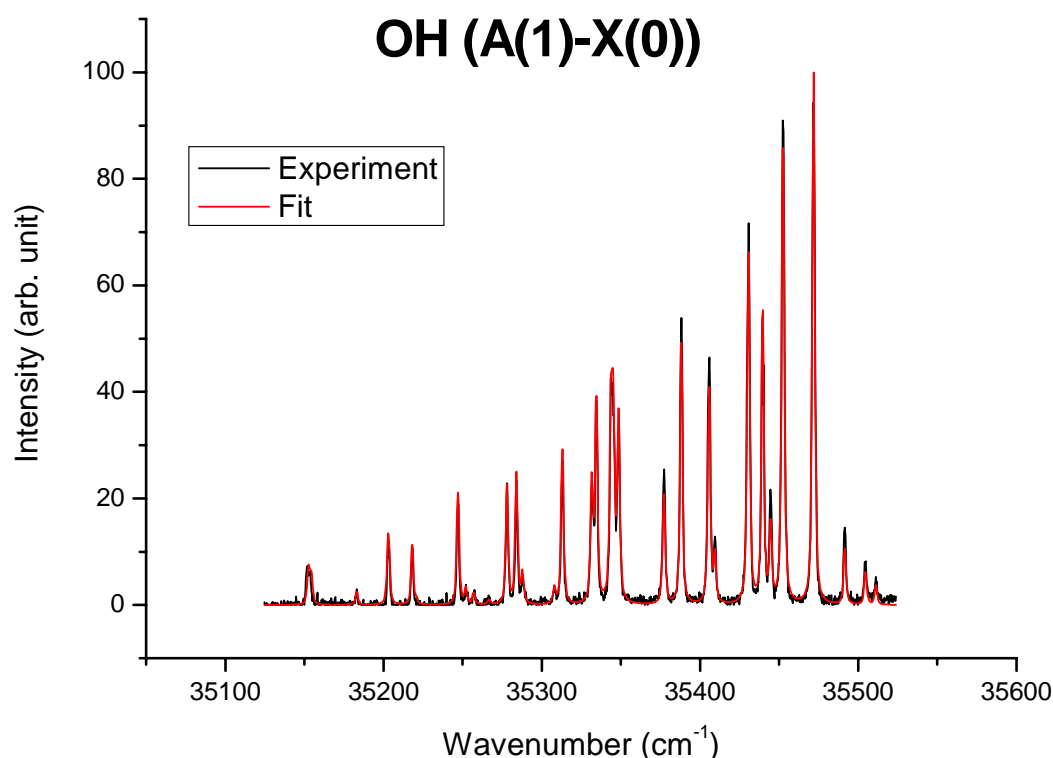


Molecular constants of diatomic molecules

The following pages are copies from the book Constants of Diatomic Molecules by Huber and Herzberg, containing information about the most common diatomic molecules of interest in combustion.

For each molecule there is information in a large number of columns. The parameters T_e , ω_e , $\omega_e x_e$, B_e , D_e , and v_{00} are given in the unit cm^{-1} (which is a way to express energy). The column with references needs no attention for the course exercises.



State	T_e	w_e	$w_e x_e$	B_e	α_e	D_e (10^{-4} cm^{-1})	r_e (\AA)	Observed Transitions Design. ν_{00}	References
NH									
$\mu = 0.94016028$ $D_0^0 = 3.47 \text{ eV}^a$ I.P. = (13.63) eV^b									
Theoretical potential functions and spectroscopic constants for the ground and excited states: (21)(44)(46); (32)(36)(51).									
d 1E^+	83160	2672.6	Z 71.2	14.390 ^c	0.621	16.0 ^d	1.116 ₃	$d^+ \rightarrow c, m$ V 39512.2 ₆ Z	(5)* (19) (23)(27) (55)*
c 1E	(43744)	[2122.6 ₄]	Z f	14.537 ^{ghi}	0.593 ^j	[22.0] ^k	1.110 ₆	$d^+ \rightarrow b, m$ R 61619.6 ₀ Z $c^+ \rightarrow b, m$ R 22106.6 ₂ Z $c^+ \rightarrow a, m$ R 30755.5 ₄ Z	(29)(55)* (4)* (26)* (1)(2)* (6) (14)*
A 2D_1	29807.4 ⁿ	3231.2	Z 98.6	16.6745 ^{op1}	0.7454	[17.80] ^q	1.0369 ₈	$A^+ \rightarrow X, m$ * 29776.76 Z	(3)(8)* (15) (35)
b 1E^+	21202	3352.4	Z 74.2 ^t	16.705 ^u	0.591	16.0 ^v	1.0360	$b^+ \rightarrow X, m$ * 21238 Z	(45)
a 1E	(12566)*	[3188]	Z (68) ^y	[16.439] ^l	0.66	[16.2]	1.034 ₁	(a-X) 12589 ^z	
X 2E^+	0	3282.2 ₇	Z 78.3 ₅	16.6993 ^{a'}	0.6490	[17.097] ^{b'}	1.0362 ₁	Rotation sp. c^+ Fundamental b. in matrices (45a)(55a)	(48)(54)

State	T_e	w_e	$w_e x_e$	B_e	α_e	D_e (10^{-6} cm^{-1})	r_e (\AA)	Observed Transitions Design. ν_{00}	References
CN									
$\mu = 6.46219329$ $D_0^0 = 7.76 \text{ eV}^a$ I.P. = 14.1 ₇ eV^b									
Theoretical work (24)(46)(48)(50)(63). RKR potential curves (15)(44)(53). Franck-Condon factors (21a)(53), and ref. in (27)(44).									
J 2D_1	65258.19 ^c	1121.76	Z 14.203 ^d	1.305 ₂	0.020 ₈	5.8	1.413 ₇	J \rightarrow A, R 55667.14 ^e Z	(8)*
K $2\text{D}(r)$	[61969.7] ^f			[1.520]			[1.310]	H \rightarrow B, R 35140.8 ₄ * Z	(7)
G 2D	[61655.0] ^{gh}			[1.085] ^h			[1.551]	G \rightarrow B, R 34826.1 ₆ * Z	(7)(43)
F 2D_r	60095.6 ₄ ⁱ	1239.5 ₀	Z 12.7 ₃	1.383 ₄	0.0187	7	1.373 ₂	F \rightarrow A, R 50563.8 ₀ * Z	(7)* (22) (47)
E 2E^+	59151.1 ₈	1681.4 ₃	Z 3.60 ^j	1.4871	0.0064 ₃ ^k	5.0	1.324 ₅	E \rightarrow A, R 49842.4 ₂	(8)* (47)
D 2D_1	54486.3 ^l	1004.7 ₁ ^m	Z 8.7 ₈	1.162 ⁿ	0.013	7	1.498	E \rightarrow X, R 58959.5 ₅ Z D \rightarrow A, R 44838.0 ₈ * Z	(8)* (43)* (7)*
A $4\text{E}^+(*)$	(32400) ⁿ	2163.9 ^o	Z 20.2 ^p	1.973 ^o	0.023 ^p	[6.6]	1.150	D \rightarrow X, R 53955.4 ₆	(7)*
B 2E^+	25752.0							B \rightarrow A, V 16680.4 ₆	(31)(54)* (1)* (2)* (4a)(53)(7)
[For a comprehensive review of molecular data on the B-X system including a bibliography of references prior to 1972 see (44).]									
A 2D_1	9245.28 ^q	1812.5 ₆ ^r	Z 12.60 ₂ ^t	1.7151 ^u	0.01708 ^u	5.93 ^{uv}	1.2333	B ⁺ \rightarrow A, X, V 25797.84 Z A ⁺ \rightarrow X, X, R 9117.35 ^z Z	(9)(32)(54) (3)* (3a) (4b)(7)(9) (15)* (25) (32)(44)*
[For a useful compilation of molecular data on the A-X system and a bibliography of references prior to 1971 see (44).]									
X 2E^+	0	2068.59	Z 13.087 ^y	1.8997 ₄ ^z	0.01736 ₉ ^{a'}	6.40 ^{b'}	1.1718 ₂	IR fundamental b. Microwave sp. c^+ ESR sp. d^+	(62)(70) (66)(71) (41)(73)

State	T_e	w_e	$w_e x_e$	B_e	α_e	D_e (10^{-6} cm^{-1})	r_e (\AA)	Observed Transitions Design. ν_{00}	References
CH									
$\mu = 0.92974056$ $D_0^0 = 3.46 \text{ eV}^a$ I.P. = 10.64 eV^b									
Several unassigned absorption bands ^c above 80000 cm^{-1} . Rydberg series joining on to G, $v = 85850 - R/(n - 0.09)^2$; $n = 3, 4, 5, 6$									
G d	[74373]							G \rightarrow X, 72960	(26)
F 2E^+	[65943]							F \rightarrow X, R 64531.5 ^f Z	(26)*
E 2E	[65625]							E \rightarrow X, R 64211.7 ^f Z	(26)*
D 2D_1	[60394] ^h							D \rightarrow B, V 33282.8 ^f Z	(26)*
C 2E^+	31801.5	2840.2	Z 125.9 ₄ ⁱ	14.603 ₃ ^k	0.7185 ^l	[15.55] ^m	1.114 ₃	D \rightarrow X, 58981.0 ^f Z	(26)*
B 2E^+	(26044)	[1794.9] ^p	Z	[12.645] ^q ^k	F	[22.2] ⁿ	1.1975 ₅	C ⁺ \rightarrow X, v_R 31778.1 ^f Z	(11)(3)(10)* (26)
A 2E	23189.8 ^u	2930.7	Z 96.65	14.934 ^v ^k	0.697	15.4 ^w	1.1019	B ⁺ \rightarrow X, v_R 25698.2 ^f Z	(2)* (3) (10)* (25)*
A 4E^+	(5844)	(3145)	(72)	(15.4)	(0.55)		(1.08 ₅)	A ⁺ \rightarrow X, v_R 23217.5 ^f Z	(3)(6)(10)* (12)* (48)
X 2D_r	0 ^t	2858.5	Z 63.0 ₂	14.457 ^{a'} ^{b'} ^k	0.534	14.5	1.1199	5985 ^y A doubling sp. a^+ Ab initio calculations; ref. in (44)	(39)

State	T_e	w_e	$w_e x_e$	B_e	α_e	D_e (10^{-2} cm^{-1})	r_e (\AA)	Observed Transitions Design. ν_{00}	References
H₂									
Q (1D_g) ^a	(113163)	[742]		[(16.3)]			[(1.43)]	Q \rightarrow B, R 21151.1	(4)(24)
B ⁺ 1E^+ $3p6$	111642.8 ^b	2039.52	83.406 ^c	26.70 ₅ ^e	2.781 ^d	[1.2] ^f	1.119 ₂	B ⁺ \rightarrow E, P 11311.5 ^g	(34)
F ¹ 1E^+ $2p6^2$	100911 ^j	[1199] ⁱ		$B_4 = 6.24k$			$r_4 = 2.31k$	B ⁺ \rightarrow X, h R 110478.2	(40)(44)(46)
E ¹ 1E^+ $2s6$	100082.3 ^m	2588.9 ^m	130.5 ^m	32.68 ^m	1.818 ^m	[2.28] ^m	1.011 ₈	F \rightarrow B, l R $v_{40} = 13635.1$	(14)(34)
C 1D_u $2p\pi$	100089.8 ⁿ	2443.77	69.524 ^{op}	31.362 ₉ ^p	1.664 ₇ ^q	2.23 ^r	1.0327 ₉	E \rightarrow B, l V 8961.23	(8)(22)(24) (34)
B 1E^+ $2p6$	91700.0 ^a	1358.09	20.888 ^b	20.015 ₄ ^c	1.1845 ^d	1.625 ^e	1.2928 ₂	C ⁺ \rightarrow X, t R 99120.1 ⁿ	(12)(37)(44) (129)
X 1E_g $1s6^2$	0	4401.21 ₃	121.33 ₆ ⁱ	60.853 ₀ ^k	3.062 ₂ ^j	4.71 ^l	0.74144	Werner b. Lyman b. Quadrupole ^m and field-induced sp. ⁿ Raman sp. ^o Rotational ^p and nuclear rf magn. reson.	(25)(77)(129) (15)(48) (26)(56)(78) (23)(56) (17a)(21) (17)(19)

State	T_e	w_e	$w_e x_e$	B_e	α_e	D_e (10^{-6}cm^{-1})	r_e (\AA)	Observed Transitions Design. ν_{00}	References
OH									
		$\mu = 0.94808710$		$D_0^0 = 4.392 \text{ eV}^a$		$I.P. = 12.90 \text{ eV}^b$			MAY 1977 Δ
		Theoretical potential functions for 48 states (92); for $X^2\text{H}$ and $A^2\text{E}^+$ see (112)(113)(118).							
C 2E^+	89459.1	1232.9	Z 19.1	4.247 ^{c,d}	0.078	2	2.046 ₁	C \rightarrow A, R 55820.7 Z (C \rightarrow X) E (88223)	(20)(42)(64)* (20)(45)* (106)
D 2E^-	(82130)	(2954)		[15.2179] ^f		[16.16]	[1.0809 ₃]	D \rightarrow X, E 81759.7 ^h Z	(16)(20)(33) (42)* (33)* (64)* (73)
B 2E^+	69774	[660.0]	Z 1	[5.086] ^{j,d}	K	[9.29] ^f	[1.869 ₈]	B \rightarrow A, R 35965.5 Z	(16)(20)(33) (42)* (33)* (64)* (73)
A 2E^+	32684.1	3178.8 ₆	Z 92.91 ⁿ	17.358 ^{OPQd}	0.786 ₈ ^f	[20.39] ^g	1.0121	A \rightarrow X, U R 32402.3 ^h Z	(3)(57)(64)* (69a)(90)
X 2D_1	0 ^v	3737.76 ₁	Z 84.881 ^w	18.910 ₈ ^{xd}	0.7242 ^f	19.38 ^t	0.95966	$\frac{1}{2} + \frac{3}{2}$ A \rightarrow 126.2	

State	T_e	w_e	$w_e x_e$	B_e	α_e	D_e (10^{-6}cm^{-1})	r_e (\AA)	Observed Transitions Design. ν_{00}	References
O₂									
F $^+$	[87510]	Group of mix line-like features similar to F $^+$ -X.							
F 3D_u	(85868)	[2008]	H v=1	[1.434]		[11]	[1.212]	F \rightarrow X, 85720	(17)(173)*
F 3D_g	(85780)	[2000]	H diffuse	[1.395]		[6.0]	[1.228]	F \rightarrow X, 85992.6 ^a Z	(136)(151) (173)*
E 1E^-	(79883)	[2001]	H b	[1.352]		[5.3]	[1.249]	E \rightarrow X, R 80369 ^b	(6)* (12) (150)(173)*
F 1E^-	76091	[2547]						f \rightarrow b, V 63141.5 Z	(84)
		1927	19.0	1.703 ^d	0.020	*	1.113	f \rightarrow X, V 76262.4 ^f	(84)(111)* (173)*
D 3E^+	(75260)	1957	19.7	1.73 ^h	0.025	i	1.10 ₄	D \rightarrow X, V (75450)	(84)(111)* (173)*
e $(^1\text{D}_{2u})$	(75254)	[1830]	H	[1.682]	(diffuse lines)		[1.119]	e \rightarrow a, V 67499.6 ^h Z	(84)(118) (146)(152)
e $(^1\text{D}_{2g})$	(74915)	[2052]	H j					e \rightarrow a, V 67272 ^k H	
d $(^1\text{E}_g)$	(69180)	[1860]	i					(d \rightarrow X) 69320 ^m	(192)
C $(^3\text{E}_g)$	(65530)	[1840]	n					(C \rightarrow X) 55670 ⁿ	(150)(171) (5)* (7)* (21)* (78)* (156)* (115)(168)
B 3E^-	49793.28	709.31 ^o	Z 10.65 ^o	0.81902 ^{OPQ}	0.01206 ^o	4.55 ^f	1.60426	B \rightarrow X, E^+ R 49358.15 Z Schumann-Runge b.	
A 3E^+	35397.8	799.07	Z 12.16 ^a	0.9106	0.0141 ₆ ^a	4.7 ^b	1.5215	(A \rightarrow b) ^c (21886) (A \rightarrow a) ^c (27125) A \rightarrow X, de R 35007.1 ₅ Z Hertzberg I b.	(16)* (22)* (89)*
A $^+$ 3D_u	(34690) ^f	(850) ^g	(20) ^g	(0.96) ^h	(0.026 ₂) ^h		(1.48)	(A \rightarrow a) ^c (26440) A \rightarrow X, I^+ R (34320) ^g Hertzberg III b.	(19)*
c 1E^-	33057.3	794.2 ₉	Z 12.73 ^k	0.915 ₅	0.0139 ₁ ^k	[7.4]	1.517 ₄	c \rightarrow a, I^+ (24782) c \rightarrow X, R 32664.1 Z Hertzberg II b.	(188) (19)* (87)
b 1E^+	13195.1	1432.77 ^a	Z 14.00 ^a	1.4003 ^a	0.01820 ^a	5.351 ^b	1.22888	b \rightarrow a, C 5238.5 b \rightarrow X, de R 13120.91 ^f Z Atmospheric oxygen b.	(40) (12)*
a 1E_g	7918.1	[1483.5 ₀]	Z (12.9)	1.4264	0.0171	[4.86]	1.215 ₅	a \rightarrow X, he R 7882.39 Z IR atmos. oxygen b.	(10)*
X 3E^-	0	1580.19 ₃	Z 11.98 ⁱ	[1.4376766] ^j $\frac{1}{2} + \frac{1}{2}$ = 1.4456	0.0159 ₃ ^{kl}	[4.839] ^{jl}	1.20752	Rot. vibr. sp. (collision induced) Rotation sp. ^{mn} Spin reorientation (fine structure) sp. ^{no} Raman sp. ^p EPR sp.	(12a)(75a) (142) (94)(105) (20)(41)(76) (120)(159) (38)* (124)* (152)(183)* (25)(138)(154)

State	T_e	w_e	$w_e x_e$	B_e	α_e	D_e (10^{-6}cm^{-1})	r_e (\AA)	Observed Transitions Design. ν_{00}	References
C₂									
		$\mu = 6.00000000$		$D_0^0 = 6.21 \text{ eV}^a$		$I.P. = 12.15 \text{ eV}^b$			JUL 1976
		Theoretical work and potential functions (16)(29)(35)(49).							
F 1D_u	[75456.9]	[1557.5]	Z	1.645	0.019	6	1.307	F \rightarrow X, R 74532.9 Z	(51)*
E 3D_g	[73183.6] ^c	[1458.06]	Z	1.5238	0.0170	6.6	1.3579	E \rightarrow a, R 71649.6 Z	(51)*
f 3E^-	71045.8	1360.5	Z 14.8	1.448 ^d	0.040 ^e	10	1.393	f \rightarrow a, R 70188.4 Z	(51)*
E 1E^+	55034.7	1671.5 ₀	Z 40.02 ^f	1.7897	0.0387 ^g	8.3 ^h	1.2529	E \rightarrow A, I^+ V 46668.3 Z	(10)*
D 1E^-	43239.4 ₄	1829.57	Z 13.94	1.8332 ^j	0.0196	7.32 ^j	1.2380	D \rightarrow X, I^+ 43226.7 ₄ ^j Z Mulliken b.	(2)* (11) (46)
e 3D_g	40796.65 ⁱ	1106.56	Z 39.260 ⁿ	1.1922	0.0242	6.3 ⁿ	1.5351	e \rightarrow a, O R 39806.46 Z Fox-Hertzberg b.	(7)*
C 1E_g		Preliminary constants from perturbations in C 1E_g see (40).							
c 1E_g	34261.3	1809.1	Z 15.81 ^p	1.783 ₆	0.0180 ^p	6.8	1.2552	C \rightarrow A, Q V _R 25969.19 Z Deelandres-d'Azambuja b.	(13)(3)(8) (50)
d 3E_g	20022.50 ^f	1788.22	Z 16.440 ^a	1.7527 ^t	0.01608 ^u	6.7 ₄ ^v	1.2661	d \rightarrow a, V V _R 19378.44 Z Swan b.	(6)* (25) (42)* (45)
c 3E^+	13312.1	1961.6 ^y	13.7	1.87 ^y			1.2	(d \rightarrow X) ^z	
a 1E_u	8391.00	1608.35	Z 12.07 ^z	1.6163 ₄ ^{a*}	0.0168 ₆ ^{a*}	6.44 ^t	1.3184 ₃	A \rightarrow X, b^+ R 8268.16 Z Phillips b.	(18)
b 3E^-	6434.2 ₇	1470.4 ₅	Z 11.19 ^{c*}	1.4985 ^{d*}	0.0163 ₄ ^{a*}	6.22	1.3592 ₈	b \rightarrow a, f^+ R 5532.7 Z Bellis-Ransay b.	(17)*
a 3E_u	716.2 ₄ ^{e*}	1641.35	Z 11.67	1.6324 ₆ ^{h*}	0.01661	6.44	1.3119 ₀		
i 1E^-	0	1854.71	Z 13.30 ^{1*}	1.8198 ₄	0.0176 ₅ ^{1*}	6.921 ^{1*}	1.2425 ₃		

State	T_e	w_e	w_e^{vib}	B_e	α_e	D_e (10^{-6}cm^{-1})	r_e (Å)	Observed Transitions		References
								Design.	v_{00}	
N₂										
B' $^3\Sigma_u^-$	66272.4 ₇	1516.88	Z 12.18 ^a	1.473 ₃ ^b	0.0166 ₆ ^c	(5.56)	1.278 ₄	B' \rightarrow B, R "r" bands, d	6545.5 (Z)	(32)(36) (182)
W $^3\Delta_u$	59808	1501.4	(Z) 11.6					B' \leftarrow X, ^e R Ogawa-Tanaka- Wilkinson b.	65852.35 Z	(30)* (35) (66)* (149) (155)
B $^3\Delta_g$	59619.3 ₅ ^e	1733.39	Z 14.122 ^h	1.6374 ₅ ⁱ	0.0179 ₁ ^j	[5.9]	1.2126 ₀	W \leftarrow B, R, V 73 Wu-Beneesch b. W \leftarrow X, ^f R 59380 Saun-Beneesch b.	9552.0 ₃ Z	(102)(124) (131)(157) (123)* (155)
A $^3\Sigma_u^+$	50203.6 ₃	1460.64	Z 13.87 ₂ ⁿ	1.4546 ^o	0.0180 ^p	[6.1 ₃]	1.2866	B' \leftarrow A, ^f V 1st pos. gr.	59306.81 Z	(29)(196)
X $^1\Sigma_g^+$	0	2358.57	Z 14.324 ^r	1.99824 ₁ ^s	0.017318 ^m	[5.76]	1.09768 ₅	B \leftarrow X, ^m R Wilkinson b. A' \leftarrow X, ^m R 49754.78 Z Vegard-Kaplan b.	49754.78 Z	(40) (29)(70)(85)
								Rot.-vibr. ^t and rot. sp. ⁱ - pressure induced - el. field induced Raman spectra ^u Mol. beam magn. reson. ^v		(14)(59)(63) (86)(135) (141)(181) (185) (20)(134) (167)

State	T_e	w_e	$w_e x_e$	B_e	α_e	D_e (10^{-6}cm^{-1})	r_e (Å)	Observed Transitions		References
								Design.	v_{00}	
NO										
C $^2\Delta_g$ 3p π	52126 ^a	2395 ^b	15 ^b	2.000 ^{bc}	0.030 ^b		1.062	C \rightarrow A, d	8172	(45)(117)* (187)*
								C \leftarrow X, e, f v 8 bands	52251 52371	(1)* (3)(4)* (6)(30)(42)* (90)(117)* (144)
b ($^4\Gamma^-$)	(48680)	1206 ^e	H 15					b \rightarrow a, v	10395 10375 10350 10325 10300 10272	(18)*(19)*
B $^2\Delta_g$	45942.6 ^h 45913.6	1039.8 ⁱ 1037.2 ⁱ	Z 8.3 ₂ ^j Z 7.7 ₀ ^k	1.152 ⁱ 1.0921 ^l	0.012 0.012	4.9	1.416 ₇	B \leftarrow X, n f R 8 bands	45392.1 ^o Z 45481.7 ^o Z	(1)*(2)*(6)* (3)(15)(18)* (24)(31)(37) (42)* (73)* (52)(90)(127)
A $^2\Gamma_g^+$ 3s f	43965.7	2374.31	Z 16.106 ^p	1.9965 ^q ^m	0.01915 ^q	5.4	1.0634	A \leftarrow X, u v f v 7 bands (a \rightarrow X) ^x M bands	44080.5 Z 44200.2 ^w Z (38000)	(1)* (3)(8) (24)(37)(46) (127) (48)(66)
a ($^4\Delta_1$)	(38440)	1017	H 11							
X $^2\Delta_g$	119.82 ^a 0	1904.04 ₀ 1904.20 ₄	Z 14.100 ^b Z 14.075 ^b	[1.72016] ^c [1.67195] ^{cd}	0.0182 0.0171	[10.2 ₃] ^c [0.34 ₃] ^c	1.15077	$\frac{1}{2} \leftarrow \frac{1}{2}$, * Rotation-vibration sp., 4 \rightarrow 2	119.73 ^f Z	(148) (63)

State	T_e	w_e	w_e^{ex}	B_e	α_e	D_e (10^{-6}cm^{-1})	r_e (Å)	Observed Transitions		References
								Design.	v_{00}	
CO										
D $^1\Delta$	65928	1094.0	10.20	1.257	0.017		1.399	D \rightarrow X, R	65391 ^m	(43)* (96) (109)
I $^1\Gamma^-$	65084.4 ₀	1092.22	Z 10.70 ₄ ^b	1.2705 ^c	0.01848 ^d	D ₂ = 9.0	1.3911	I' \leftarrow X, R	64546.2 ₆ Z	(39)* (95) (105)*
A $^1\Delta$	65075.7 ₇	1518.2 ₄	Z 19.40 ^f	1.6115 ^e	0.0232 ₅ ^h	7.33 ⁱ	1.2353	A' \leftarrow X, ^{kl} R	64748.48 ^m Z 4th positive gr.	(63)* (109)*
e $^3\Gamma^-$	64230.2 ₄	1117.7 ₂	Z 10.686 ⁿ	1.2836 ^o	0.01753 ^p	6.77 ^q	1.3840	e \rightarrow a, ^r R	15231.6	(19)* (26)
								e \rightarrow X, R	63704.8 ₅ Z	(23)* (96) (105)*
d $^3\Delta_1$	61120.1 ^s	1171.94	Z 10.635 ^t	1.3108 ^u	0.01782 ^v	6.59 ^q	1.3696	d' \rightarrow a, ^r R	12148.7	(13)(29)(76)
								Triplet b.		
								d \rightarrow X, R	60621.9 ^t Z	(71)(109)*
a' $^3\Gamma^+$	55825.4 ₉	1228.60	Z 10.468 ^x	1.3446 ^y	0.0189 ₂ ^z	6.41 ^q	1.3523	a' \rightarrow a, ^r R	6882.4	(1)(11)(14)
								Assund b.		
								a' \rightarrow X, ^r R	55355.6 Z	(23)* (96) (105)*
								Hopfield-Birge b.		
a $^3\Delta_g$	48686.70 ^a	1743.4 ₁	14.3 ₆ ^b	1.69124 ^c	0.01904 ^d	6.36 ^e	1.20574	f \leftarrow X, ^g R	48473.22 ^b Z	(20)* (104) (109)*
X $^1\Gamma^+$	0	2169.81358	Z 13.28831 ^h	1.93128087 ⁱ	0.01750441 ^j	6.1214 ₇ ^k	1.12832 ^l	Rot.-vibr. sp. ^m , 3-0 2-0 ^o 1-0 ^p Rotation sp. ^r Far IR sp. ^q Microwave sp. Mol. beat el. reson. ^t Mol. beam magn. reson. ^u		(45)(48)(125) (135) (26)(41) (25) (24)(40)(137) (141) (45)(54)