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ULTRAVIOLET ABSORPTION CROSS-SECTIONS OF SOME
CARBONYL COMPOUNDS AND THEIR TEMPERATURE DEPENDENCE.

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ABSTRACT

Ultraviolet absorption cross-section of phosgene (CCl_2O), trichloroacetylchloride ($\text{CCl}_3\text{-CClO}$) and trichloroacetaldehyde ($\text{CCl}_3\text{-CHO}$) have been measured between 170 and 320 nm for temperature ranging from 210 to 295 K with a classical double beam equipment.

These data are compared with other available determinations performed at room temperature. Photodissociation coefficients are estimated and their temperature dependence is discussed. Impact of the photodissociation on the total atmospheric destruction of these compounds is illustrated.

I. INTRODUCTION.

Chemical degradation of alternative hydrochloro-fluorocarbons in troposphere produces a series of carbonyl compounds like phosgene, halo-aldehydes or halo-ketones, which could themselves be removed in three ways : (1) Photodissociation by solar UV radiations to produce potential odd chlorine precursors,
(2) Reaction with H_2O and
(3) Reaction with OH.

In order to determine the lifetimes for the photolysis processes, an accurate knowledge of the UV absorption cross-sections is required as a function of wavelength and temperature.

II. EXPERIMENTAL.

Ultraviolet absorption cross sections of phosgene (CCl_2O), trichloroacetylchloride ($\text{CCl}_3\text{-CClO}$) and trichloroacetaldehyde ($\text{CCl}_3\text{-CHO}$) have been measured between 170 and 320 nm for temperature ranging from 210 to 295 K with a classical double beam equipment. (Gillotay et al., 1989). The purity of the three compounds is better than 99.5 % as determined by gas phase chromatography.

III. RESULTS.

Numerical values of absorption cross-sections for wavenumber intervals of 500 cm^{-1} are given in tables I-III. The absorption spectra are illustrated in Figures 1-3 for selected temperatures namely 295, 250 and 210 K and compared with other available data at 295 K.

In all cases, Beer-Lambert's law was verified for absorption ranging from 10 to 85 %. In such conditions, and according to the error budget previously published, (Simon et al., 1988), the absorption cross-sections reported here are determined with an accuracy of $\pm 2 \%$ at room temperature and of ± 3 to $\pm 4 \%$ at the lowest temperature.

Carbonyl compounds display a continuous absorption in the 170-335 nm range. The presence of two maxima and the temperature dependence observed near the maxima and for the longest wavelengths, seems to indicate that there are two continua, one corresponding to the absorption of the C-Cl bond, with a maximum around 170 nm and the other centred respectively at 240 nm for phosgene, around 260 nm for trichloroacetylchloride and at 290 nm for trichloroacetaldehyde, corresponding to the absorption of the C=O bond.

Absorption cross-sections values change with temperature by a factor, which depends on both the wavelength and the chemical composition of the compound. For each wavelength, an exponential dependence of the absorption cross-section versus temperature is clearly established, with a decrease of absorption cross-sections in the region of low absorptions (up to 80 % at 305 nm and 210 K in the case of CCl_2O) and a small increase near the maximum of absorption (up to 5 %). This effect is the most important at the lowest temperature.

Discrepancies observed between the different set of available data at room temperature have to be discussed in more details in terms of experimental conditions.

IV. DISCUSSION.

Photodissociation coefficients of the molecules have been calculated, neglecting the effects of multiple scattering, for given altitude (z), zenith angle (χ) and wavelengths intervals according to the relation :

$$J(z) = \sigma_{\lambda} q_{\lambda}(z) \quad ; \quad q_{\lambda}(z) = q_{\lambda}(\infty) e^{-\tau_{\lambda}(z)}$$

$$\tau_{\lambda}(z) = \int_z^{\infty} [n(\text{O}_2) \sigma_{\lambda}(\text{O}_2) + n(\text{O}_3) \sigma_{\lambda}(\text{O}_3) + n(\text{air}) \sigma_{\text{scatt}}] \sec \chi dz$$

where

z is the altitude,

σ are the absorption cross-sections,

$q(z)$ and $q(\infty)$ are the solar irradiance at altitude z or extraterrestrial ($z = \infty$)

n is the number of particles per volume unit.

Calculations are made for solar zenith angle of 0° and 60° ($\sec = 1$ and 2), taking into account the values of $\sigma(O_2)$ and $\sigma(O_3)$ from WMO and Kockarts (1976), σ_{scatt} from Nicolet (1984) and the values of $q(\infty)$ from WMO (1986) and by taking into account the actual values of cross-sections corresponding to the temperature conditions at each altitude.

These photodissociation coefficients are illustrated in Figures 4-6 and compared with those calculated with values of absorption cross-sections measured at room temperature.

Stratospheric photodissociation coefficients (for altitude ranging from 15 to 50 km) calculated the temperature dependent absorption cross-sections, are smaller than those calculated with the room temperature values in the 20-35 km region, due to the decrease in the absorption cross-sections in the 200 nm region and the influence of the wavelengths longer than 280nm in the low stratosphere.

Tropospheric photodissociation coefficients for phosgene are very low (between 10^{-9} and $10^{-11} \text{ sec}^{-1}$) and are reduced down to 20 % of their room temperature values, using the temperature dependent cross-sections. For the two other compounds, tropospheric photodissociation coefficients are relatively high (between 10^{-7} and 10^{-4} sec^{-1}) and show a small temperature dependence. Photolysis is, for these two molecules, a non negligible mechanism for their tropospheric removal.

In conclusion, this work presents a new set of experimental data on the absorption cross-sections of carbonyl compounds in atmospheric temperature condition and highlights a non negligible temperature dependence of their photolysis.

V. REFERENCE.

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Table I PHOSGENE (CCl_2O)

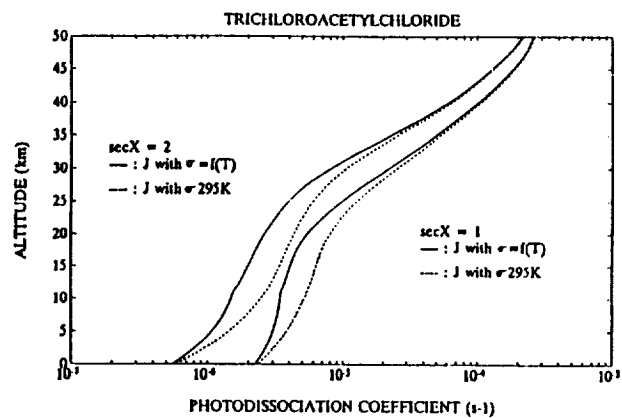
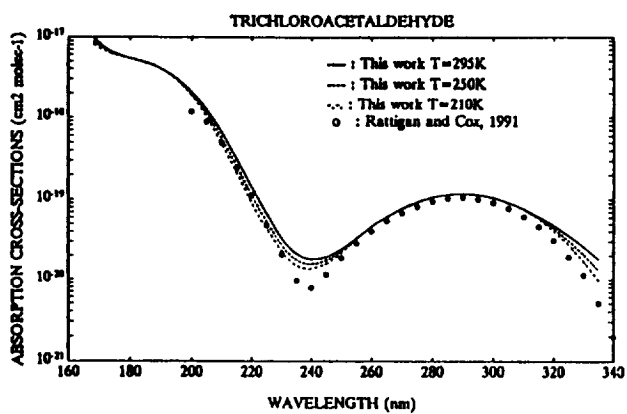
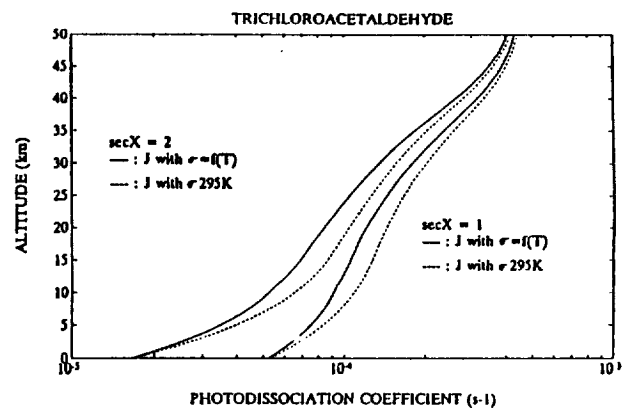
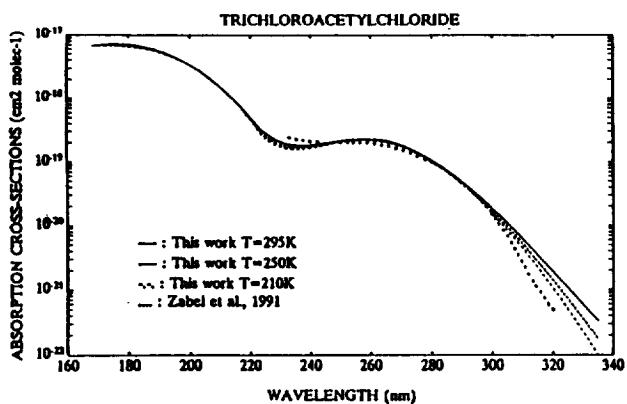
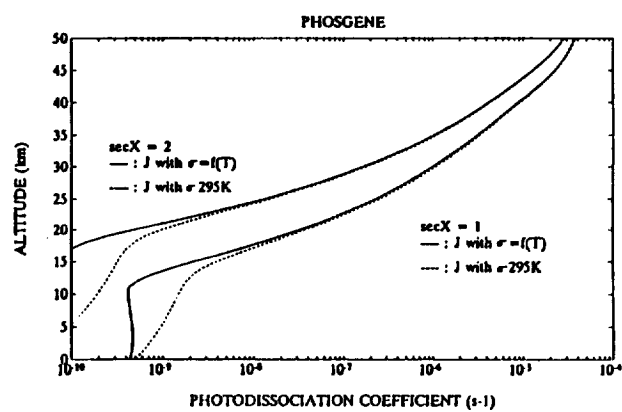
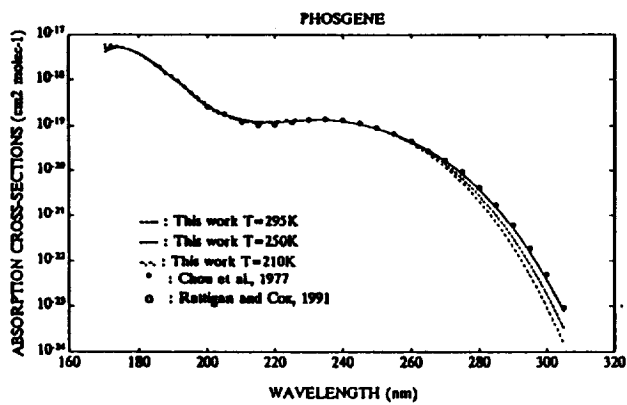
N°	(nm)	$\sigma(\lambda) \times 10^{21} (\text{cm}^2 \text{ molec}^{-1})$				
		295K	270K	250K	230K	210K
42	164.7-169.5	3010	2600	4160	4610	5560
43	169.5-172.4	4330	4760	5130	5530	5960
44	172.4-173.9	4930	3160	5360	5560	5770
45	173.2-175.4	5030	3190	5240	5330	5430
46	175.4-177.0	4890	4810	4920	4940	4930
47	177.0-178.6	4540	4490	4460	4420	4280
48	178.6-180.2	4040	3960	3900	3850	3790
49	180.2-181.8	3470	3390	3320	3270	3160
50	181.8-183.5	2890	2820	2760	2710	2660
51	183.5-185.2	2340	2280	2240	2200	2150
52	185.2-186.9	1860	1820	1790	1760	1730
53	186.9-188.7	1460	1440	1410	1390	1376
54	188.7-190.5	1160	1140	1120	1110	1090
55	190.5-192.3	903	888	876	865	854
56	192.3-194.2	715	698	685	673	660
57	194.2-196.1	524	509	498	487	476
58	196.1-198.0	399	386	376	367	357
59	198.0-200.0	312	302	293	285	277
60	200.0-202.0	252	243	236	229	223
61	202.0-204.1	209	202	196	191	185
62	204.1-206.2	179	173	168	163	159
63	206.2-208.3	158	153	149	145	141
64	208.3-210.5	143	139	136	132	129
65	201.9-212.8	133	129	127	124	121
66	212.8-215.0	126	124	121	119	117
67	215.0-217.4	123	121	119	117	115
68	217.4-219.8	122	120	119	117	116
69	219.8-221.2	122	121	120	119	118
70	222.2-224.7	124	124	123	123	122
71	224.2-227.3	127	127	127	127	127
72	227.3-229.9	131	131	131	131	132
73	229.9-232.6	134	134	135	135	135
74	232.6-235.3	136	137	137	137	137
75	235.3-238.1	131	132	132	133	133
76	238.1-241.0	125	125	126	126	127
77	241.0-243.9	116	116	117	117	118
78	243.9-246.9	105	104	104	106	104
79	246.9-250.0	93.1	92.9	92.8	92.7	92.5
80	250.0-253.2	80.5	79.8	78.3	78.8	78.2
81	253.2-256.4	66.5	65.3	64.4	63.4	62.5
82	256.4-259.7	53.4	51.8	50.6	49.4	48.2
83	259.7-263.2	40.3	38.4	37.0	35.7	34.4
84	263.2-266.7	28.9	27.1	25.8	24.9	23.2
85	266.7-270.3	19.8	18.1	16.9	15.8	14.7
86	270.3-274.0	12.6	11.3	10.3	9.45	8.64
87	274.0-277.8	7.34	6.39	5.71	5.10	4.55
88	277.8-281.7	3.92	3.29	2.86	2.48	2.16
89	281.7-285.7	1.93	1.56	1.31	1.11	0.933
90	285.7-289.9	0.831	0.643	0.524	0.426	0.347
91	289.9-294.1	0.316	0.233	0.193	0.143	0.112
92	294.1-298.5	0.105	0.0732	0.0569	0.0413	0.0310
93	298.5-303.0	0.0318	0.0190	0.0136	0.00970	0.00694
94	303.0-307.7	0.00491	0.00298	0.00200	0.00134	0.000899

Table III TRICHLOROACETYLCHLORIDE (CCl₃CClO)

N°	$\epsilon(\lambda) \times 10^{21} \text{ (cm}^2 \text{ molec.}^{-1}\text{)}$				
	(nm)	295K	270K	250K	210K
42	166.7-169.5	6710	6710	6700	6690
43	169.5-172.4	6990	6910	6840	6720
44	172.4-173.9	7160	7000	6900	6810
45	173.9-175.4	7160	7020	6880	6690
46	175.4-177.0	7170	7010	6890	6650
47	177.0-178.6	7140	6970	6840	6710
48	178.6-180.2	7060	6890	6750	6620
49	180.2-181.8	6940	6780	6640	6510
50	181.8-183.5	6790	6640	6500	6380
51	183.5-185.2	6590	6410	6300	6160
52	185.2-186.9	6390	6200	6090	5950
53	186.9-188.7	5990	5800	5700	5560
54	188.7-190.5	5670	5480	5380	5240
55	190.5-192.3	5280	5100	5000	4860
56	192.3-194.2	4870	4700	4590	4450
57	194.2-196.1	4450	4280	4170	4030
58	196.1-198.0	4020	3850	3740	3600
59	198.0-200.0	3580	3410	3300	3160
60	200.0-202.0	3140	3000	2890	2750
61	202.0-204.1	2720	2580	2470	2330
62	204.1-206.2	2310	2170	2060	1920
63	206.2-208.3	1940	1800	1690	1550
64	208.3-210.5	1600	1460	1350	1210
65	210.5-212.8	1290	1150	1040	900
66	212.8-215.0	1020	880	770	630
67	215.0-217.4	789	649	538	398
68	217.4-219.8	595	455	344	204
69	219.8-222.2	440	300	189	79
70	222.2-224.7	318	178	67	27
71	224.7-227.3	238	107	32	12
72	227.3-229.9	189	72	24	8
73	229.9-232.6	152	57	18	6
74	232.6-235.3	125	42	13	4
75	235.3-238.1	100	32	10	3
76	238.1-241.0	81	25	8	2
77	241.0-243.9	67	19	6	2
78	243.9-246.9	56	14	5	1
79	246.9-250.0	47	11	4	1
80	250.0-253.2	39	8	3	1
81	253.2-256.4	32	6	2	1
82	256.4-259.7	26	5	2	1
83	259.7-263.2	22	4	1	1
84	263.2-266.7	19	3	1	1
85	266.7-270.3	16	2	1	1
86	270.3-274.0	13	1	1	1
87	274.0-277.8	10	1	1	1
88	277.8-281.7	7	1	1	1
89	281.7-285.7	5	1	1	1
90	285.7-289.9	4	1	1	1
91	289.9-294.1	3	1	1	1
92	294.1-298.5	2	1	1	1
93	298.5-303.0	1	1	1	1
94	303.0-307.7	1	1	1	1
95	307.7-312.5	1	1	1	1
96	312.5-317.5	1	1	1	1
97	317.5-322.5	1	1	1	1
98	322.5-327.5	1	1	1	1
99	327.5-332.5	1	1	1	1
100	332.5-337.5	1	1	1	1

Table II TRICHLOROACETALDEHYDE (CCl₃CHO)

N°	$\epsilon(\lambda) \times 10^{21} \text{ (cm}^2 \text{ molec.}^{-1}\text{)}$				
	(nm)	295K	270K	250K	210K
42	166.7-169.5	9880	9210	8880	8670
43	169.5-172.4	7850	7440	7110	6900
44	172.4-173.9	6890	6500	6170	5960
45	173.9-175.4	6420	6030	5700	5490
46	175.4-177.0	6040	5650	5320	5110
47	177.0-178.6	5720	5330	5000	4790
48	178.6-180.2	5480	5090	4760	4550
49	180.2-181.8	5250	4860	4530	4320
50	181.8-183.5	5030	4640	4310	4100
51	183.5-185.2	4820	4430	4100	3890
52	185.2-186.9	4590	4200	3870	3660
53	186.9-188.7	4370	3980	3650	3440
54	188.7-190.5	4150	3760	3430	3220
55	190.5-192.3	3740	3350	3020	2810
56	192.3-194.2	3400	2990	2660	2450
57	194.2-196.1	3030	2620	2290	2080
58	196.1-198.0	2650	2240	1900	1690
59	198.0-200.0	2270	1860	1520	1310
60	200.0-202.0	1890	1480	1140	930
61	202.0-204.1	1540	1130	780	570
62	204.1-206.2	1210	800	450	240
63	206.2-208.3	936	520	170	60
64	208.3-210.5	711	300	80	20
65	210.5-212.8	511	180	40	10
66	212.8-215.0	365	100	20	5
67	215.0-217.4	265	50	10	2
68	217.4-219.8	177	20	5	1
69	219.8-222.2	121	10	2	1
70	222.2-224.7	81	5	1	1
71	224.7-227.3	57	3	1	1
72	227.3-229.9	39	2	1	1
73	229.9-232.6	28	1	1	1
74	232.6-235.3	22	1	1	1
75	235.3-238.1	19	1	1	1
76	238.1-241.0	17	1	1	1
77	241.0-243.9	16	1	1	1
78	243.9-246.9	15	1	1	1
79	246.9-250.0	14	1	1	1
80	250.0-253.2	13	1	1	1
81	253.2-256.4	12	1	1	1
82	256.4-259.7	11	1	1	1
83	259.7-263.2	10	1	1	1
84	263.2-266.7	9	1	1	1
85	266.7-270.3	8	1	1	1
86	270.3-274.0	7	1	1	1
87	274.0-277.8	6	1	1	1
88	277.8-281.7	5	1	1	1
89	281.7-285.7	4	1	1	1
90	285.7-289.9	3	1	1	1
91	289.9-294.1	2	1	1	1
92	294.1-298.5	1	1	1	1
93	298.5-303.0	1	1	1	1
94	303.0-307.7	1	1	1	1
95	307.7-312.5	1	1	1	1
96	312.5-317.5	1	1	1	1
97	317.5-322.5	1	1	1	1
98	322.5-327.5	1	1	1	1
99	327.5-332.5	1	1	1	1
100	332.5-337.5	1	1	1	1



Figures 1-3 : Ultraviolet absorption cross-sections at 295 K, 250 K and 210 K.

Figures 4-6 : Photodissociation coefficients as a function of altitude.