-08	5.29E-07	2.97E-06	1.14E-05	3.31E-05	1.64E-04	5.99E-04	1.18E-03	2.28E-03	3.82E-03
64091 1.61E-20	1.62E-16	7.53E-14	6.04E-12	1.62E-10	2.08E-09	1.60E-08	3.38E-07	2.98E-06	1.51E-05	5.31E-05	1.45E-04
75256 3.14E-24	1.57E-19	2.14E-16	3.68E-14	1.75E-12	3.51E-11	3.85E-10	1.40E-08	1.80E-07	1.22E-06	5.37E-06	1.75E-05
1C5801 3.00E-34	1.22E-27	3.09E-23	4.32E-21	9.05E-18	6.71E-16	1.96E-14	3.07E-12	1.12E-10	1.67E-09	1.35E-08	7.20E-08
97878 1.49E-31	1.94E-25	2.30E-21	1.67E-19	2.34E-16	1.41E-14	3.19E-13	3.42E-11	5.56E-10	1.16E-08	8.01E-08	3.75E-07
119878 1.20E-38	3.69E-31	3.62E-26	1.32E-22	6.23E-20	7.44E-18	3.41E-16	1.05E-13	6.24E-12	1.33E-10	1.43E-09	9.50E-09
15C0CO 0.	3.27E-38	5.77E-32	1.67E-27	3.68E-24	1.47E-21	1.76E-19	2.29E-16	3.83E-14	1.77E-12	3.47E-11	3.74E-10
1580CC 0.	0.	6.92E-34	3.45E-29	1.15E-25	6.31E-23	9.77E-21	1.87E-17	4.11E-15	2.33E-13	5.36E-12	6.57E-11
1a10CC 0.	0.	2.24E-39	5.41E-34	5.88E-30	6.08E-27	2.61E-24	1.51E-20	7.27E-18	7.45E-16	2.71E-14	4.80E-13
60777 1.4CE-19	8.73E-16	2.95E-13	1.89E-11	4.26E-10	4.79E-09	3.32E-08	5.55E-07	4.71E-06	2.2CE-05	7.24E-05	1.87E-04
69722 6.73E-22	1.52E-17	1.21E-14	1.43E-12	5.12E-11	8.23E-10	7.58E-09	2.1CE-07	2.25E-06	1.32E-05	5.26E-05	1.55E-04
78426 2.14E-24	1.69E-19	3.11E-16	6.67E-14	3.72E-12	8.49E-11	1.03E-09	4.35E-08	6.25E-07	4.59E-06	2.15E-05	7.39E-05
78184 5.09E-25	3.89E-20	6.99E-17	1.47E-14	8.13E-13	1.83E-11	2.21E-10	9.22E-09	1.31E-07	9.59E-07	4.48E-06	1.53E-05
80866 2.22E-25	2.49E-20	5.80E-17	1.47E-14	9.29E-13	2.33E-11	3.07E-10	1.45E-08	2.27E-07	1.78E-06	8.75E-06	3.12E-05
83850 4.32E-26	7.46E-21	2.31E-17	7.17E-15	5.29E-13	1.50E-11	2.17E-10	1.18E-08	2.05E-07	1.73E-06	9.05E-06	3.39E-05
84CCC 5.43E-26	9.58E-21	3.01E-17	9.44E-15	7.02E-13	2.00E-11	2.91E-10	1.60E-08	2.78E-07	2.36E-06	1.24E-05	4.64E-05
1160CC 1.56E-36	2.75E-29	1.86E-24	5.22E-21	2.01E-18	2.06E-16	8.32E-15	2.13E-12	1.11E-10	2.13E-09	2.12E-08	1.33E-07
125000 0.	4.64E-31	7.44E-26	3.88E-22	2.37E-19	1.87E-15	7.37E-15	5.22E-13	5.22E-11	2.27E-C9	1.51E-10	1.09E-07
1320CC 0.	1.38E-32	4.32E-27	3.63E-23	3.18E-20	6.17E-18	4.16E-16	2.29E-13	2.06E-11	6.00E-10	8.22E-09	6.64E-08
136169 0.	2.50E-34	1.17E-24	1.31E-21	1.42E-19	3.26E-19	2.51E-17	1.69E-14	1.75E-12	5.67E-11	8.44E-10	7.29E-09
1380CC 0.	7.27E-35	4.05E-29	5.14E-25	6.13E-22	1.51E-19	1.23E-17	9.06E-15	1.00E-12	3.40E-11	5.25E-10	4.67E-09
1450CC 0.	3.88E-36	4.23E-30	8.68E-26	1.48E-22	4.83E-20	4.94E-19	5.07E-15	7.14E-13	2.90E-11	5.14E-10	5.12E-09
154000 0.	3.64E-38	9.42E-32	3.58E-27	9.71E-24	4.93E-21	6.18E-19	9.77E-16	1.87E-13	9.57E-12	2.33E-10	2.34E-09
1580CC 0.	0.	4.43E-32	2.21E-27	7.37E-24	4.0CE-21	6.25E-19	1.20E-15	2.63E-13	1.49E-11	3.43E-10	4.20E-09
1840CC 0.	0.	5.74E-38	1.70E-28	3.34E-25	2.16E-21	1.19E-22	7.93E-19	4.24E-16	4.69E-14	1.81E-12	3.37E-11
184000 0.	0.	0.	4.96E-34	1.05E-29	2.43E-26	1.19E-23	1.28E-19	5.65E-17	1.38E-14	6.52E-13	1.42E-11
17C0CC 0.	1.58E-35	1.79E-30	1.11E-26	9.80E-24	2.23E-21	7.58E-18	2.51E-15	1.94E-13	5.67E-12	8.41E-11	
1750CC 0.	3.74E-37	7.08E-32	7.73E-23	9.8CE-25	2.97E-22	1.56E-18	7.02E-16	6.03E-14	2.39E-12	4.10E-11	
2160CC 0.	0.	2.20E-38	1.44E-33	8.03E-30	7.94E-27	2.46E-22	3.93E-19	5.9CE-17	7.26E-15	2.25E-13	
2250CC 0.	0.	0.	1.01E-34	8.03E-31	1.06E-27	5.05E-23	1.10CE-19	3.49E-17	3.0CE-15	1.09E-13	
2240CC 0.	0.	0.	2.03E-34	1.56E-30	1.99E-27	9.02E-23	1.90E-19	5.87E-17	5.05E-15	1.78E-13	
233000C 0.	1.42E-35	1.56E-31	2.65E-28	1.05E-23	1.05E-20	2.07E-17	2.07E-17	2.07E-17	2.13E-15	6.65E-14	
243000C 0.	1.31E-37	2.15E-33	5.03E-30	5.68E-25	2.29E-21	1.16E-18	1.45E-16	6.93E-15	1.64E-16	6.93E-15	
252000 0.	0.	0.	2.15E-34	6.71E-31	1.17E-22	6.41E-22	6.41E-22	4.07E-19	6.13E-17	3.37E-15	

Table 4. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF C⁺

STATE	LF LEVEL (CM-1)	LF LEVEL (EV)	STAT. WT.	TEMPERATURE (DEG K)						ICCO	ICCO	
				200	300	400	500	600	800			
2s ² 2p ¹ P _{1/2}	0	0.0000	2	4.42E-01	4.22E-01	4.05E-01	3.75E-01	3.47E-01	3.19E-01	3.04E-01	3.04E-01	
2s ² 2p ³ P _{1/2}	64	0.0079	4	5.58E-01	5.8CE-01	5.95E-01	6.14E-01	6.25E-01	6.41E-01	6.46E-01	6.53E-01	
2s ² 2p ³ P _{3/2}	43033	5.3353	12	0.	0.	0.	0.	0.	0.	0.	0.	
2s ² 2D ³ D	74932	9.2901	10	0.	0.	0.	0.	0.	0.	0.	0.	
2s ² 2s ²	96494	11.9634	2	0.	0.	0.	0.	0.	0.	0.	0.	
2p ¹ 1P _{1/2}	110653	13.7189	6	0.	0.	0.	0.	0.	0.	0.	0.	
2p ¹ 1S _{1/2}	142024	17.0083	4	0.	0.	0.	0.	0.	0.	0.	0.	
2p ¹ 1D _{3/2}	150465	18.6548	1C	0.	0.	0.	0.	0.	0.	0.	0.	
2p ¹ 1D _{5/2}	168744	2C.9210	6	0.	0.	0.	0.	0.	0.	0.	0.	
2s ² 1S _{1/2}	116538	14.4485	2	0.	0.	0.	0.	0.	0.	0.	0.	
3p	131725	16.3314	6	0.	0.	0.	0.	0.	0.	0.	0.	
3d	145550	18.0454	1C	0.	0.	0.	0.	0.	0.	0.	0.	
4s	157235	19.4942	2	0.	0.	0.	0.	0.	0.	0.	0.	
4p	162523	20.1498	6	0.	0.	0.	0.	0.	0.	0.	0.	
4d	168125	20.8443	10	0.	0.	0.	0.	0.	0.	0.	0.	
2s ² 2p ¹ P _{1/2}) _{3s}	4f	168979	20.9502	14	0.	0.	0.	0.	0.	0.	0.	0.
3p	170643	21.1565	18	0.	0.	0.	0.	0.	0.	0.	0.	
3d	184786	22.9100	54	0.	0.	0.	0.	0.	0.	0.	0.	
3d	197742	24.5163	9C	0.	0.	0.	0.	0.	0.	0.	0.	
4s	21CCC0*	26.0360	18	0.	0.	0.	0.	0.	0.	0.	0.	
4p	215730**	26.7464	54	0.	0.	0.	0.	0.	0.	0.	0.	
4d	220465	27.3335	9C	0.	0.	0.	0.	0.	0.	0.	0.	
4f	221458	27.4566	126	0.	0.	0.	0.	0.	0.	0.	0.	
(1P ⁰) _{3s}	219CC0*	27.1518	6	0.	0.	0.	0.	0.	0.	0.	0.	
3p	2340C0*	29.0116	18	0.	0.	0.	0.	0.	0.	0.	0.	
3d	246CC0*	30.4993	3C	0.	0.	0.	0.	0.	0.	0.	0.	
4s	26CCCC*	32.2351	6	0.	0.	0.	0.	0.	0.	0.	0.	
4p	265CC0*	32.8550	18	0.	0.	0.	0.	0.	0.	0.	0.	
4d	27C5CC*	33.5369	3C	0.	0.	0.	0.	0.	0.	0.	0.	
4f	27140C*	33.6484	42	0.	0.	0.	0.	0.	0.	0.	0.	
?p ¹ P _{1/2}) _{3s}	257050*	31.8631	18	0.	0.	0.	0.	0.	0.	0.	0.	
3p	271CC0*	33.5989	54	0.	0.	0.	0.	0.	0.	0.	0.	
3d	283CC0*	35.0866	9C	0.	0.	0.	0.	0.	0.	0.	0.	
4	3045C0*	37.7522	268	0.	0.	0.	0.	0.	0.	0.	0.	
(1D) _{3s}	2670C0*	33.1029	1C	0.	0.	0.	0.	0.	0.	0.	0.	
3p	281CC0*	34.8387	3C	0.	0.	0.	0.	0.	0.	0.	0.	
3d	292CC0*	36.2025	5C	0.	0.	0.	0.	0.	0.	0.	0.	
4	313CCC*	38.8061	16C	0.	0.	0.	0.	0.	0.	0.	0.	
(1S) _{3s}	315CCC*	39.5499	18	0.	0.	0.	0.	0.	0.	0.	0.	
4	3490CC*	43.2694	32	0.	0.	0.	0.	0.	0.	0.	0.	

*ESTIMATED.

**INCLUDES ESTIMATED SUBLVELS.

ACNSTARREC ENERGY LEVELS FROM MCCRE (39) AND GLAC (39).
 NOTE: STATES INVOLVING ELECTRONS WITH PRINCIPAL QUANTUM NUMBERS ABOVE N = 4 ARE NOT INCLUDED.
 ALL LEVELS ABOVE 196659 CP-1 ARE SUBJECT TO AUTICNIZATION.

Table 4 (Cont.). ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF C⁺

LEVEL (CP-1)	TEMPERATURE (ICCG K)										CCCC ICCC
	2CCC	2CC0	3CC0	35CC	4CCC	4CC0	5CC0	6000	7000	8C00	
C	3.44E-01	3.42E-01	3.4CE-01	3.39E-C1	3.38E-01	3.37E-01	3.36E-01	3.35E-01	3.34E-01	3.34E-01	6.56E-01
6.	6.56E-01	6.58E-01	6.60E-01	6.61E-C1	6.62E-01	6.62E-01	6.63E-01	6.64E-01	6.64E-01	6.64E-01	6.58E-01
43033	7.40E-14	3.60E-11	2.22E-09	4.23E-CR	3.05E-C7	2.15E-C6	8.48E-06	6.67E-05	2.91E-04	8.77E-04	2.07E-03
74932	6.67E-24	3.19E-19	4.20E-16	7.11E-14	3.33E-12	6.65E-11	7.29E-10	2.65E-08	3.44E-C7	2.36E-06	1.05E-05
96694	2.45E-31	2.6CF-25	2.71E-21	2.01E-18	2.86E-16	1.35E-14	2.95E-13	3.01E-11	8.18E-10	9.75E-C9	6.69E-08
1116553	2.77E-35	2.26E-28	9.15E-24	1.79E-20	5.26E-18	4.37E-16	1.50E-14	3.02E-12	1.34E-10	2.29E-09	2.09E-C8
42024	0.	2.17E-36	1.78E-30	2.97E-26	4.41E-23	1.28E-20	1.2CE-18	1.CSE-15	1.41E-13	5.42E-12	9.23E-11
4.21E-38	7.78E-32	2.33E-27	2.39E-24	2.65E-21	3.60E-19	6.23E-16	2.97E-14	2.97E-12	5.93E-11	6.62E-10	6.62E-10
68744	0.	7.27E-36	7.61E-31	4.43E-27	3.75E-24	8.26E-22	2.70E-18	6.72E-16	6.65E-14	1.93E-12	2.06E-11
1.34E-37	2.54E-30	1.81E-25	5.31E-22	2.11E-19	2.22E-17	9.21E-16	2.46E-13	1.33E-11	2.65E-10	2.72E-C5	1.74E-08
131725	0.	1.22E-33	3.74E-28	3.09E-24	2.65E-21	5.19E-19	3.49E-17	1.93E-14	1.76E-12	5.18E-11	7.19E-10
1.13E-37	8.22E-31	1.75E-26	3.10E-23	1.04E-20	1.09E-18	1.17E-15	1.71E-13	7.18E-12	1.31E-10	5.34E-09	1.45550
1.57235	C.	6.05E-34	2.88E-29	9.27E-26	4.96E-23	7.56E-21	1.42E-17	3.10E-15	1.76E-13	4.06E-12	5.00E-11
62523	0.	1.44E-34	9.82E-30	4.15E-26	2.74E-23	4.95E-21	1.20E-17	2.13E-15	2.04E-13	5.23E-12	7.1CE-11
1.68125	0.	1.63E-35	1.64E-30	9.22E-27	7.63E-24	1.65E-21	5.21E-18	1.65E-15	1.24E-13	3.56E-12	5.22E-11
686979	C.	1.52E-35	1.61E-30	9.5CE-27	8.13E-24	1.80E-21	5.95E-18	1.94E-15	1.49E-13	4.35E-12	6.46E-11
8.70E-36	1.05E-3C	6.71E-27	6.14E-24	1.44E-21	5.13E-18	1.77E-15	1.42E-13	4.28E-12	6.53E-11	6.53E-11	6.53E-11
2.98E-38	9.37E-33	1.24E-28	2.CCE-25	7.35E-23	5.18E-19	3.10E-16	3.34E-14	1.34E-12	2.56E-11	2.56E-11	2.56E-11
1.9742	C.	7.60E-35	1.96E-3C	5.30E-27	2.95E-24	3.86E-20	3.37E-17	5.42E-15	2.81E-13	6.62E-12	7.1CE-11
211000	0.	9.84E-38	4.77E-33	2.1CE-29	1.73E-26	4.C9E-22	5.43E-19	1.20E-16	7.93E-15	2.27E-13	2.27E-13
0.	0.	2.80E-38	1.82E-33	1.C1E-29	9.59E-27	3.10E-22	5.C2E-19	1.28E-16	9.52E-15	4.35E-12	6.46E-11
0.	0.	5.53E-34	3.70E-3C	4.26E-27	1.66E-22	3.16E-19	9.10E-17	7.44E-15	4.28E-12	6.53E-11	6.53E-11
0.	0.	5.42E-34	3.78E-30	4.48E-27	1.83E-22	3.61E-19	1.07E-16	8.89E-15	3.34E-12	5.67E-11	5.67E-11
0.	0.	6.25E-35	3.55E-31	4.33E-23	2.57E-23	2.85E-20	2.90E-18	6.27E-16	2.C7E-15	3.67E-13	5.67E-13
0.	0.	8.5CE-37	9.78E-33	1.73E-29	1.29E-24	3.91E-21	1.60E-18	1.71E-16	7.18E-15	2.27E-13	2.27E-13
0.	0.	1.89E-38	3.52E-34	9.15E-31	1.21E-25	5.54E-22	3.07E-19	4.11E-17	2.13E-15	2.99E-13	2.99E-13
0.	0.	0.	8.CCE-37	3.26E-33	8.46E-28	6.23E-24	4.95E-21	8.92E-19	5.68E-17	6.53E-16	6.53E-16
0.	0.	4.85E-37	2.32E-33	7.65E-28	6.69E-24	6.05E-21	1.20E-18	6.30E-17	6.30E-17	6.30E-17	6.30E-17
0.	0.	1.39E-37	7.93E-34	3.41E-28	3.6CE-24	3.75E-21	8.33E-19	6.27E-17	6.27E-17	6.27E-17	6.27E-17
0.	0.	1.46E-37	8.57E-34	3.85E-28	4.19E-24	4.46E-21	1.01E-18	7.11E-17	7.11E-17	7.11E-17	7.11E-17
0.	0.	0.	6.26E-36	2.32E-32	5.21E-27	3.4CE-23	2.55E-20	4.33E-18	2.62E-16	1.55E-17	1.38E-17
0.	0.	0.	2.14E-37	1.24E-33	5.44E-28	4.85E-24	6.17E-21	1.38E-18	1.05E-16	4.46E-20	4.74E-18
0.	0.	0.	0.	6.52E-35	5.11E-29	8.27E-25	1.11E-21	3.39E-19	3.11E-17	3.98E-19	3.51E-19
0.	0.	0.	4.29E-37	9.42E-31	3.19E-26	7.95E-23	3.47E-20	4.52E-18	3.66E-22	3.14E-22	3.14E-22
0.	0.	0.	1.42E-37	7.24E-34	2.63E-28	2.46E-24	2.34E-21	4.86E-19	3.46E-17	3.46E-17	3.46E-17
0.	0.	0.	0.	3.87E-35	2.75E-29	4.16E-25	5.67E-22	1.55E-19	1.55E-17	1.55E-17	1.55E-17
0.	0.	0.	0.	2.72E-36	3.28E-30	7.23E-26	1.31E-22	4.46E-20	4.46E-18	4.46E-18	4.46E-18
0.	0.	0.	0.	0.	6.82E-32	3.C9E-27	9.39E-25	1.1CE-21	3.98E-24	4.98E-22	4.98E-22
0.	0.	0.	0.	0.	1.82E-33	1.C9E-28	3.1CE-26	5.66E-25	3.66E-25	3.66E-25	3.66E-25
0.	0.	0.	0.	2.43E-36	3.78E-31	9.78E-29	1.01E-22	4.01E-20	4.01E-18	4.01E-18	4.01E-18
0.	0.	0.	0.	0.	81000	0.	0.	0.	0.	0.	0.
0.	0.	0.	0.	0.	22000	0.	0.	0.	0.	0.	0.
0.	0.	0.	0.	0.	183000	0.	0.	0.	0.	0.	0.
0.	0.	0.	0.	0.	185000	0.	0.	0.	0.	0.	0.
0.	0.	0.	0.	0.	167000	0.	0.	0.	0.	0.	0.
0.	0.	0.	0.	0.	1649000	0.	0.	0.	0.	0.	0.

Table 5. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF N

STATE (CP-1)	LEVEL (EV)	STAT. WT.	TEMPERATURE (DEG K)						ICCO	1500
			200	250	300	400	500	600		
2 ¹ 2p 3 ⁰ 3D [*]	192220	C	0.0000	4	1.00E 00					
2 ¹ p 3 ⁰	288339	2.3039	1C	0	0	0	0	2.30E-30	2.30E-24	2.42E-15
2 ¹ 2p 3 ¹ P	88132	3.5755	6	0	0	0	0	0	0	2.44E-12
2 ¹ D	1210CC0*	10.9267	12	0	0	0	0	1.37E-36	1.39E-30	1.43E-18
2 ¹ S	14211C*	17.6169	2	0	0	0	0	0	0	5.80E-37
2 ¹ P	1582C0*	19.6138	6	0	0	0	0	0	0	0
2 ¹ D	2329C0*	28.8752	6	0	0	0	0	0	0	0
2 ¹ S 2 ¹ p (3P) 3P	8428R	1C.4501	18	0	0	0	0	0	0	0
	95780	11.8749	54	0	0	0	0	0	0	0
3d	104861	13.0008	9C	0	0	0	0	0	0	0
4s	103861	12.8768	18	0	0	0	0	0	0	0
4p	107420**	13.3180	54	0	0	0	0	0	0	0
4d	11C315	13.6770	90	0	0	0	0	0	0	0
4f	110441	13.6926	126	0	0	0	0	0	0	0
(1D) 3s	99664	12.3564	1C	0	0	0	0	0	0	0
3p	11C973	13.7585	3C	0	0	0	0	0	0	0
3d	121CC0**	15.0017	5C	0	0	0	0	0	0	0
4s	1246C0**	15.4480	16C	0	0	0	0	0	0	0
(1S) 3s	116279	14.4164	2	0	0	0	0	0	0	0
3p	1284C0*	15.9192	6	0	0	0	0	0	0	0
3d	1375CC*	17.0474	1C	0	0	0	0	0	0	0
4s	142CCC*	17.6053	32	0	0	0	0	0	0	0
2 ¹ 2p 3 ¹ S	153CC0*	18.9691	9C	0	0	0	0	0	0	0
4	156CC0*	19.3410	15C	0	0	0	0	0	0	0
(1D*) 3	192LC0*	23.8044	27C	0	0	0	0	0	0	0
4	2C2CCC*	25.0442	48C	0	0	0	0	0	0	0
(1p) 3-4	215CC0*	26.6559	45C	0	0	0	0	0	0	0
(1S'U, 1P)	3-4	32.1111	55C	0	0	0	0	0	0	0
2 ¹ p (1D, 1S) 3-4	331CCC*	41.0377	75C	0	0	0	0	0	0	0

*ESTIMATED.

**INCLUDES FESTIVALE SUMLEVELS.

NONSTARRED ENERGY LEVELS FROM MCCREY (40) AND ERIKSSON (43).

NON-STATES INVOLVING ELECTRONS WITH PRINCIPAL QUANTUM NUMBERS ABOVE N = 4 ARE NOT INCLUDED.

ALL LEVELS ABOVE 117214 CP-1 ARE SUBJECT TO ALTCINIZATION.

Table 5 (Cont.). ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF N

LEVEL (E-V)	7CC0	2500	3CC0	35C0	40C0	45C0	50C0	6C00	7000	8000	9C00	1CC00
C	1.00E+00	1.00E+00	9.99E-01	9.97E-01	9.95E-01	9.93E-01	9.90E-01	9.74E-01	5.51E-01	9.20E-01	8.85E-01	8.47E-01
10220	2.46E-06	3.91E-05	2.47E-04	9.22E-C4	2.47E-C4	5.32E-03	9.78E-03	2.42E-02	4.27E-02	1.02E-01	1.33E-01	1.33E-01
28839	1.47E-09	9.29E-08	1.48E-06	1.06E-C5	4.68E-C5	1.48E-04	3.69E-04	1.45E-03	3.80E-03	7.71E-C3	1.22E-02	2.00E-02
88132	8.75E-28	2.81E-22	1.32E-18	5.53E-16	5.11E-14	1.73E-12	2.88E-11	1.94E-09	3.87E-08	3.61E-07	2.02E-06	7.90E-06
1210CC	3.93E-38	1.43E-30	1.57E-25	6.24E-22	3.13E-19	3.92E-17	1.87E-15	6.10E-13	3.76E-11	8.14E-10	8.79E-09	5.84E-08
142110	0.	1.51E-36	1.26E-34	2.13E-32	3.11E-29	9.2CE-23	9.2CE-21	8.61E-19	7.73E-16	9.8CE-14	3.65E-12	6.01E-11
158200	0.	1.68E-33	8.55E-30	2.9CE-29	2.9CE-25	1.61E-22	2.52E-20	6.69E-17	1.0CE-14	6.07E-13	1.38E-11	1.62E-10
232900	0.	0.	0.	6.20E-37	6.82E-33	1.16E-29	8.13E-25	2.31E-21	8.88E-19	8.97E-17	3.56E-15	
84288	2.09E-26	3.85E-21	1.25E-17	4.03E-15	3.08E-13	8.05E-12	1.30E-10	7.31E-09	1.28E-07	1.08E-06	5.60E-06	2.06E-05
5578C	1.61E-29	1.55E-23	1.52E-19	1.07E-16	1.47E-14	6.73E-13	1.43E-11	1.39E-09	3.62E-08	4.10E-07	2.57E-06	1.18E-05
104861	3.9CE-32	1.39E-25	3.24E-21	4.27E-18	9.34E-16	6.15E-14	1.75E-12	2.63E-10	5.33E-09	1.34E-07	1.04E-06	5.34E-06
1C3861	1.60E-32	4.94E-26	1.05E-21	1.29E-18	2.68E-16	1.69E-14	4.67E-13	6.69E-11	2.29E-09	3.25E-09	2.45E-07	1.23E-06
107420	3.71E-33	1.91E-26	5.70E-22	8.96E-19	2.23E-16	1.63E-14	5.03E-13	8.55E-11	3.31E-09	5.06E-08	4.16E-07	2.22E-06
11C315	7.70E-34	6.02E-27	2.37E-22	4.54E-19	1.31E-16	1.08E-14	3.8E-13	7.12E-11	3.04E-09	5.01E-08	4.36E-07	2.44E-06
110641	9.85E-34	7.84E-27	3.13E-22	6.04E-19	1.76E-16	1.45E-14	4.92E-13	9.67E-11	4.15E-09	6.85E-08	5.99E-07	3.35E-06
596664	1.02E-31	3.07E-25	6.36E-21	6.02E-18	6.73E-16	3.60E-14	8.60E-13	1.02E-10	3.02E-09	3.4E-08	2.66E-07	1.25E-06
110973	1.60E-34	1.37E-27	5.77E-23	3.15E-19	3.45E-17	7.91E-15	1.00E-13	2.32E-11	6.35E-10	1.31E-08	1.31E-07	7.35E-07
1210CC	1.96E-37	7.14E-30	7.84E-25	3.12E-21	1.56E-18	1.96E-16	9.35E-15	3.05E-12	1.88E-10	4.07E-09	4.39E-08	2.91E-07
124600	0.	2.88E-30	4.46E-25	2.27E-21	1.37E-18	1.59E-16	1.06E-14	4.12E-12	2.87E-10	6.82E-09	7.91E-08	5.55E-07
116279	2.34E-37	4.32E-30	3.02E-25	8.69E-22	3.42E-19	3.55E-17	1.46E-15	3.78E-13	1.98E-11	3.81E-10	3.74E-09	2.3CE-08
1284CC	0.	1.21E-32	2.70E-27	1.79E-23	1.31E-20	2.21E	1.33E-16	6.21E-14	4.92E-12	1.29E-11	1.61E-09	1.20E-08
137500	0.	1.07E-34	5.73E-29	7.07E-25	8.27E-22	2.01E-19	1.02E-17	1.17E-15	1.27E-14	1.26E-12	4.19E-11	6.28E-10
1420CC	0.	2.58E-35	2.12E-29	3.56E-25	5.24E-22	1.52E-19	1.42E-17	1.27E-15	1.60E-13	5.96E-11	9.79E-10	9.08E-09
1530CC	0.	1.29E-37	3.05E-31	1.09E-26	2.82E-23	1.27E-20	1.69E-18	2.55E-15	4.70E-13	2.32E-11	4.75E-10	5.24E-09
1550CC	0.	1.29E-31	5.63E-27	1.70E-23	8.67E-21	1.26E-18	2.21E-15	4.51E-13	2.40E-11	5.22E-10	6.05E-09	
1520CC	0.	0.	0.	3.56E-33	6.84E-29	1.47E-25	6.77E-23	6.65E-19	4.66E-16	6.26E-14	2.79E-12	5.75E-11
2C20CC	0.	0.	C.	1.04E-34	3.33E-30	1.07E-26	1.07E-19	1.06E-16	1.06E-14	1.06E-12	2.00E-11	2.42E-10
2150CC	0.	0.	0.	4.64E-37	2.91E-32	1.56E-28	1.51E-25	4.46E-21	6.87E-18	1.67E-15	1.16E-13	3.50E-12
2550CC	0.	0.	0.	0.	0.	1.49E-34	5.04E-31	1.43E-25	9.92E-22	7.45E-19	1.27E-16	7.62E-15
3310CC	0.	0.	0.	0.	0.	0.	C.	6.17E-33	5.06E-28	2.42E-24	1.73E-21	3.30E-19

Table 6. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF N.

STAT.	(CP-1) LEVEL (EV)	STAT. WT.	TEMPERATURE (LOG K)									
			2CC	25C	3CC	4CC	50C	600	800	1000	1500	2000
2s ² 2p ⁸ 2p ⁰	0	0.0000	1	1.77E-01	1.76E-01	1.51E-01	1.31E-01	1.26E-01	1.21E-01	1.20E-01	1.20E-01	1.20E-01
	4p ⁻⁷	C.C060	3	4.17E-C1	4.03E-01	3.93E-01	3.79E-01	3.70E-01	4.07E-01	4.87E-01	5.13E-01	5.22E-01
	130 ⁻⁸	0.0162	2	3.65E-C1	4.11E-01	4.42E-01	4.70E-01	4.87E-01	5.13E-01	5.22E-01	5.33E-01	5.44E-01
	15316	1.898 ₇	1	0.	4.65E-39	1.04E-32	6.93E-25	5.14E-20	7.65E-17	7.07E-13	1.69E-10	2.52E-01
	32389	4.0528	1	0.	0.	0.	0.	0.	0.	0.	0.	2.92E-11
2s ² 2p ⁸ 2p ⁰	46785	5.0005	5	C.	0.	0.	0.	0.	0.	0.	0.	0.
	92245	11.4366	15	C.	0.	0.	0.	0.	0.	0.	0.	0.
	1C9218	13.5410	9	C.	0.	0.	0.	0.	0.	0.	0.	0.
	155127	19.2328	3	C.	0.	0.	0.	0.	0.	0.	0.	0.
	1441e8	17.8766	5	C.	0.	0.	0.	0.	0.	0.	0.	0.
	166766	20.6758	3	C.	0.	0.	0.	0.	0.	0.	0.	0.
	2180CC0*	27.0279	9	C.	0.	0.	0.	0.	0.	0.	0.	0.
	2290CC0*	28.3916	5	C.	0.	0.	0.	0.	0.	0.	0.	0.
	2640CC0*	32.7310	1	C.	0.	0.	0.	0.	0.	0.	0.	0.
	149056	16.4801	12	C.	0.	0.	0.	0.	0.	0.	0.	0.
	3C	169022	20.9555	36	C.	0.	0.	0.	0.	0.	0.	0.
	3d	187693	23.2704	6C	C.	0.	0.	0.	0.	0.	0.	0.
	4C	196955	24.4187	12	C.	0.	0.	0.	0.	0.	0.	0.
	4P	2C3384	25.2158	36	C.	0.	0.	0.	0.	0.	0.	0.
	4G	210286	26.0712	6C	C.	0.	0.	0.	0.	0.	0.	0.
2s ² 2p ⁸ (1P) 3s	4f	211271	26.1936	84	C.	0.	0.	0.	0.	0.	0.	0.
	3p	207974	25.7048	24	C.	0.	0.	0.	0.	0.	0.	0.
	3p	220455	28.0761	72	C.	0.	0.	0.	0.	0.	0.	0.
	3p	2445C0	30.3134	12C	C.	0.	0.	0.	0.	0.	0.	0.
	4	2661C0	32.9913	384	C.	0.	0.	0.	0.	0.	0.	0.
	2s ² 2p ⁸ (1D) 3s	252CCC	31.2412	2C	C.	0.	0.	0.	0.	0.	0.	0.
	3p	27C0CC	33.4749	5C	C.	0.	0.	0.	0.	0.	0.	0.
	3p	298CCC	35.7065	1CC	C.	0.	0.	0.	0.	0.	0.	0.
	4	309800	38.409?	32C	C.	0.	0.	0.	0.	0.	0.	0.
	(1P) 3	3231C0	40.0583	1C8	C.	0.	0.	0.	0.	0.	0.	0.
	4	3548CC	43.9885	192	C.	0.	0.	0.	0.	0.	0.	0.
	(2S) 3	3081C0	38.1985	3C	C.	0.	0.	0.	0.	0.	0.	0.
	4	3358CC	42.1287	64	C.	0.	0.	0.	0.	0.	0.	0.
	4	364CCC	45.1296	72	C.	0.	0.	0.	0.	0.	0.	0.
	2P ⁸ (1S) 3	396CCC	49.0964	12C	C.	0.	0.	0.	0.	0.	0.	0.
	4	3ACCCC	47.1124	1AC	C.	0.	0.	0.	0.	0.	0.	0.
	3	4120CC	51.502	32C	C.	0.	0.	0.	0.	0.	0.	0.
	(8P) 3	4C8CCC	50.5042	1C8	C.	0.	0.	0.	0.	0.	0.	0.
	4	439CCC	54.4277	192	C.	0.	0.	0.	0.	0.	0.	0.

•ESTIMATED.
ENERGY LEVELS FROM FIGURE (3) AND ERIKSSON (42).
NOTE: STATES INVOLVING ELECTRONS WITH PRINCIPAL QUANTUM NUMBERS ABOVE $n = 4$ ARE NOT INCLUDED.
ALL LEVELS ABOVE 238751 CM $^{-1}$ ARE SUBJECT TO AUTIONIZATION.

Table 6 (Cont.) ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF N⁺

LEVEL (CP-1)	2CCO	2SCO	3CCO	3SCC	4CCO	4SCCO	5CCO	6CCO	7CCO	8CCO	9CCO	10CCO
C 1.18E-01	1.17E-01	1.16E-01	1.15E-01	1.14E-01	1.14E-01	1.13E-01	1.12E-01	1.10E-01	1.09E-01	1.07E-01	1.06E-01	1.05E-01
48 3.43E-01	3.41E-01	3.40E-01	3.38E-01	3.37E-01	3.36E-01	3.35E-01	3.32E-01	3.28E-01	3.24E-01	3.20E-01	3.15E-01	3.15E-01
13C 5.39E-01	5.42E-01	5.44E-01	5.45E-01	5.46E-01	5.46E-01	5.46E-01	5.42E-01	5.38E-01	5.32E-01	5.26E-01	5.19E-01	5.19E-01
15316 9.71E-C6	8.68E-05	8.74E-04	1.06E-03	2.32E-03	4.25E-03	6.90E-03	1.42E-02	2.37E-02	3.47E-02	4.64E-02	5.84E-02	5.84E-02
32689 7.25E-12	7.90E-10	1.80E-C8	1.68E-07	8.96E-07	3.25E-06	9.30E-C6	4.41E-05	1.33E-C4	3.05E-04	5.77E-04	9.59E-04	9.59E-04
46785 1.43E-15	1.18E-12	1.04E-10	2.56E-09	2.81E-08	1.81E-07	8.05E-07	7.51E-06	3.68E-05	1.21E-04	3.03E-04	6.31E-04	6.31E-04
92245 2.69E-29	1.54E-23	1.06E-19	5.87E-17	6.68E-15	2.65E-13	5.03E-12	4.15E-10	5.66E-09	1.02E-07	6.35E-07	2.73E-07	2.73E-07
1C9218 8.03E-35	5.29E-28	3.2P-23	8.94E-18	6.94E-16	2.29E-14	4.25E-12	1.77E-10	2.89E-09	2.52E-08	1.43E-07	1.43E-07	1.43E-07
155127 0.	0.	1.70E-33	6.97E-29	2.01E-25	9.04E-23	1.39E-17	2.35E-15	5.71E-13	2.50E-13	5.47E-12	6.43E-11	6.43E-11
144188 C.	5.34E-37	5.38E-31	1.04E-26	1.71E-23	5.41E-21	5.41E-19	5.39E-16	7.43E-14	2.98E-12	5.24E-11	5.17E-10	5.17E-10
166766 0.	0.	6.40E-36	5.83E-31	3.05E-27	2.38E-24	4.90E-22	1.44E-18	4.30E-16	3.08E-14	8.50E-13	1.21E-11	1.21E-11
2180CC 0.	0.	0.	0.	9.08E-35	5.49E-31	5.81E-28	1.99E-23	3.45E-20	9.20E-18	7.07E-16	2.27E-14	2.27E-14
225CCC 0.	0.	0.	0.	9.65E-37	9.05E-33	1.36E-29	7.92E-25	2.CCE-21	7.07E-19	6.77E-17	2.60E-15	2.60E-15
264CCC 0.	0.	0.	0.	0.	2.50E-38	1.15E-34	3.59E-29	3.COE-25	2.61E-22	5.03E-20	3.37E-18	3.37E-18
149056 0.	7.79E-38	1.25E-31	3.3ME-27	7.13E-24	2.74E-21	3.20E-19	4.03E-16	6.56E-14	2.98E-12	5.77E-11	6.6E-10	6.6E-10
169022 0.	0.	2.60E-35	2.77E-30	1.63E-26	1.39E-23	3.C7E-21	1.01E-17	3.25E-15	2.46E-13	7.12E-12	1.05E-10	1.05E-10
187693 0.	0.	0.	2.14E-33	3.28E-29	5.91E-26	2.37E-23	1.91E-19	1.17E-16	1.43E-14	5.99E-13	1.19E-11	1.19E-11
196955 0.	0.	0.	0.	9.50E-36	2.35E-31	6.12E-28	3.31E-25	4.13E-21	3.48E-18	5.40E-16	2.73E-14	2.66E-13
2C3384 0.	0.	0.	0.	2.03E-36	6.97E-32	2.35E-26	1.56E-25	2.65E-21	2.78E-18	5.10E-16	2.93E-14	7.45E-13
210284 0.	0.	0.	1.92E-37	9.71E-33	4.31E-29	3.57E-26	8.33E-27	7.48E-23	1.12E-18	2.46E-16	1.62E-14	4.60E-13
211271 C.	0.	0.	1.85E-37	9.53E-33	4.4CE-29	3.76E-26	9.35E-22	1.28E-18	2.88E-16	1.94E-14	5.59E-13	5.59E-13
2C7974 0.	0.	0.	2.05E-37	8.92E-33	3.61E-29	2.77E-26	5.52E-22	7.22E-19	1.49E-16	9.37E-15	2.57E-13	2.57E-13
226455 0.	0.	0.	0.	3.47E-35	2.94E-31	4.08E-26	2.10E-23	4.85E-20	1.61E-17	1.46E-15	5.39E-14	5.39E-14
2445CC 0.	0.	0.	0.	8.78E-38	1.53E-33	3.78E-30	4.62E-25	1.98E-21	1.04E-18	1.36E-16	6.70E-15	6.70E-15
2661CC 0.	0.	0.	0.	0.	4.90E-36	2.42E-32	8.33E-27	7.48E-23	6.87E-20	1.38E-17	9.58E-16	9.58E-16
2520CC C.	0.	0.	0.	0.	2.32E-35	7.20E-32	1.20E-26	7.C7E-23	4.52E-20	6.85E-18	3.79E-16	3.79E-16
27C0CC 0.	0.	0.	0.	0.	2.20E-37	1.23E-33	5.11E-28	5.24E-24	5.33E-21	1.16E-18	A.54E-17	A.54E-17
2880CC 0.	0.	0.	0.	0.	0.	1.15E-35	1.14E-29	2.16E-25	3.9E-22	1.08E-19	1.07E-17	1.07E-17
3059CC 0.	0.	0.	0.	0.	0.	6.96E-38	1.95E-31	7.83E-27	2.21E-23	1.06E-20	1.08E-18	1.08E-18
3231CC 0.	0.	0.	0.	0.	0.	0.	2.71E-33	1.72E-28	6.83E-25	4.28E-22	7.39E-20	7.39E-20
3548CC C.	0.	0.	0.	0.	0.	0.	2.41E-36	4.52E-31	4.06E-27	4.80E-24	1.37E-21	1.37E-21
3C80CC C.	0.	0.	0.	0.	0.	0.	3.30E-32	1.25E-27	3.38E-24	2.13E-19	2.13E-19	2.13E-19
3398CC C.	0.	0.	0.	0.	0.	0.	2.93E-35	3.25E-30	2.01E-26	1.76E-23	3.96E-21	3.96E-21
364CCC C.	0.	0.	0.	0.	0.	0.	9.95E-38	2.56E-32	2.91E-28	4.13E-25	1.37E-22	1.37E-22
3560CC 0.	0.	0.	0.	0.	0.	0.	6.33E-35	1.64E-30	4.41E-27	2.44E-4	2.44E-4	2.44E-4
38CCCC C.	0.	0.	0.	0.	0.	0.	2.35E-33	4.09E-29	4.06E-27	4.80E-26	3.43E-23	3.43E-23
412CCC C.	0.	0.	0.	0.	0.	0.	5.90E-36	2.30E-31	8.54E-28	6.10E-25	1.07E-25	1.07E-25
1C80CC 0.	0.	0.	0.	0.	0.	0.	4.53E-36	1.60E-31	5.46E-28	3.66E-25	1.07E-30	1.07E-30
4390CC 0.	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.

Table 7. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF O⁻

STATE	(CP-1)	LEVEL	STAT. #T.	TEMPERATURE (DEG K)				600	800	1000	1500
				200	250	300	400				
2s ² 2p ³ P _{1/2}	0	0.0353	4	9.40E-01	9.12E-01	8.87E-01	8.48E-01	8.20E-01	7.98E-01	7.70E-01	7.44E-01
	205	0.0353	2	6.05E-C2	8.04E-02	1.13E-01	1.52E-01	1.80E-01	2.02E-01	2.30E-01	2.49E-01
ENERGY LEVELS FROM HERRY ET AL. (43).											
LEVEL	(CP-1)	2CCC	2500	3CCC	3500	4000	4500	5000	6000	7000	9000
C	7.11E-01	7.02E-01	6.96E-01	6.92E-C1	6.89E-01	6.87E-01	6.85E-01	6.82E-01	6.78E-01	6.77E-01	6.76E-01
205	2.89E-01	2.94E-01	3.04E-01	3.08E-C1	3.11E-01	3.13E-01	3.15E-01	3.18E-01	3.20E-01	3.22E-01	3.24E-01

Table 8. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF 0

STATE (CP-1)	LEVEL (EV)	STAT. WT.	TEMPERATURE (CEG K)						1500
			200	250	300	400	500	600	
2s 2p ¹ P ^o	0	C.CCCC	5	8.12E-01	7.42E-01	7.00E-01	6.74E-01	6.55E-01	6.16E-01
	158.4	C.G196	3	1.56E-C1	1.86E-01	2.08E-01	2.38E-01	2.69E-01	2.95E-01
2p ¹ P ^o	226.5	0.0281	1	3.18E-02	4.19E-02	5.01E-02	6.22E-02	7.40E-02	8.90E-02
D ^o	15868	1.9673	5	C.	0.	6.6CE-34	1.14E-25	9.95E-21	1.95E-17
s	13792	4.1896	1	0.	0.	0.	0.	0.	1.46E-07
									1.00E-15
2s 2p ¹ P ^o	1263C4	15.6593	9	0.	0.	0.	0.	0.	0.
	189837	23.5362	3	0.	0.	0.	0.	0.	0.
2p ¹ S ^o	277CC0*	34.3427	1	C.	0.	0.	0.	0.	0.
2s 2p ¹ (S ^o)	3d	74903	9.2865	8	0.	0.	0.	0.	0.
		87379	10.8333	24	0.	0.	0.	0.	0.
									2.94E-32
3d	97443	12.0811	4C	0.	0.	0.	0.	0.	0.
	95757	11.8720	8	0.	0.	0.	0.	0.	0.
	99314	12.3130	24	0.	0.	0.	0.	0.	0.
4p ¹	1029C0**	12.7576	116	0.	0.	0.	0.	0.	0.
(D ^o)	3s	101523	12.5869	20	0.	0.	0.	0.	0.
									1.14E-36
3d	1136CC**	14.0842	60	0.	0.	0.	0.	0.	0.
	12394C**	15.3662	1CC	0.	0.	0.	0.	0.	0.
	1287C0**	15.9584	320	0.	0.	0.	0.	0.	0.
(P ^o)	3s	114416	14.1854	12	0.	0.	0.	0.	0.
	3p	1279CC**	15.8572	36	C.	0.	0.	0.	0.
									0.
3d	138CC0*	17.1094	6C	0.	0.	0.	0.	0.	0.
	1421CC*	17.6177	192	0.	0.	0.	0.	0.	0.
2s 2p ¹ (P ^o)	3	212CC0*	25.2840	216	0.	0.	0.	0.	0.
	4	222CC0*	27.5238	384	0.	0.	0.	0.	0.
(D ^o)	3	258CC0*	31.5871	18C	0.	0.	0.	0.	0.
									0.
4	268CCC*	33.6269	320	0.	0.	0.	0.	0.	0.
(S ^o)	3	287CCC*	35.5825	36	0.	0.	0.	0.	0.
	4	298CC0*	36.5463	64	0.	0.	0.	0.	0.
(P ^o)	3	304CC0*	37.6902	108	0.	0.	0.	0.	0.
	4	314CC0*	38.9300	192	0.	0.	0.	0.	0.
2p ¹ (P ^o)	3	409CC0*	5C.7082	1C8	0.	0.	0.	0.	0.
	4	419CC0*	51.9480	192	0.	0.	0.	0.	0.

*ESTIMATED.

**INCLUDES ESTIMATED SUBLVELS.

UNSTARRED ENERGY LEVELS FROM MCCREY (3) AND ACMEN (44).
NOTE: STATES INVOLVING ELECTRONS WITH PRINCIPAL QUANTUM NUMBERS ABOVE N = 4 ARE NOT INCLUDED.

ALL LEVELS ABOVE 109837 CP-1 ARE SUBJECT TO ALTCIONIZATION.

Table 8 (Cont.). ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF O

LEVEL (CP-1)	TEMPERATURE (CEG K)										ICCO
	2CC0	25CC	3CC0	35CC	4000	45CC0	5000	6CC0	700C	80CC	
0	5.86E-01	5.80E-01	5.76E-01	5.73E-C1	5.67E-01	5.65E-01	5.59E-01	5.52E-01	5.45E-01	5.38E-01	5.31E-01
15A	3.14E-C1	3.18E-01	3.20E-01	3.22E-C1	3.23E-01	3.24E-01	3.23E-01	3.21E-01	3.15E-01	3.11E-01	3.07E-01
226	9.96E-C2	1.02E-01	1.03E-01	1.04E-01	1.05E-01	1.06E-01	1.06E-01	1.05E-01	1.04E-01	1.03E-01	1.02E-01
15868	6.46E-06	6.22E-05	2.85E-04	8.42E-04	1.88E-03	3.55E-03	5.87E-03	1.24E-02	2.12E-02	3.14E-02	4.26E-02
33772	3.25E-12	4.15E-10	1.06E-07	6.00E-07	2.30E-06	6.76E-06	1.38E-05	3.38E-05	1.06E-04	2.5CE-04	4.85E-04
126304	0.	2.82E-32	5.11E-27	2.91E-23	1.91E-20	2.96E-18	1.67E-16	7.06E-14	5.28E-12	1.34E-10	1.65E-09
169937	0.	0.	C.	4.41E-35	7.56E-31	1.48E-27	6.39E-25	5.69E-21	3.75E-18	4.87E-16	2.13E-14
277000	0.	0.	0.	0.	0.	0.	2.73E-36	1.55E-30	2.07E-26	2.52E-23	6.21E-21
74903	3.72E-24	1.76E-19	2.31E-16	3.89E-14	1.82E-12	3.61E-11	3.94E-10	1.42E-08	1.62E-07	1.23E-06	1.77E-05
87379	1.41E-27	4.03E-22	1.75E-18	6.91E-16	6.13E-14	2.C0E-12	3.26E-11	2.13E-09	4.2CE-08	3.92E-07	2.22E-06
97443	1.69E-30	2.05E-24	2.33E-20	1.84E-17	2.73E-15	1.34E-13	3.C0E-12	3.10E-11	8.05E-09	1.07E-07	7.39E-07
95757	1.14E-30	1.08E-24	1.05E-20	7.35E-16	1.00E-15	4.59E-14	9.75E-13	9.53E-11	2.5CE-09	2.99E-C8	1.54E-C7
69314	2.64E-31	4.19E-25	5.70E-21	5.11E-18	8.37E-16	4.41E-14	1.05E-12	1.22E-10	3.62E-09	4.58E-08	3.29E-07
1C29CC	9.65E-32	2.57E-25	4.93E-21	5.66E-18	1.11E-15	6.77E-14	1.81E-12	2.49E-11	8.36E-CS	1.16E-07	8.96E-07
1C1523	4.48E-32	9.79E-26	1.65E-21	1.72E-18	3.15E-16	1.81E-14	4.64E-13	5.98E-11	1.91E-09	2.57E-08	1.92E-07
113600	2.27E-35	2.81E-28	1.51E-23	3.60E-20	1.23E-17	1.14E-15	4.31E-14	9.9CE-12	4.8CE-11	8.77E-09	8.38E-08
12394C	2.22E-38	1.22E-30	1.76E-25	8.55E-22	4.96E-19	7.C0E-17	3.66E-15	1.38E-12	9.54E-11	2.28E-09	2.67E-08
12877C	0.	5.25E-31	5.76E-26	3.R7E-22	4.87E-19	4.89E-17	2.98E-15	4.41E-12	1.15E-10	3.C9E-C9	3.08E-07
3.52E-36	3.52E-29	2.04E-24	5.15E-21	1.83E-18	1.76E-16	6.81E-15	6.63E-12	8.11E-11	1.51E-09	1.47E-08	9.C3E-08
0.	5.25E-32	9.51E-27	6.04E-23	4.3CE-20	7.10E-18	4.22E-16	1.93E-13	1.52E-11	4.02E-10	5.11E-09	3.89E-08
0.	0.	0.	C.	0.	0.	0.	3.C5E-35	1.16E-31	2.72E-26	1.05E-22	1.38E-19
0.	2.24E-34	1.25E-28	1.58E-24	1.69E-21	4.68E-19	3.84E-17	2.85E-14	3.18E-12	1.C9E-10	1.69E-09	1.52E-08
14210C	6.78E-35	5.59E-29	9.40E-25	1.39E-21	4.C4E-19	3.78E-17	3.41E-14	4.35E-12	1.67E-11	2.02E-09	2.69E-08
212CCC	C.	0.	3.51E-37	1.80E-32	6.94E-29	7.82E-26	2.C2E-21	2.94E-18	6.51E-16	4.77E-14	1.30E-13
2222000	0.	0.	0.	0.	9.15E-34	6.5CE-3C	7.82E-27	3.26E-22	6.47E-19	1.92E-16	5.48E-13
2580C0	0.	0.	C.	0.	0.	0.	3.C5E-35	1.16E-31	2.72E-26	1.05E-22	1.38E-19
1380CC	0.	0.	0.	0.	0.	0.	2.22E-36	1.16E-32	6.4CE-27	4.22E-23	6.C0E-2C
14210C	0.	0.	0.	0.	0.	0.	5.52E-36	5.19E-30	9.56E-26	1.50F-22	4.59E-20
2870CC	0.	0.	C.	0.	0.	0.	4.14E-37	6.6CE-31	1.27E-26	3.7CE-23	1.41E-20
2S80CC	0.	0.	0.	0.	0.	0.	1.24E-37	2.64E-31	8.71E-27	2.12E-23	9.10E-21
3C4000	0.	0.	0.	0.	0.	0.	4.27E-32	1.98E-27	6.24E-24	3.27E-21	4.88E-19
31400C	0.	0.	C.	0.	0.	0.	0.	0.	0.	0.	0.
4CSCCC	C.	0.	C.	0.	0.	0.	0.	0.	3.69E-36	1.33E-31	4.67E-28
419000	0.	0.	C.	0.	0.	0.	0.	0.	0.	3.93E-32	1.68E-28

Table 9. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF O^+

STATE	(CM-1)	LEVEL (EV)	STAT- N.	TEMPERATURE (EFG K)				ABC	AGC	IGCC	ISCO	
				2CC	2SC	3CC	4CC					
2s ² 2p ² ¹ S ⁰	C	0.0000	4	1.00E 0C	1.00E 0C	1.00E 0C	1.00E 0C	1.00E CC	1.00E CC	1.00E CO	1.00E CO	
	26819	5.3250	1C	C-	C-	C-	C-	0.	7.62E-34	2.94E-28	2.82E-17	
2s ² 2p ² ³ D ⁰	4C467	5.0171	6	C-	C-	C-	C-	0.	0.	0.	1.36E-17	
	119933	14.8694	12	C-	C-	C-	C-	0.	0.	0.	7.72E-32	
2s ² 2p ² ¹ P ⁰	165941	20.5797	1C	C-	C-	C-	C-	0.	0.	0.	2.08E-17	
	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.	0.	
2s ² 2p ² ³ S ⁰	195710	24.2643	2	C-	C-	C-	C-	0.	0.	0.	0.	
	212650	26.3646	6	C-	C-	C-	C-	0.	0.	0.	0.	
2s ² 2p ² ³ P ⁰	3174C0*	39.3516	6	C-	C-	C-	C-	0.	0.	0.	0.	
	1866C4	23.01354	1B	C-	C-	C-	C-	0.	0.	0.	0.	
2s ² 2p ² ³ P ¹	2C92C8	25.9378	54	0.	C-	C-	C-	0.	0.	0.	0.	
	3d	232563	28.8334	90	C-	C-	C-	C-	0.	0.	0.	0.
	4s	239348	29.6746	1B	C-	C-	C-	C-	0.	0.	0.	0.
	4p	246P60**	3C-6C59	54	C-	C-	C-	C-	0.	0.	0.	0.
	4d	255CC6	31.6159	90	0.	0.	0.	0.	0.	0.	0.	0.
	4f	255G32	31.7307	126	C-	C-	C-	C-	0.	0.	0.	0.
('D) 3s	206772	25.6606	1C	C-	C-	C-	C-	0.	0.	0.	0.	
	3p	229898	28.5C30	3C	0.	0.	0.	0.	0.	0.	0.	0.
	3d	252571	31.3140	5C	0.	0.	0.	0.	0.	0.	0.	0.
	4s	27293C**	33.8381	16C	C-	C-	C-	C-	0.	0.	0.	0.
('S) 3s	226851*	28.1252	2	C-	C-	C-	C-	0.	0.	0.	0.	
	3p	25C251*	31.0264	6	C-	C-	C-	C-	0.	0.	0.	0.
	3d	27591*	34.2227	1C	0.	0.	0.	0.	0.	0.	0.	0.
	4s	296CCC*	36.6984	32	0.	0.	0.	0.	0.	0.	0.	0.
2s ² 2p ² ¹ S ⁰) 3	2A1CC0*	34.8387	9C	0.	0.	0.	0.	0.	0.	0.	0.	
	313CC0*	36.8C61	16C	C-	C-	C-	C-	0.	0.	0.	0.	
(³ D ⁰) 3	34CCCC*	42.1515	27C	0.	0.	0.	0.	0.	0.	0.	0.	
	4	373CCC*	46.2645	48C	C-	C-	C-	C-	0.	0.	0.	0.
(³ P ⁰) 3-4	3A3CCC*	47.4847	45C	0.	0.	0.	0.	0.	0.	0.	0.	
(³ S ⁰ , ¹ D ⁰ , ³ P ⁰) 3-4	437CCC*	54.1797	55C	C-	C-	C-	C-	0.	0.	0.	0.	
2p ² (³ S ⁰) 4	534CC0*	66.2059	75C	0.	0.	0.	0.	0.	0.	0.	0.	

*ESTIMATED.

**INCLUDES ESTIMATED SUBLVELS.

ACSF: STARRED ENERGY LEVELS FROM MCCREY (3) AND ERIKSSON (45).

ACTF: STATES INVOLVING ELECTRONS WITH PRINCIPAL QUANTUM NUMBERS ABOVE N = 4 ARE NOT INCLUDED.

ALL LEVELS ABCVE 283244 CM-1 ARE SUBJECT TO AUTONORMALIZATION.

Table 9 (Cont.). ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF O^+

LEVEL ICP-1)	2CCC	2CCO	3CCC	3CCO	4CCC	4CCO	5CCC	5CCO	6CCC	6CCO	8CCC	8CCO	10CCC	
C	1.CCE CC	1.CCC CO	1.CCE CC	1.CCE CC	1.CCE CC	1.CCE CC	9.SSE-CI	9.26E-01	S.SCE-CI	9.79E-01	9.65E-01	9.46E-01		
26815	1.C4E-CR	4.95E-07	6.45E-06	4.07E-05	1.62E-04	4.72E-14	1.11E-C3	4.C1E-03	9.59E-03	1.97E-02	3.31E-02	4.95E-02		
46667	1.41E-13	1.15E-10	5.59E-09	8.94E-CB	7.15E-07	3.60E-16	1.31E-05	9.11E-05	3.62E-05	1.01E-04	2.24E-03	4.20E-03		
11931	1.C2E-37	3.14E-25	1.16E-21	5.52E-15	6.56E-17	3.C8E-15	9.66E-13	5.84E-11	1.26E-09	1.36E-08	5.1CE-08			
16991	C.	6.47E-35	5.80E-30	2.94E-26	2.23E-22	4.5CCE-21	1.29E-17	3.77E-15	2.65E-13	7.21E-12	1.CCE-10			
145710	C.	0.	0.	5.74E-36	1.34E-31	1.33E-28	1.74E-25	2.07E-21	1.68E-18	2.53E-16	1.25E-14	2.79E-13		
212650	C.	C.	C.	1.63E-38	5.C6E-34	4.44E-30	3.98E-27	1.07E-22	1.55E-19	3.61E-17	2.49E-15	7.32E-14		
317400	C.	C.	C.	C.	0.	0.	C.	1.32E-33	6.SCE-29	2.38E-25	1.33E-22	2.08E-20		
186604	C.	C.	C.	C.	2.18E-33	3.18E-29	5.52E-26	2.15E-23	1.65E-19	9.80E-17	1.17E-14	4.81E-13	9.31E-12	
209208	C.	C.	C.	C.	6.C3E-37	2.81E-32	1.20E-28	5.46E-26	2.19E-21	2.82E-18	6.C2E-16	3.89E-14	1.C8E-12	
232563	C.	0.	C.	C.	1.C5E-35	1.15E-31	1.94E-28	1.35E-23	3.87E-20	1.51E-17	1.55E-15	6.25E-14		
279348	C.	C.	C.	C.	1.83E-37	2.42E-33	5.51E-30	5.31E-25	1.52E-21	8.9CE-19	1.C5E-16	4.71E-15		
248860	C.	C.	C.	C.	3.65E-38	7.11E-34	1.90E-30	2.63E-25	1.23E-21	6.91E-16	9.45E-15	4.80E-15		
245506	C.	C.	C.	C.	C.	8.76E-35	3.04E-31	6.21E-26	3.84E-22	2.66E-19	4.28E-17	2.48E-15		
255932	C.	0.	C.	C.	9.12E-35	3.26E-31	6.97E-26	4.45E-22	3.16E-19	5.17E-17	3.C9E-15			
206672	C.	C.	C.	C.	2.80E-37	1.16E-32	4.55E-29	3.40E-26	6.94E-22	H.28E-19	1.67E-16	1.C3E-14	2.76E-13	
229898	C.	0.	C.	C.	9.15E-36	9.SCE-32	1.39E-28	8.53E-24	2.23E-20	8.11E-18	7.9CE-16	3.C6E-14		
252571	C.	0.	C.	C.	C.	1.C6E-34	3.41E-31	6.13E-26	3.52E-22	2.29E-19	3.51E-17	1.95E-15		
272920	C.	0.	C.	C.	C.	5.C5E-37	3.11E-33	1.5CE-27	1.72E-23	1.80E-20	4.34E-18	3.34E-16		
226851	C.	0.	C.	C.	1.82E-36	1.58E-32	2.23E-29	1.18E-24	2.78E-21	9.36E-19	8.58E-17	3.16E-15		
250251	C.	C.	C.	C.	C.	2.67E-35	7.97E-32	1.30E-26	6.8CE-23	4.17E-18	6.11E-18	3.27E-16		
275951	C.	0.	C.	C.	C.	1.2CE-38	8.15E-35	4.55E-29	5.76E-25	6.84E-22	1.67E-19	1.35E-17		
256000	C.	C.	C.	C.	C.	0.	8.15E-37	1.17E-30	2.59E-26	5.45E-23	2.17E-20	2.42E-18		
2PLCCC	0.	C.	C.	C.	C.	1.72F-34	1.22E-2H	1.84E-24	2.48E-21	6.72E-15	5.88E-17			
3.30CC	C.	C.	C.	C.	C.	C.	1.C1E-31	4.54E-27	1.4CE-23	7.17E-21	1.C5E-18			
34CL00	C.	0.	C.	C.	C.	0.	C.	2.62E-34	2.98E-29	1.84E-25	1.61E-22	3.C3E-20		
37CCCC	C.	C.	C.	C.	C.	C.	0.	6.C1E-32	8.43E-28	1.47E-24	5.CCE-22			
35CCCC	C.	C.	C.	C.	C.	C.	C.	7.21E-31	1.34E-28	2.78E-25	1.24E-22			
4100CC	C.	0.	C.	C.	C.	C.	C.	0.	9.9CE-33	6.C6E-25	6.44E-26			
514GCC	C.	C.	C.	C.	C.	C.	C.	C.	1.52E-35	7.61E-32				

Table 10. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF Ar

STATE	(CP-1) LEVEL (EV)	STAT. WT.	2CC	25C	3CC	7C ATURE (DEG K)	600	8CO	10CO	15CC
$3s^2 3p^4 (^1S)$	C	C.CCCO	1	1.CCE CC	1.CCE OC	1.CCE 00	1.CCE CC	1.CCE OC	1.CCE OC	1.CCE OC
$3s^2 3p^4 (^3P_{1/2})$ 3d	114854	14.2397	2C	C.	C.	0.	0.	0.	0.	C.
4s	95186	11.8015	4	0.	C.	0.	0.	0.	0.	C.
4p	107421	13.3182	12	C.	C.	C.	C.	C.	C.	C.
4d	120721	14.9671	2C	0.	0.	0.	C.	C.	C.	C.
3s ² 3p ⁴ (^3P _{3/2}) 4f	121654	15.0828	28	0.	0.	0.	0.	0.	0.	0.
4s	112948	14.0394	4C	0.	0.	0.	0.	0.	0.	C.
4s	93371	11.5762	8	C.	0.	0.	0.	0.	0.	C.
4p	105631	13.0962	24	C.	C.	C.	C.	C.	C.	C.
4d	119211	14.7799	4C	0.	0.	0.	0.	0.	0.	0.
3s ² 3p ⁴ (^1S)	4f	120222	14.952	56	C.	C.	0.	0.	0.	C.
3d	222000*	27.5238	2C	C.	C.	0.	0.	0.	0.	C.
4s	203CCC*	25.1661	4	C.	C.	0.	0.	0.	0.	C.
4p	215CCC*	26.6559	12	C.	C.	C.	C.	C.	C.	C.
4d	2275CC*	28.2553	2C	0.	C.	C.	0.	0.	0.	C.
4f	2289CC*	28.3793	2H	C.	0.	0.	C.	C.	0.	C.

*ESTIMATED.

ENERGY LEVELS FROM BURRS AND ADAMS (146) AND HUMPHREYS AND PAUL (47).

NOTE: STATES INVOLVING ELECTRONS WITH PRINCIPAL QUANTUM NUMBERS ABOVE $n = 4$ ARE NOT INCLUDED.
ALL LEVELS ABOVE 127110 CP-1 ARE SUBJECT TO ARTIFICIALIZATION.

Table 10 (Cont.). ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF Ar

LEVEL (C _{w-1})	TEMPERATURE (E _G K)									
	2000	2500	3000	3500	4000	4500	5000	6000	7000	8000
C 1.00E+00	1.00E+00	1.00E+00	1.00E+00	1.00E+00	1.00E+00	1.00E+00	1.00E+00	1.00E+00	1.00E+00	1.00E+00
114054 C 2.61E-35	1.93E-29	6.39E-23	2.25E-20	7.29E-17	2.29E-14	2.25E-12	8.66E-11	2.15E-11	1.12E-10	2.14E-09
95188 C 7.29E-30	6.46E-24	5.97E-20	4.66E-17	5.46E-15	2.42E-13	5.08E-12	4.85E-11	1.27E-10	1.47E-09	1.21E-08
1C7421 C 1.29E-33	1.70E-26	9.07E-22	7.97E-19	1.99E-16	1.46E-14	4.51E-13	7.80E-12	3.05E-11	4.51E-10	4.51E-09
120121 C 3.44E-37	1.34E-29	1.43E-24	5.61E-21	2.77E-18	3.45E-16	1.64E-14	5.76E-13	1.35E-12	4.88E-11	4.18E-10
121654 C 2.75E-37	1.10E-29	1.28E-24	5.35E-21	2.77E-18	3.59E-16	1.75E-14	5.99E-12	3.87E-11	6.31E-10	7.41E-09
112948 C 2.06E-34	2.95E-27	1.19E-22	2.74E-19	9.08E-17	6.29E-15	3.07E-13	6.91E-11	3.31E-10	5.76E-09	3.50E-08
93371 C 5.39E-29	3.68E-23	2.85E-19	1.71E-16	2.08E-14	8.67E-13	1.72E-11	1.51E-10	3.70E-09	4.08E-08	2.63E-07
1C56- C 2.39E-32	9.52E-26	2.39E-21	3.33E-18	7.57E-16	5.16E-14	1.51E-12	2.4CE-11	8.53E-10	1.35E-09	1.11E-08
119211 C 2.28E-36	6.40E-29	5.92E-24	2.09E-20	9.54E-18	1.12E-15	5.06E-14	1.54E-11	5.13E-10	1.15E-09	2.12E-08
12C222 C 1.54E-36	5.01E-29	5.10E-24	1.93E-20	9.28E-18	1.13E-15	5.29E-14	1.69E-11	1.04E-10	2.52E-09	1.71E-08
222000 C. 0.	0.	C.	C.	4.18E-34	2.98E-30	3.01E-27	1.52E-22	3.05E-19	3.05E-17	7.72E-15
2C3000 C. 0.	0.	C.	C.	2.29E-36	7.77E-32	2.59E-29	1.71E-25	2.85E-21	3.08E-18	3.02E-16
215000 C. 0.	0.	C.	C.	4.95E-38	3.11E-33	1.68E-29	1.62E-26	4.88E-22	7.71E-19	1.42E-16
227900 C. C. 0.	0.	C.	C.	5.C1E-35	4.522E-31	6.61E-28	3.69E-23	5.C7E-20	1.16E-17	3.C1E-15
226900 C. C. 0.	0.	C.	C.	4.085E-35	4.66CE-31	6.44E-28	4.06E-23	1.03E-19	1.7CE-17	3.59E-15

Table 11. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF Ar⁺

STATE	(CP-1)	LEVEL	LEVEL	STATISTICAL WEIGHT						TEMPERATURE (CEG K)					
				2CC	2CC	250	100	400	500	600	800	1000	1500	1000	1500
3 ² 3p ² 3p _{3/2}	C	0.CCCC	4	1.00E+00	9.59E-01	9.97E-01	9.84E-01	9.63E-01	9.40E-01	9.17E-01	8.88E-01	8.63E-01	8.39E-02	8.05E-02	7.71E-02
3 ² 3p ² 3p _{1/2}	C	1432.0	C-1775	2	1.6AE-05	1.32E-04	5.20E-04	2.89E-03	8.05E-03	3.67E-02	1.56E-02	5.99E-02	C.	C.	C.
3 ² 3p ² 3p _{3/2}	C	108723	134796	2	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
3 ² 3p ² (1P)3d ¹ D	C	132476	164245	2C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
3d other	C	146319	181408	70	C.	C.	C.	C.	C.	C.	C.	C.	C.	C.	C.
4s	C	136028	168649	18	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4p	C	158023	195918	54	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4d	C	186693	231464	90	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4f	C	191567	242466	176	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
(D)3d	C	164082	203431	5C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4s	C	148754	184427	1C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4p	C	171631	211303A	3C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4d	C	199657	247537	5C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4f	C	209029	259156	7C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
(S)3d	C	179728	222829	1C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4s	C	167109	2017431	2	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4p	C	192095	2388161	6C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4d	C	227758	2772758	1C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4f	C	228000	282677	14	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
3e 3p ² (1P)3d	C	269000	333509	9C	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4s	C	245000	3078713	18	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4p	C	273000	338468	54	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4d	C	301000	373183	90	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.
4f	C	309000	383101	126	0.	0.	0.	0.	0.	0.	0.	0.	C.	C.	C.

*ESTIMATED.

ENERGY LEVELS FROM MCCREY (31) AND MINNICHEN (48).

NOTE: STATES INVOLVING ELECTRONS WITH PRINCIPAL QUANTUM NUMBERS ABOVE N = 4 ARE NOT INCLUDED.
ALL LEVELS ABOVE 22248 CP-1 ARE SUBJECT TO ALIGNMENT.

Table 11 (Cont.). ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF Ar⁺

LEV [L] (CP-L)	2CC0	2SC0	3CC0	3SCC	4CCCC	4SCCO	5CCC	6CCCO	7CCC	8OCC	9CCO	10CCO
0	8.49E-01	8.20E-01	7.59E-01	7.03E-01	7.7CE-01	7.51E-01	7.38E-01	7.29E-01	7.21E-01	7.15E-01	7.11E-01	7.11E-01
1432	1.51E-01	1.80E-01	2.01E-01	2.17E-01	2.3CE-01	2.49E-01	2.62E-01	2.79E-01	2.95E-01	2.89E-01	2.89E-01	2.89E-01
1C0723	4.56F-35	2.74E-28	9.04E-24	1.52E-20	3.95E-10	3.04E-16	9.72E-15	1.76E-12	7.8E-11	1.16E-09	1.01E-08	5.72E-08
132416	C.	3.17E-33	1.02E-27	8.74E-24	7.78E-21	1.53E-18	1.4CE-16	5.9CE-14	5.44E-12	1.62E-10	2.27E-09	1.87E-08
146319	0.	3.85E-36	4.67E-30	1.01E-25	1.87E-22	6.40E-20	6.81E-18	7.47E-15	1.11E-12	4.70E-11	8.69E-10	8.95E-09
136028	C.	3.70E-34	1.67E-28	1.81E-24	1.95E-21	4.42E-19	3.38E-17	2.26E-14	2.36E-12	7.70E-11	1.16E-09	1.11E-08
158023	0.	0.	1.31E-32	6.49E-28	7.14E-24	1.17E-21	1.81E-19	3.48E-16	7.71E-14	4.42E-12	1.03E-10	1.24E-09
186693	0.	0.	0.	8.23E-33	1.19E-28	2.04E-25	7.88E-23	5.99E-19	3.54E-16	4.25E-14	1.76E-12	3.45E-11
155567	C.	C.	C.	3.C01E-34	6.03E-30	1.67E-26	8.59E-24	9.99E-20	8.C1E-17	1.21E-14	5.96E-13	1.35E-11
164082	0.	0.	0.	6.66E-34	4.98E-29	2.25E-25	1.56E-22	2.93E-20	7.53E-17	2.C5E-14	1.38E-12	3.63E-11
148734	C.	1.35E-37	2.08E-31	5.42E-27	1.11E-23	4.20E-21	4.83E-19	5.55E-16	5.59E-14	4.34E-12	8.41E-11	9.01E-10
171831	0.	0.	9.72E-36	1.23E-30	8.3CE-27	7.86E-24	1.89E-21	7.05E-18	2.51E-15	2.05E-13	6.3CE-12	9.77E-11
199657	0.	0.	0.	2.22E-35	6.22E-31	1.79E-27	1.05E-24	1.49E-20	1.37E-17	1.23E-15	2.2E-13	2.97E-12
2C9029	0.	0.	0.	6.59E-37	2.99E-32	1.25E-28	9.51E-26	2.2CE-21	2.8CE-18	5.95E-16	3.05E-14	1.09E-12
179728	C.	0.	7.34E-38	1.60E-32	1.62E-28	2.10E-25	6.5CE-23	3.54E-19	1.65E-16	1.65E-14	5.95E-13	1.05E-11
1.7309	0.	0.	5.67E-36	5.28E-31	2.81E-27	2.23E-24	4.63E-22	1.39E-18	4.23E-16	3.00E-14	8.66E-13	1.25E-11
192095	0.	0.	5.95E-35	1.13E-30	2.41E-27	1.11E-24	1.05E-22	7.78E-18	1.02E-15	4.94E-14	1.02E-12	1.02E-12
22C00C	0.	0.	0.	8.76E-35	5.37E-31	6.02E-28	2.26E-23	4.19E-20	1.18E-17	9.51E-16	3.16E-14	3.16E-14
22BCCC	C.	C.	C.	6.51E-36	5.82E-32	8.44E-29	4.65E-24	1.13E-20	3.32E-18	3.71E-16	1.41E-14	1.41E-14
26900C	0.	0.	0.	0.	0.	7.59E-37	4.08E-33	1.61E-27	1.59E-23	1.58E-20	3.39E-18	2.48E-16
249CCC	0.	0.	C.	0.	0.	9.C9E-35	2.58E-31	3.85E-26	1.15E-19	1.66E-17	8.93E-16	8.93E-16
27300C	0.	0.	0.	0.	0.	1.27E-37	7.74E-34	3.69E-28	4.20E-24	4.63E-21	1.C7E-18	8.38E-17
3C1CCC	0.	0.	C.	0.	0.	4.C9E-37	7.47E-31	2.22E-26	5.01E-23	2.04E-20	2.49E-18	2.49E-18
3C900C	0.	0.	0.	0.	0.	5.73E-38	1.54E-31	5.59E-27	1.66E-23	7.93E-21	1.1CE-18	1.1CE-18

Table 12. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF CO

		STATE							
		$\chi \ 1\Sigma^+$	$a \ 3\Pi$	$a' \ 3\Sigma^+$	$d \ 3\Delta$	$e \ 3\Sigma^-$	$\iota \ 1\Sigma^-$	$A \ 1\Pi$	
ENERGY (CM ⁻¹)	(EV)	0	49474	55354	60647	63709	64547	64747	
		0.0000	6.0099	6.8628	7.5191	7.8987	8.0026	8.0273	
VIB. INT. (CM ⁻¹)	(EV)	2143	1715	1209	1138	1094	1071	1481	
INT.		0.2657	0.2126	0.1439	0.1410	0.1357	0.1328	0.1836	
TEMP. (DEG K)		FRACTIONAL POPULATION							
200	1.00E 00	0.	0.	C.	0.	0.	0.	0.	
250	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
300	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
400	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
500	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
600	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
800	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
1000	1.00E 00	3.68E-30	1.30E-34	0.	0.	0.	0.	0.	
1500	1.00E 00	4.77E-20	4.85E-23	6.43E-25	1.78E-26	2.73E-27	2.98E-27		
2000	1.00E 00	5.49E-15	3.03E-17	1.44E-18	8.34E-20	1.57E-20	1.73E-20		
2500	1.00E 00	5.99E-12	9.17E-14	9.35E-15	8.48E-16	1.80E-16	1.99E-16		
3000	1.00E 00	6.37E-10	1.33E-11	3.28E-12	4.01E-13	9.20E-14	1.02E-13		
3500	1.00E 00	1.79E-08	8.85E-10	2.17E-10	3.28E-11	7.94E-12	8.83E-12		
4000	1.00E 00	2.19E-07	1.56E-08	5.04E-09	8.96E-10	2.25E-10	2.51E-10		
4500	1.00E 00	1.53E-06	1.46E-07	5.84E-08	1.18E-08	3.03E-09	3.40E-09		
5000	1.00E 00	7.29E-06	8.77E-07	4.15E-07	9.26E-08	2.42E-08	2.73E-08		
6000	1.00E 00	7.58E-05	1.29E-05	7.91E-06	2.04E-06	5.45E-07	6.21E-07		
7000	9.99E-01	4.05E-07	8.87E-05	6.48E-05	1.86E-05	4.97E-06	5.73E-06		
8000	9.98E-01	1.42E-03	3.74E-04	3.12E-04	9.60E-05	2.57E-05	2.99E-05		
9000	9.94E-01	3.77E-03	1.14E-03	1.05E-03	3.40E-04	9.08E-05	1.07E-04		
10000	9.85E-01	8.15E-03	2.75E-03	2.011E-03	9.17E-04	2.45E-04	2.89E-04		

Based on energy-level data from references 49 and 57.

Table 13. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF N₂

		STATE							
		X 1Σ ⁺ _g	A 3Σ ⁺ _u	3Δ _u	B 3Π _g	B' 3Σ ⁻ _u	a' 1Σ ⁻ _u	a 1Π _g	w 1Δ _u
ENERGY (CM ⁻¹)	(EV)	0	49756	~58CCC	59310	65852	67739	68951	71698
	0.000	6.1688	~7.1909	7.3533	8.1644	8.3984	8.5486	8.88c2	
VIE. (CM ⁻¹)	2330	1433	~1464	1705	1493	1506	1666	1535	
INT. (EV)	0.2888	0.1777	~0.1815	C.2114	0.1851	0.1867	0.2066	0.19c4	
TEMP. (DEG K)		FRACTIONAL POPULATION							
200	1.00E 00	0.	C.	C.	0.	0.	0.	0.	0.
250	1.00E 00	0.	0.	C.	0.	0.	0.	0.	C.
300	1.00E 00	C.	0.	0.	U.	0.	0.	0.	0.
400	1.00E 00	0.	0.	0.	0.	0.	0.	0.	C.
500	1.00E 00	C.	C.	U.	0.	0.	0.	0.	0.
600	1.00E 00	0.	0.	C.	0.	0.	0.	0.	0.
800	1.00E 00	0.	0.	0.	0.	0.	0.	0.	0.
1000	1.00E 00	3.72E-31	0.	0.	0.	0.	0.	0.	0.
1500	1.00E 00	9.34E-21	6.77E-24	1.60E-24	1.78E-27	9.62E-29	5.27E-29	4.22E-30	
2000	1.00E 00	1.51E-15	7.87E-18	2.51E-18	1.36E-20	1.15E-21	8.33E-22	1.30E-22	
2500	1.00E 00	2.03E-12	3.46E-14	1.31E-14	1.84E-16	2.05E-17	1.75E-17	4.0C9E-18	
3000	1.00E 00	2.49E-10	9.33E-12	3.97E-12	1.05E-13	1.40E-14	1.34E-14	4.0C9E-15	
3500	1.00E 00	7.75E-09	5.1CE-10	2.35E-10	9.83E-12	1.49E-12	1.54E-12	5.69E-13	
4000	1.00E 00	1.02E-07	1.03E-08	5.04E-09	2.96E-10	4.94E-11	5.40E-11	2.31E-11	
4500	1.00E 00	7.64E-07	1.06E-07	5.47E-08	4.19E-09	7.54E-10	8.62E-10	4.13E-10	
5000	1.00E 00	3.82E-06	6.91E-07	3.69E-07	3.50E-08	6.68E-09	7.92E-09	4.15E-C9	
6000	1.00E 00	4.28E-05	1.15E-05	5.47E-06	8.48E-07	1.77E-07	2.21E-07	1.33E-07	
7000	1.00E 00	2.40E-04	8.56E-05	5.02E-05	8.29E-06	1.84E-06	2.39E-06	1.59E-06	
8000	9.98E-01	6.67E-04	3.86E-04	2.33E-04	4.59E-05	1.07E-05	1.43E-05	1.02E-05	
9000	9.95E-01	2.33E-03	1.24E-03	7.64E-04	1.74E-04	4.19E-05	5.76E-05	4.34E-05	
10000	9.89E-01	5.08E-03	3.13E-03	1.96E-03	5.00E-04	1.25E-04	1.75E-04	1.38E-04	

Tables 13 to 18 are based on energy-level data from reference 50.

Table 14. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF N_2^+

ENERGY (CM ⁻¹) (EV)	VIB. INT. (EV)	STATE						TEMP. (DEC K)	FRACTIONAL POPULATION
		X 2Σ ⁺ g	A 2Π _u	B 2Σ ⁺ _u	4 Σ ⁺ _u	4 Δ _u	D 2Π _g	4 Σ ⁻ _u	C 2Σ ⁺ _u
0	0.0000	9016	25566	~44328	~51328	51203	~60328	~7.4795	64542 8.C020
1.1178	3.1697	~5.4958	~6.3637	6.3482	6.3482	~6.3482	~7.4795	64542 8.C020	
2175	1873	2371	~1668	~1572	889	~1472	2051		
0.2696	0.2323	0.2940	~0.2068	~0.1949	0.1102	~0.1825	0.2543		
200	1.00E 00	1.50E-28	0.	0.	0.	0.	0.	0.	
250	1.00E 00	6.44E-23	0.	0.	0.	0.	0.	0.	
300	1.00E 00	3.67E-19	0.	0.	0.	0.	0.	0.	
400	1.00E 00	1.84E-14	0.	0.	0.	0.	0.	0.	
500	1.00E 00	1.19E-11	1.04E-32	0.	0.	0.	0.	0.	
600	1.00E 00	9.05E-10	2.19E-27	0.	0.	0.	0.	0.	
800	1.00E 00	2.03E-07	9.87E-21	5.92E-35	0.	0.	0.	0.	
1000	1.00E 00	5.26E-06	9.68E-17	5.08E-28	4.37E-32	4.75E-32	0.	0.	
1500	1.00E 00	4.07E-04	2.02E-11	9.09E-19	2.26E-21	2.58E-21	2.37E-25	1.67E-27	
2000	9.96E-01	3.59E-03	9.16E-09	3.88E-14	5.21E-16	6.18E-16	4.76E-19	8.84E-21	
2500	9.87E-01	1.32E-02	3.57E-07	2.32E-11	8.57E-13	1.05E-12	2.87E-15	9.49E-17	
3000	9.69E-01	3.11E-02	4.05E-06	1.63E-09	1.18E-10	1.50E-10	9.45E-13	4.56E-14	
3500	9.43E-01	5.67E-02	2.26E-05	3.34E-08	3.96E-09	5.16E-09	5.87E-11	3.70E-12	
4000	9.12E-01	8.79E-02	8.13E-05	3.18E-07	5.43E-08	7.24E-08	1.28E-09	9.88E-11	
4500	8.78E-01	1.22E-01	2.17E-04	1.81E-06	4.11E-07	5.58E-07	1.40E-08	1.25E-09	
5000	8.42E-01	1.57E-01	4.73E-04	7.17E-06	2.06E-06	2.82E-06	9.34E-08	9.46E-09	
6000	7.74E-01	2.25E-01	1.49E-03	5.49E-05	2.25E-05	3.10E-05	1.58E-06	1.90E-07	
7000	7.12E-01	2.84E-01	3.32E-03	2.27E-04	1.21E-04	1.65E-04	1.15E-05	1.57E-06	
8000	6.59E-01	3.33E-01	6.01E-03	6.37E-04	4.20E-04	5.56E-04	4.99E-05	7.40E-06	
9000	6.13E-01	3.74E-01	9.46E-03	1.39E-03	1.09E-03	1.33E-03	1.53E-04	2.41E-05	
10000	5.73E-01	4.06E-01	1.35E-02	2.52E-03	2.28E-03	2.80E-03	3.65E-04	6.05E-05	

Table 15. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF NO

		STATE							
		x 2_{Π}	a 4_{Π}	A 2_{Σ}^+	B 2_{Π}	b 4_{Σ}^-	c 2_{Π}	d 2_{Σ}^+	
ENERGY (CM $^{-1}$) (EV)	INT. (EV)	62*	~37965	44149 5.4799	45505 5.6418	~47092 ~5.8385	52380** 6.4941**	53291 6.6071	
VIB. INT. (EV)		1876	995	2342	1023	~1203	2365**	2279	
		0.2326	0.1234	0.2903	0.1268	~0.1491	0.2932**	0.2826	
TEMP. (DEG K)		FRACTIONAL POPULATION							
200	1.00E 00	0.	0.	C.	0.	0.	0.	0.	
250	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
300	1.00E 00	C.	0.	0.	0.	0.	0.	0.	
400	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
500	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
600	1.00E 00	0.	0.	0.	0.	0.	0.	0.	
800	1.00E 00	8.47E-30	1.4CE-35	0.	0.	0.	0.	0.	
1C00	1.00E 00	7.49E-24	1.08E-28	7.45E-29	5.89E-30	1.67E-33	2.26E-34		
15C0	1.00E 00	6.58E-16	1.63E-19	2.41E-19	3.99E-20	1.27E-22	2.68E-23		
2C00	1.00E 00	6.31E-12	6.26E-15	1.40E-14	3.34E-15	3.48E-17	9.07E-18		
2500	1.00E 00	1.56E-C9	3.49E-12	1.01E-11	3.04E-12	6.30E-14	1.85E-14		
3C00	1.00E 00	6.20E-08	2.34E-10	8.24E-10	2.87E-10	9.31E-12	2.93E-12		
35C0	1.00E 00	8.61E-07	4.65E-09	1.91E-08	7.43E-09	3.26E-10	1.06E-10		
4000	1.00E 00	6.17E-06	4.32E-08	2.02E-07	8.56E-08	4.65E-09	1.54E-09		
45C0	1.00E 00	2.83E-05	2.41E-07	1.27E-06	5.74E-07	3.63E-08	1.21E-08		
5C00	1.00E 00	9.52E-05	9.40E-07	5.51E-06	2.63E-06	1.86E-07	6.20E-08		
60C0	9.99E-01	5.73E-04	7.00E-06	4.98E-05	2.57E-05	2.09E-06	6.89E-07		
7000	9.98E-01	2.00E-03	2.82E-05	2.37E-04	1.29E-04	1.13E-05	3.68E-06		
8CCC	9.94E-01	4.94E-03	7.75E-05	7.56E-04	4.22E-04	3.88E-05	1.24E-05		
9C00	9.87E-01	9.73E-03	1.65E-04	1.83E-03	1.04E-03	9.85E-05	3.11E-05		
10CC0	9.77E-01	1.63E-02	2.94E-04	3.65E-03	2.11E-03	2.02E-04	6.29E-05		

* Average energy of the two spin components, above that of the lower component ($^2\Pi_z$).

** "Perturbed" values.

Table 16. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF NO⁺

TEMP. (DEG K.)	FRACTIONAL POPULATION					
	X 1 _Σ ⁺	a 3 _Σ ⁺	3 _Δ	STATE	3 _Π	A 1 _Π
ENERGY (CM-1) (EV)	0	~39934	~58804	~63764	73084	~73334
VIB. INT.	(CM-1) (EV)	0.2344 0.2907	~1572 ~0.349	~1972 ~0.2445	~1672 ~0.2073	1562 0.1937
200	1.00E 00	0.	0.	0.	0.	0.
250	1.00E 00	0.	0.	0.	0.	0.
300	1.00E 00	0.	0.	0.	0.	0.
400	1.00E 00	0.	0.	0.	0.	0.
500	1.00E 00	0.	0.	0.	0.	0.
600	1.00E 00	0.	0.	C.	0.	0.
800	1.00E 00	2.46E-31	0.	0.	0.	0.
1000	1.00E 00	4.39E-25	0.	0.	0.	0.
1500	1.00E 00	9.76E-17	2.70E-24	2.01E-26	1.06E-30	1.87E-30
2000	1.00E 00	1.48E-12	3.67E-18	9.19E-20	4.55E-23	8.79E-23
2500	1.00E 00	4.80E-10	1.76E-14	9.15E-16	1.75E-18	3.54E-18
3000	1.00E 00	2.28E-08	5.04E-12	4.25E-13	2.01E-15	4.13E-15
3500	1.00E 00	3.60E-07	2.86E-10	3.43E-11	3.10E-13	6.33E-13
4000	1.00E 00	2.86E-06	5.94E-09	9.27E-10	1.36E-11	2.72E-11
4500	1.00E 00	1.43E-05	6.31E-08	1.21E-08	2.59E-10	4.99E-10
5000	1.00E 00	5.22E-05	4.19E-07	9.45E-08	2.73E-09	5.04E-09
6000	1.00E 00	3.63E-04	7.22E-06	2.08E-06	9.30E-08	1.57E-07
7000	9.98E-01	1.45E-03	5.57E-05	1.90E-05	1.14E-06	1.75E-06
8000	9.96E-01	4.10E-03	2.58E-04	9.93E-05	7.39E-06	1.04E-05
9000	9.90E-01	9.14E-03	8.48E-04	3.56E-04	3.09E-05	4.00E-05
10000	9.79E-01	1.72E-02	2.18E-03	9.78E-04	9.52E-05	1.15E-04

Table 17. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF O_2

		STATE						
		X $^3\Sigma_g^-$	a $^1\Delta_g$	b $^1\Sigma_g^+$	c $^3\Delta_u$	A $^3\Sigma_u^+$	c $^1\Sigma_u^-$	B $^3\Sigma_u^-$
ENERGY (CM-1)	0 (EV)	0.0000	0.9773	1.13121	~34329	35004	36213	49358
VIB. INT. (CM-1)	1556 (EV)	0.1930	0.1839	1.1483	~4.2561	4.3398	4.4897	6.1194
TEMP. (DEC K)				0.1742	~0.1017	0.0960	0.0764	0.0853
					FRACTIONAL POPULATION			
200	1.00E 00	1.60E-25	0.	0.	0.	0.	0.	0.
250	1.00E 00	1.34E-20	5.53E-34	C.	0.	0.	0.	0.
300	1.00E 00	2.58E-17	1.62E-28	0.	0.	0.	0.	0.
400	1.00E 00	3.29E-13	1.10E-21	0.	0.	0.	0.	0.
500	1.00E 00	9.56E-11	1.39E-17	0.	0.	0.	0.	0.
600	1.00E 00	4.20E-09	7.54E-15	0.	0.	0.	0.	0.
800	1.00E 00	4.76E-07	1.99E-11	5.78E-27	9.21E-28	4.47E-29	0.	0.
1000	1.00E 00	8.14E-06	2.25E-09	1.43E-21	2.90E-22	2.24E-23	3.73E-31	
1500	1.00E 00	3.60E-04	1.24E-06	2.30E-14	6.30E-15	9.18E-16	8.26E-21	
2000	9.98E-01	2.39E-03	2.91E-05	9.29E-11	2.89E-11	5.68E-12	1.25E-15	
2500	9.92E-01	7.44E-03	1.94E-04	1.32E-08	4.37E-09	1.01E-09	1.57E-12	
3000	9.84E-01	1.58E-02	6.82E-04	3.47E-07	1.19E-07	3.05E-08	1.78E-10	
3500	9.72E-01	2.68E-02	1.67E-03	3.46E-06	1.21E-06	3.32E-07	5.08E-09	
4000	9.57E-01	3.98E-02	3.25E-03	1.88E-05	6.68E-06	1.92E-06	6.08E-08	
4500	9.41E-01	5.37E-02	5.43E-03	6.82E-05	2.45E-05	7.26E-06	4.08E-07	
5000	9.23E-01	6.81E-02	8.16E-03	1.86E-04	6.73E-05	2.05E-05	1.83E-06	
6000	8.88E-01	9.60E-02	1.48E-02	7.96E-04	2.90E-04	9.17E-05	1.64E-05	
7000	8.53E-01	1.21E-01	2.24E-02	2.11E-03	7.75E-04	2.51E-04	7.39E-05	
8000	8.21E-01	1.43E-01	3.00E-02	4.20E-03	5.55E-03	5.10E-04	2.19E-04	
9000	7.91E-01	1.61E-01	3.72E-02	6.94E-03	2.56E-03	8.56E-04	4.93E-04	
10000	7.65E-01	1.75E-01	4.36E-02	1.01E-02	3.74E-03	1.26E-03	9.21E-04	

Table 18. ENERGY LEVELS AND EQUILIBRIUM FRACTIONAL ELECTRONIC POPULATIONS OF O_2^+

		STATE			
		$x\ 2\Pi_g$	$a\ 4\Pi_u$	$A\ 2\Pi_u$	$b\ 4\Sigma_g^-$
ENERGY (CM ⁻¹)	98*	32571	38303	49238	
(EV)	0.0122*	4.0382	4.7488	6.1046	
VIB. INT. (CM ⁻¹)	1843	1015	872	1163	
INT. (EV)	0.2286	0.1258	0.1081	0.1441	
TEMP. (DEG K)		FRACTIONAL POPULATION			
200	1.00E 00	0.	0.	0.	0.
250	1.00E 00	0.	0.	0.	0.
300	1.00E 00	0.	0.	0.	0.
400	1.00E 00	0.	0.	0.	0.
500	1.00E 00	0.	0.	0.	0.
600	1.00E 00	5.00E-34	0.	0.	0.
800	1.00E 00	1.52E-25	2.81E-30	0.	0.
1000	1.00E 00	1.89E-20	2.81E-24	2.96E-31	
1500	1.00E 00	1.22E-13	2.92E-16	5.54E-21	
2000	1.00E 00	3.16E-10	3.06E-12	7.71E-16	
2500	1.00E 00	3.56E-08	8.02E-10	9.51E-13	
3000	1.00E 00	8.38E-07	3.31E-08	1.10E-10	
3500	1.00E 00	8.03E-06	4.72E-07	3.31E-09	
4000	1.00E 00	4.39E-05	3.46E-06	4.25E-08	
4500	1.00E 00	1.65E-04	1.67E-05	3.10E-07	
5000	9.99E-01	4.74E-04	5.54E-05	1.52E-06	
6000	9.97E-01	2.29E-03	3.41E-04	1.63E-05	
7000	9.92E-01	6.92E-03	1.21E-03	8.73E-05	
8000	9.81E-01	1.55E-02	3.01E-03	2.99E-04	
9000	9.65E-01	2.82E-02	5.94E-03	7.58E-04	
10000	9.44E-01	4.44E-02	9.93E-03	1.56E-03	

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* Average energy of the two spin components, above that of the lower component ($^2\Pi_{\frac{1}{2}}$).

Table 19. LOWER ELECTRONIC AND VIBRATIONAL ENERGY LEVELS OF SELECTED DIATOMIC MOLECULES. UNITS: cm^{-1} and ev

Molecule	State	Electronic Energy	Lowest Vibrational Interval	References
H_2	$X^1\Sigma_g^+$	0 cm^{-1} 0 ev	4161 0.516	51
H_2^+	$X^2\Sigma_g^+$	0 0	2191 0.272	14
CO^+	$X^2\Sigma^+$	0 0	2184 0.271	52
	$A^2\Pi$	20408 2.530	1535 0.190	
	$B^2\Sigma^+$	45633 5.658	1679 0.208	
NO^-	$X^3\Sigma^-$	0 0	(~1600) (~0.20)	*
O_2^-	$X^2\Pi_g$	0 0	" (0.143)	53
OH^-	$X^1\Sigma^+$	0 0	(3600) (0.446)	28
OH	$X^2\Sigma^-$	0 0	3570 0.443	52
	$A^2\Sigma^+$	32402 4.017	2989 0.371	
OH^+	$X^3\Sigma^-$	0 0	2967 0.368	52
	$A^3\Pi$	27952 3.466	1986 0.246	

*Vibrational interval estimated from O_2 .

Table 20. VIBRATIONAL SPACING OF TRIATOMIC MOLECULES

Molecule	Ground State	Vibrational Intervals (cm^{-1} ; ev)			References
		ν_1	ν_2	ν_3	
H_2O	${}^1\text{A}_1$	3657 0.453	1595 0.198	3756 0.466	32
H_2O^+	(${}^2\text{B}_1$) (~0.40)	(~3200) (~0.40)	(~1500) (~0.19)	(~3300) (~0.41)	*
CO_2	${}^1\Sigma_g^+$	1388 0.172	667 0.083	2349 0.291	32
CO_2^+	${}^2\Pi_g$	1280 0.159	(~400) (~0.05)	(1469) (0.182)	32
NO_2^-	${}^1\text{A}_1$	(1325) (0.164)	(829) (0.103)	(1270) (0.157)	54
NO_2	${}^2\text{A}_1$	(1320) (0.164)	750 0.093	1618 0.201	32
NO_2^+	(${}^1\Sigma_g^+$) (0.174)	(1400) (0.174)	(538) (0.067)	(2360) (0.293)	55
N_2O	${}^1\Sigma^+$	2224 0.276	589 0.073	1285 0.159	32
N_2O^+	${}^2\Pi$	1737 0.215	461 0.057	1126 0.140	32
O_3^-	(${}^2\text{B}_1$) (0.156)	(1260) (0.156)	(800) (0.099)	(1140) (0.141)	56
O_3	${}^1\text{A}_1$	1110 0.138	705 0.087	1042 0.129	32
O_3^+	(${}^2\text{A}_1$) (~0.16)	(~1300) (~0.16)	(~700) (~0.09)	(~1600) (~0.20)	†

*Vibrational intervals estimated from Rydberg states of H_2O .

†Vibrational intervals estimated from NO_2 .

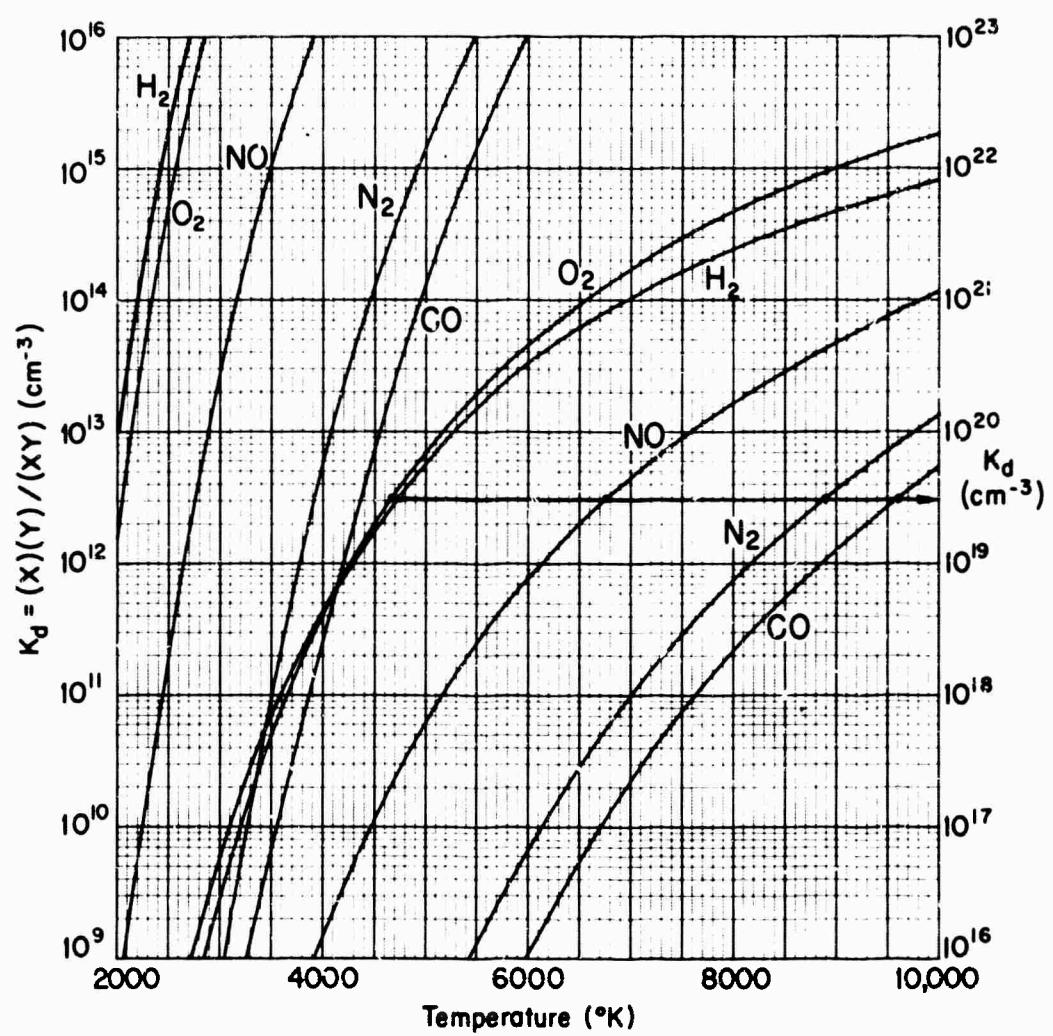


Fig. 1--Equilibrium constants for dissociation

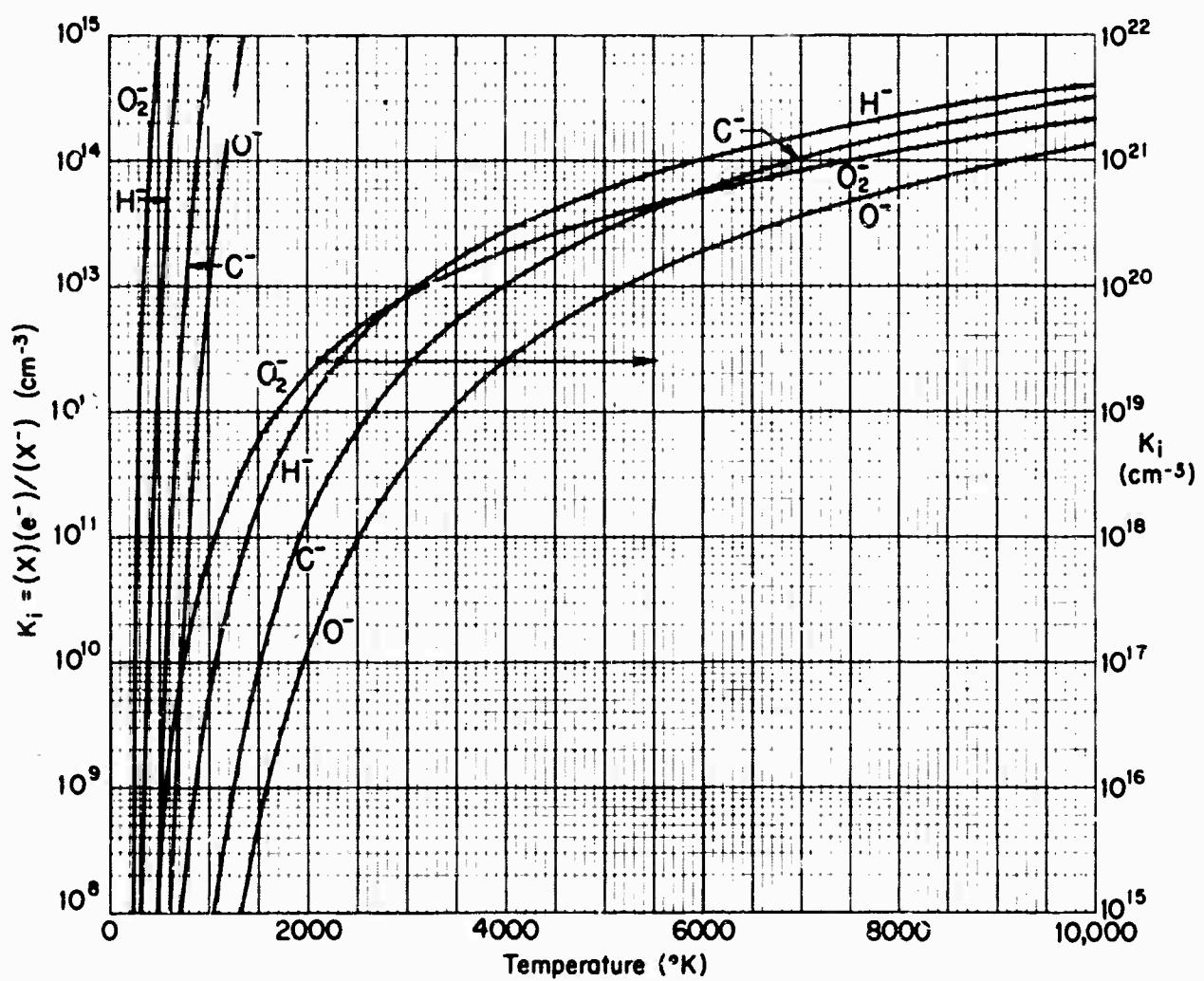


Fig. 2--Equilibrium constants for ionization (detachment) of negative ions

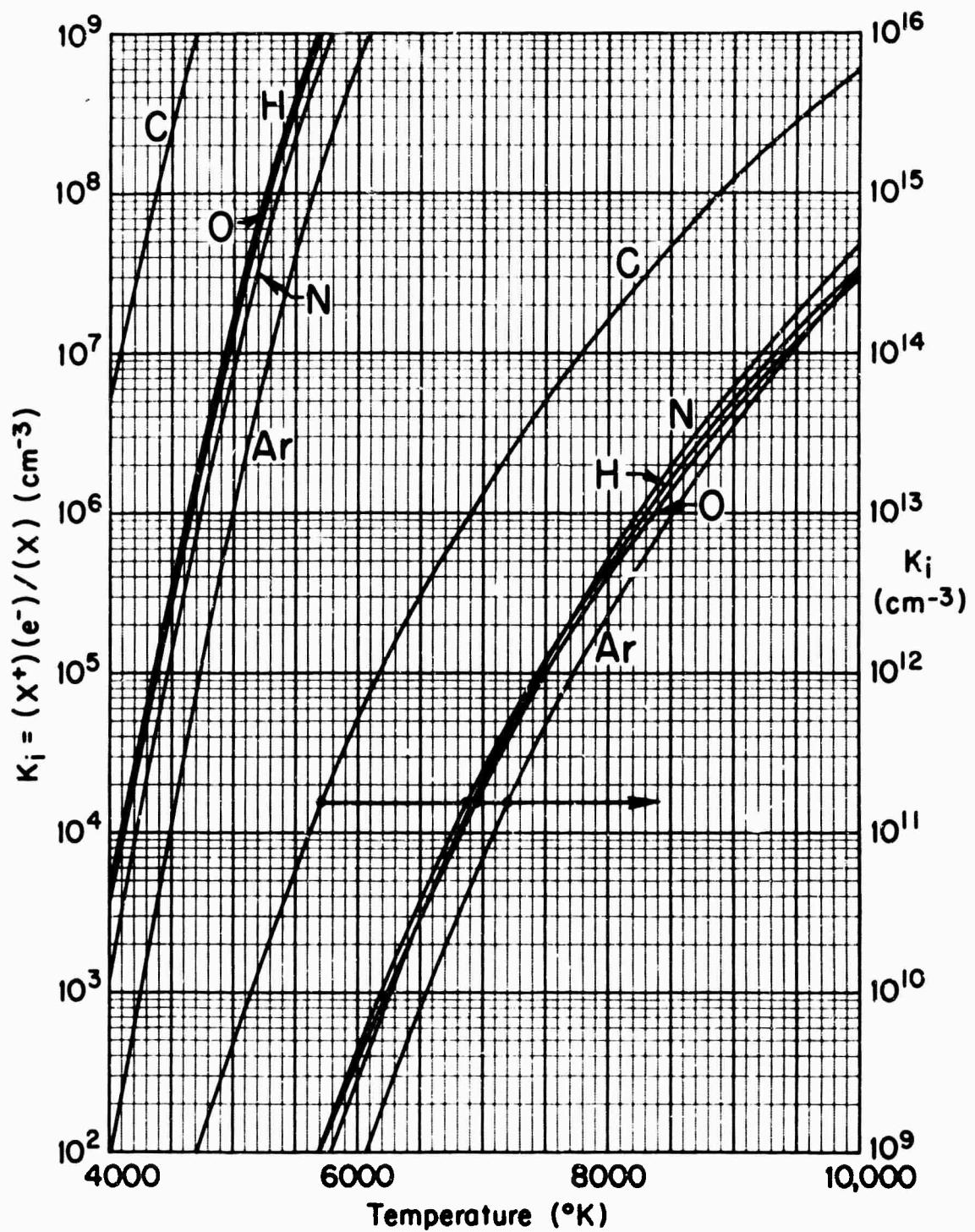


Fig. 3--Equilibrium constants for ionization of atoms

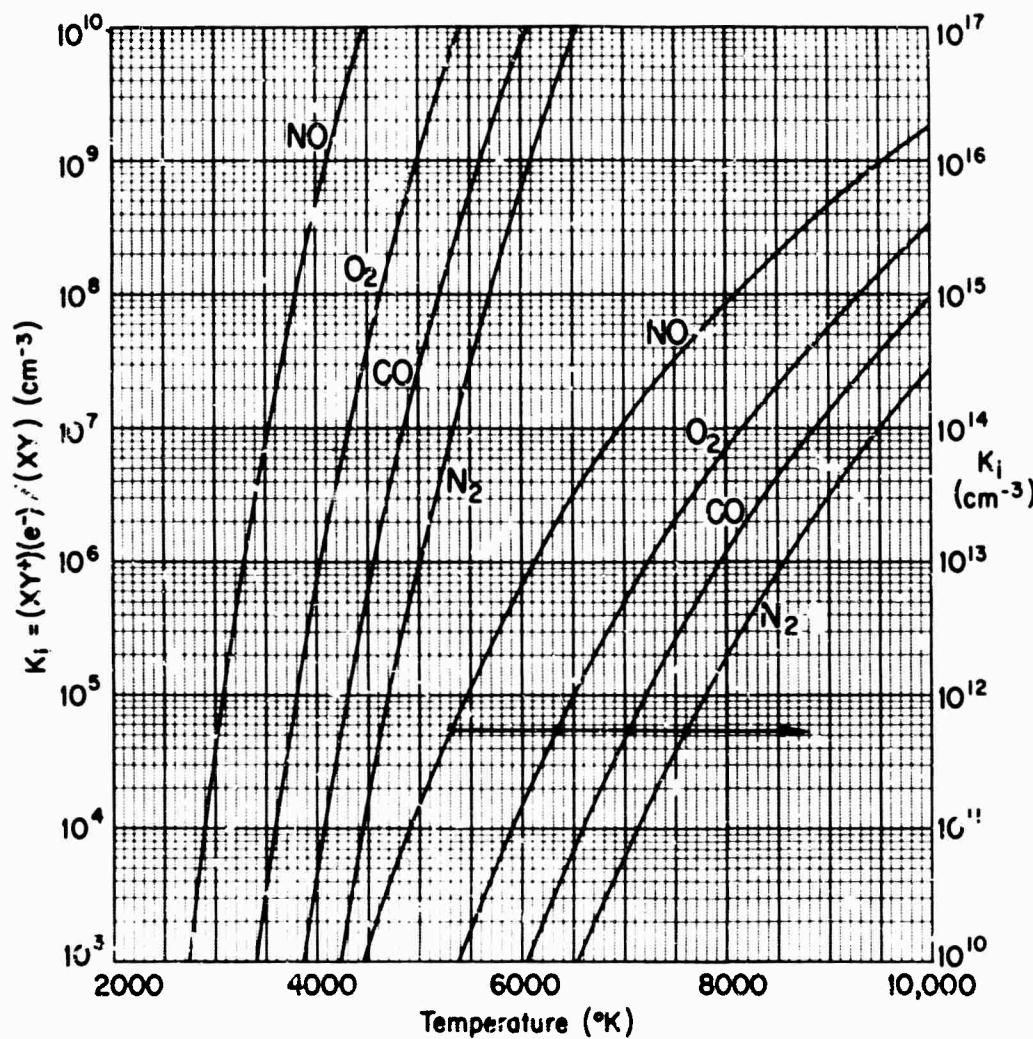


Fig. 4--Equilibrium constants for ionization of diatomic molecules

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10. ABSTRACT Tables of the formation energies, dissociation energies, ionization energies, electronic energy levels, and vibrational level spacings for most atomic, diatomic, and triatomic molecules involving hydrogen, carbon, nitrogen, oxygen, and argon are presented. Many positively and negatively charged ions are included. Tables of the equilibrium fractional electronic-state populations, and graphs of the equilibrium constants for dissociation, ionization, and detachment for most of the atomic and diatomic species are appended. A brief discussion of the significance of such data precedes the tables and graphs.		II. KEY WORDS Physics Reentry vehicles Atmosphere Radiation Ionization