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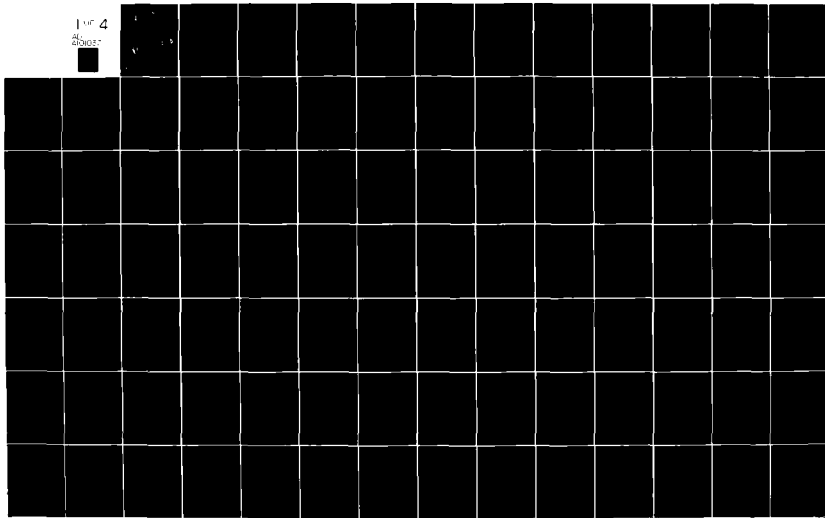
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TECHNICAL REPORT RH-81-4

COMPILATION OF ATOMIC AND MOLECULAR DATA RELEVANT TO
GAS LASERS

VOLUME VIII

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US Army Missile Laboratory

December 1980

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U.S. ARMY MISSILE COMMAND

Redstone Arsenal, Alabama 35809

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ABSTRACT (CONTINUED)

Species Index for Volumes I-V of the Compilation of Data Relevant to Gas Lasers", September 1979. These six volumes, authored by E.W. McDaniel and other personnel at Georgia Tech, Georgia State University, the Joint Institute of Laboratory Astrophysics (JILA), and the Army Missile Command (MICOM), were published as MIRADCOM Technical Report H-78-1 at Redstone Arsenal, Alabama.

Volumes I and II were prepared in the context of the two most-used techniques for gas laser pumping: electrical discharges and high intensity high energy electron and ion beams. Heavy emphasis was placed on the rare gases and halogens (atoms, molecules, and ions), and the rare gas-halides, although a significant amount of material on other species was included. Volumes III, IV, and V contain much information relevant to electrical discharges and high intensity, high energy electron and ion beams, but are oriented toward a third pumping technique: nuclear pumping. Since nuclear reactions may also become interesting in some form of hybrid laser where the excitation and ionization produced by the reaction products might be used to supply electrons for an electrical discharge laser or an initiator for a pulsed chemical laser, or as an initiator and sustainer for a continuous wave (CW) chemical laser; data relevant to these systems was also included.

The present volumes serve to update most of the areas covered in the previous documents. Those areas not treated here are considered to have been adequately dealt with earlier, as far as immediate data needs are concerned. However, even in those areas where new data are not presented here, references are given to past volumes in order to facilitate access to the previous data. Another function of the present work is to expand somewhat the scope of our data coverage, both with respect to atomic and molecular structural properties and with respect to atomic collisions. New species and sets of collision partners that have recently assumed importance are treated here, and other systems that may become important in the gas laser context are given attention. A significant amount of new material is also added to the chapter on surface impact phenomena, partly because of current interest in hollow-cathode lasers.

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SECURITY CLASSIFICATION OF THIS PAGE(When Data Entered)

PREFACE

This volume and the succeeding volume are the seventh and the eighth in a series that presents data relevant to research and development in the field of gas lasers. Volumes I and II are entitled, "Compilation of Data Relevant to Rare Gas-Rare Gas and Rare Gas-Monohalide Excimer Lasers", December 1977. Volumes III, IV, and V comprise a "Compilation of Data Relevant to Nuclear-Pumped Lasers", December 1978. Volume VI provides a "Cumulative Reactant Species Index for Volumes I-V of the Compilation of Data Relevant to Gas Lasers", September 1979. These six volumes, authored by E.W. McDaniel and other personnel at Georgia Tech, Georgia State University, the Joint Institute of Laboratory Astrophysics (JILA), and the Army Missile Command (MICOM), were published as MIRADCOM Technical Report H-78-1 at Redstone Arsenal, Alabama.

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The present volumes serve to update most of the areas covered in the previous documents. Those areas not treated here are considered to have been adequately dealt with earlier, as far as immediate data needs are concerned. Such areas include all nuclear processes, and atomic collisions occurring at "high" energies, i.e., above about 100 eV impact energy. However, even in those areas where new data are not presented here, references are given to past volumes in order to facilitate access to the previous data. Attention should also be called to another document that may prove useful to those requiring data--"Bibliography: Sources of Information on Phenomena of Interest in Gas Laser Research and Development", Technical Report RH-77-1, by E.W. McDaniel, H.W. Ellis, F.L. Eisele, and M.G. Thackston, January 1977, US Army Missile Command, Redstone Arsenal, Alabama. A second, updated edition of this bibliography will be published early in 1981.

Another function of the present volume is to expand somewhat the scope of our data coverage, both with respect to atomic and molecular structural properties and with respect to atomic collisions (by the

12 + 14 → latter term, we mean two- and three- body collisions between electrons, ions, atoms, molecules, and photons at impact energies sufficiently low that nuclear forces are unimportant). New species and sets of collision partners that have recently assumed importance are treated here, and other systems that may become important in the gas laser context are given attention. A significant amount of new material is also added to the chapter on surface impact phenomena, partly because of current interest in hollow-cathode lasers.

In conclusion, we wish to thank C.F. Barnett, former Director of the Controlled Fusion Atomic Data Center at the Oak Ridge National Laboratory, and E.C. Beaty, Chief of the Information Center at JILA, for their cooperation and the use of their facilities. In certain areas, our work would have been immensely more difficult without their assistance. Chapter D on photon collision processes in gases was put together with the aid of several scientists. Particularly significant were the contributions of Dr. Joseph Berkowitz, of Argonne National Laboratory, whose book Photoabsorption, Photoionization, and Photoelectron Spectroscopy (Academic Press, New York, 1979) provided us with a wealth of references and critically evaluated data on atoms and molecules. We gratefully acknowledge being allowed access to the manuscript prior to publication, as well as Dr. Berkowitz providing us with a number of large-size versions of figures from his book. In addition, we acknowledge the contributions of Professor C.E. Brion, of the University of British Columbia, for providing us with a complete set of reprints, spanning a decade, of his very extensive work on partial and total cross sections of atoms and molecules. Also, the expert help of Professor H.W. Ellis, of Eckerd College, St Petersburg, Florida, on the transport properties of electrons, ions, and neutrals in gases is gratefully acknowledged.

C. ELECTRON-HEAVY PARTICLE COLLISIONS

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The views, opinions, and/or findings contained in this report are those of the authors and should not be construed as an official Department of the Army position, policy, or decision, unless so designated by other documentation.

C-1. ELECTRON SCATTERING: ELASTIC, TOTAL, AND MOMENTUM TRANSFER

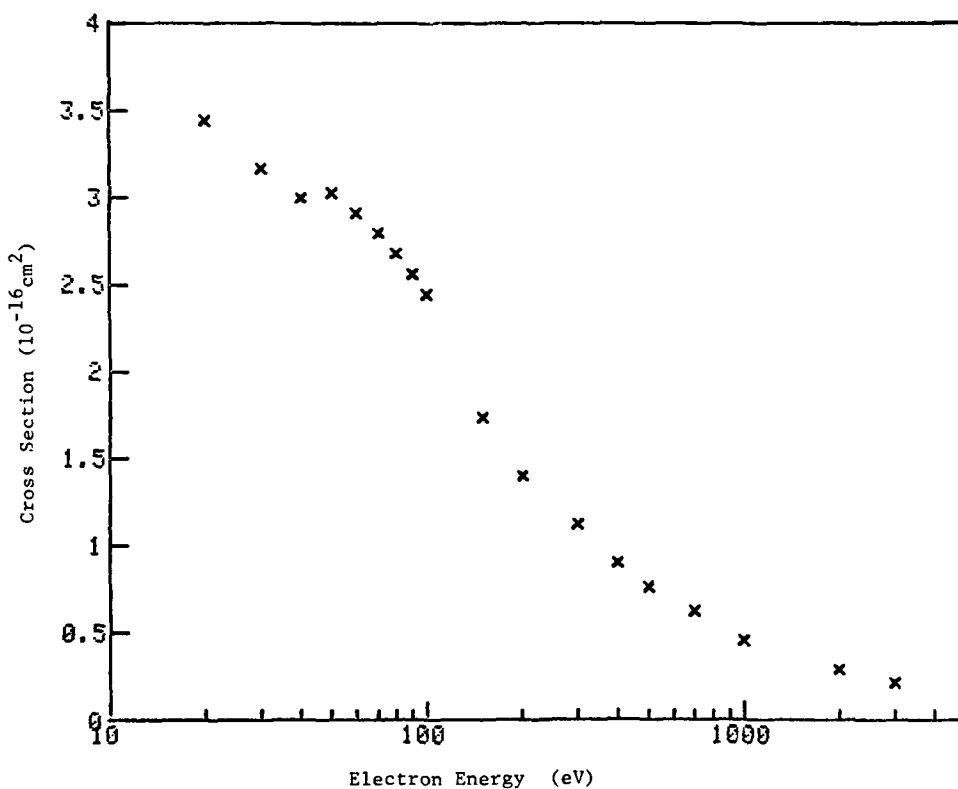
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Tabular and Graphical Data C-1.1. Semi-empirical (average) cross sections
for elastic scattering of electrons in Ne.

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
20	3.444	200	1.398
30	3.170	300	1.120
40	2.999	400	0.9003
50	3.030	500	0.7620
60	2.912	700	0.6191
70	2.797	1000	0.4556
80	2.680	2000	0.2856
90	2.563	3000	0.2095
100	2.447		
150	1.735		

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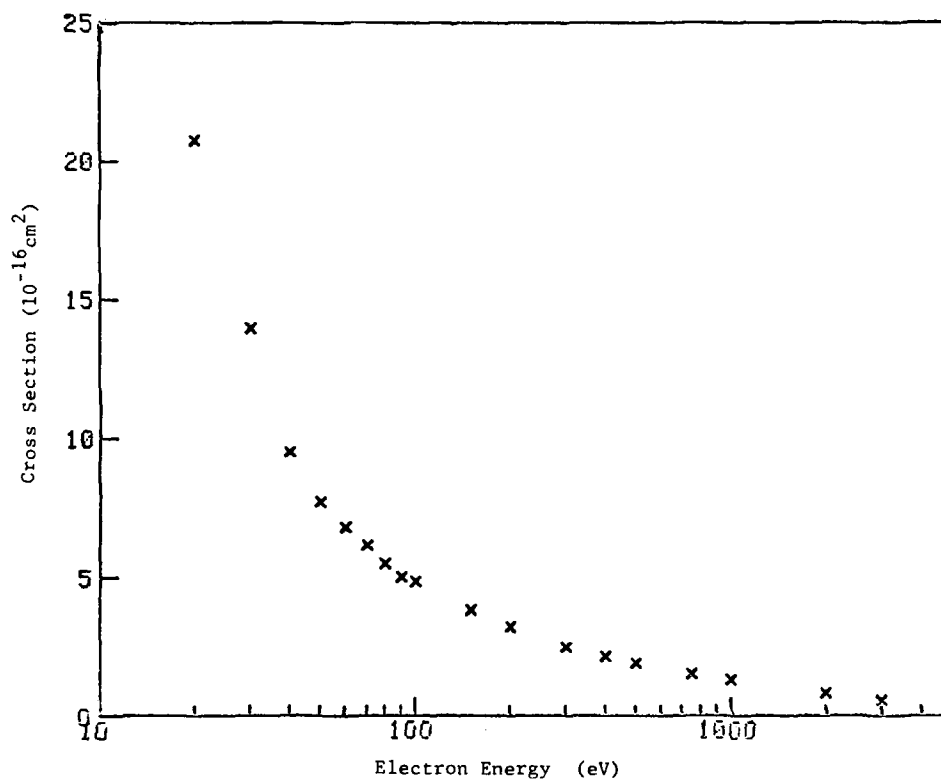


Reference: F. J. de Heer, R. H. J. Jansen, W. van der Kaay, J. Phys. B 12,
979 (1979).

Tabular and Graphical Data C-1.2. Semi-empirical (average) cross sections
for elastic scattering of electrons in Ar.

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
20	20.76	200	3.201
30	13.95	300	2.467
40	9.510	400	2.116
50	7.743	500	1.886
60	6.799	750	1.493
70	6.147	1000	1.274
80	5.489	2000	0.8037
90	5.010	3000	0.5741
100	4.861		
150	3.794		

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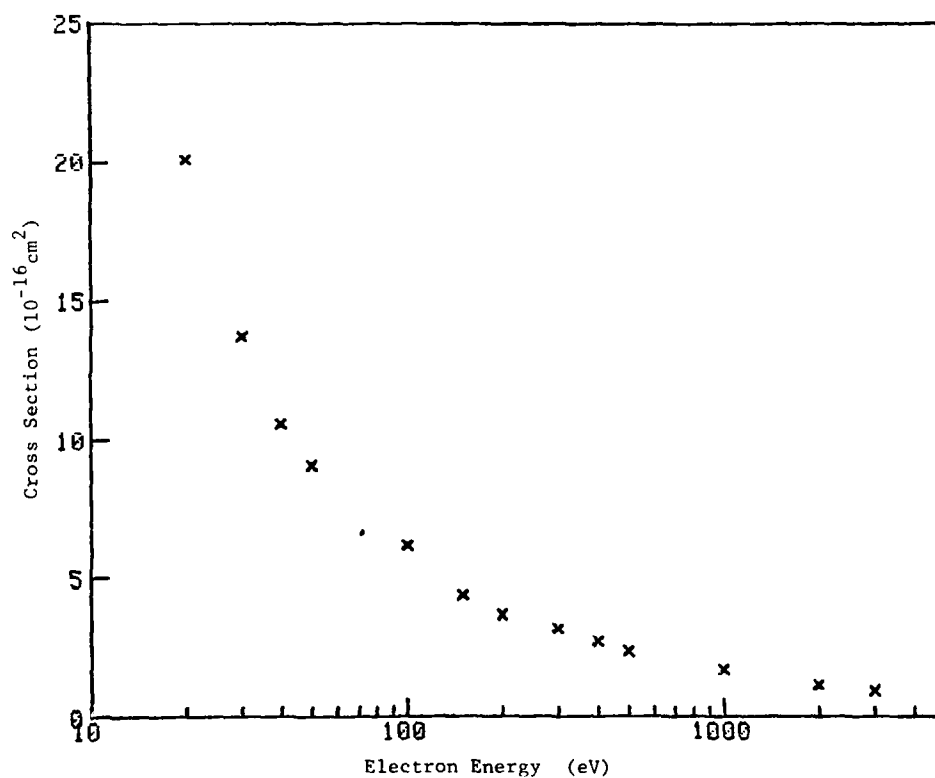


Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12,
979 (1979).

Tabular and Graphical Data C-1.3. Semi-empirical (average) cross sections
for elastic scattering of electrons in Kr.

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
20	20.11	300	3.184
30	13.74	400	2.721
40	10.58	500	2.373
50	9.090	1000	1.695
100	6.169	2000	1.126
150	4.591	3000	0.9417
200	3.677		

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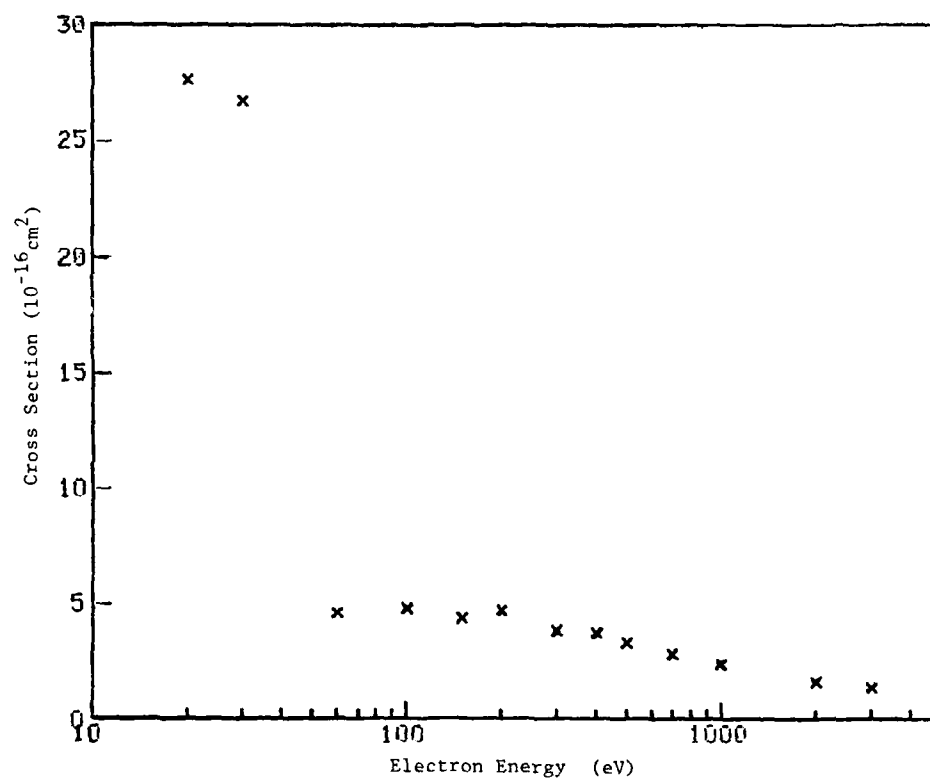


Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12,
979 (1979).

Tabular and Graphical Data C-1.4. Semi-empirical (average) cross sections
for elastic scattering of electrons in Xe.

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
20	27.66	400	3.752
30	26.75	500	3.338
60	4.587	700	2.834
100	4.825	1000	2.395
150	4.385	2000	1.649
200	4.724	3000	1.379
300	3.834		

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Reference : F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12,
979 (1979).

Tabular Data C-1. 5. Cross sections for elastic scattering
for electrons incident on Hg.

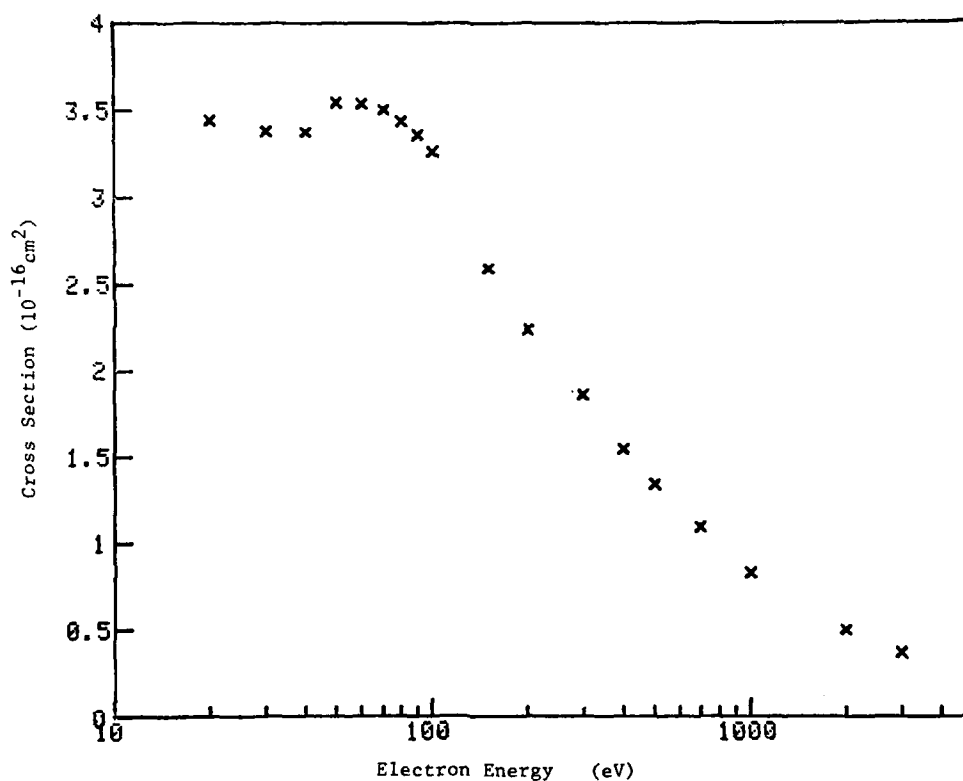
Electron Energy eV	Cross Section 10^{-16} cm^2
300	4.76
400	3.86
500	3.33

Reference: K. Jost, and B. Ohnemus, Phys. Rev. A 19,
611 (1979).

Tabular and Graphical Data C-1.6. Semi-empirical total scattering cross sections
for electrons incident on Ne.

Electron Energy eV	Cross Section 10^{-16}cm^2	Electron Energy eV	Cross Section 10^{-16}cm^2
20	3.444	200	2.234
30	3.360	300	1.854
40	3.377	400	1.544
50	3.548	500	1.336
60	3.540	700	1.088
70	3.506	1000	0.8236
80	3.439	2000	0.4979
90	3.360	3000	0.3632
100	3.265		
150	2.590		

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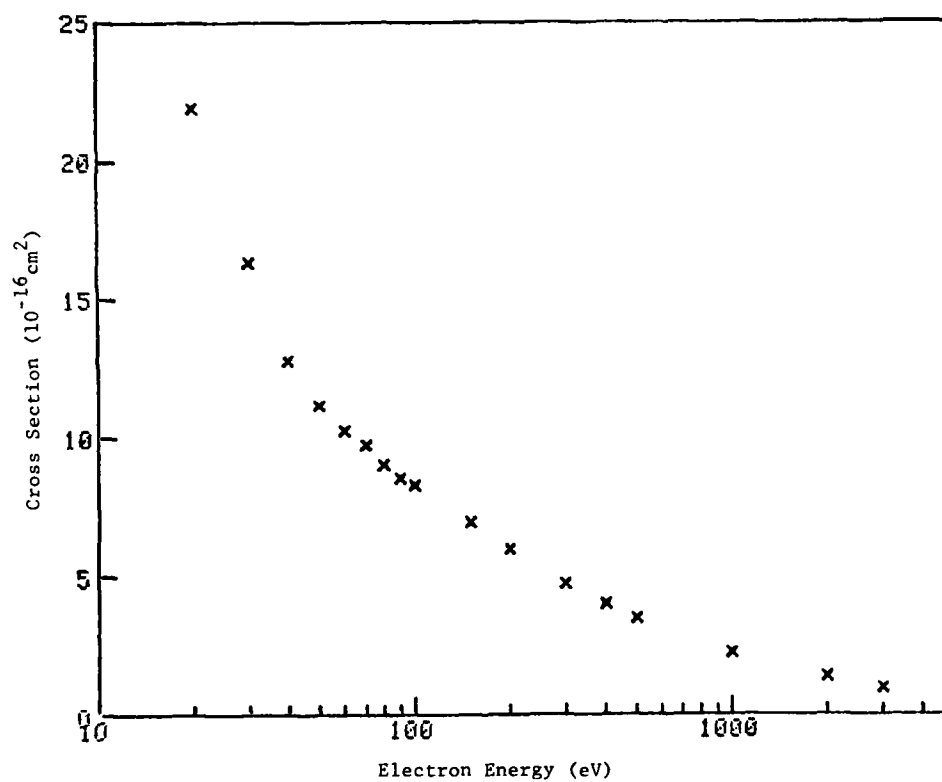


Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12,
979 (1979).

Tabular and Graphical Data C-1.7. Semi-empirical total scattering cross sections for electrons incident on Ar.

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
20	21.92	150	6.931
30	16.33	200	5.981
40	12.76	300	4.707
50	11.15	400	3.996
60	10.25	500	3.503
70	9.706	1000	2.229
80	9.003	2000	1.350
90	8.538	3000	0.9717
100	8.278		

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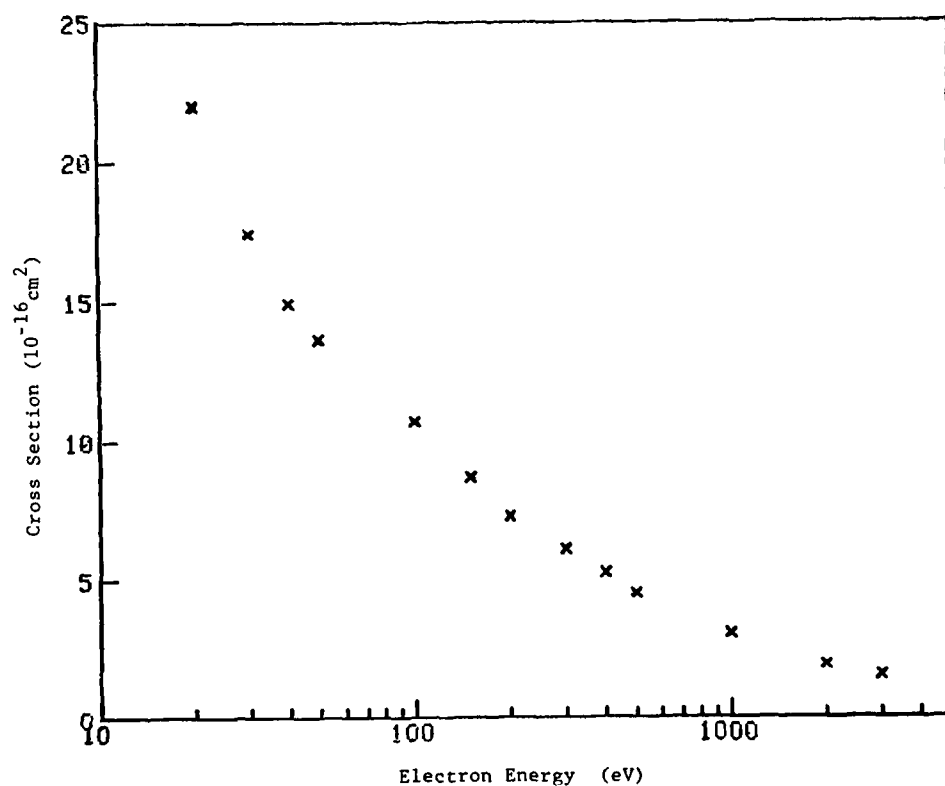


Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12, 979 (1979).

Tabular and Graphical Data C-1.8. Semi-empirical total scattering cross sections for electrons incident on Kr.

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16}cm^2	eV	10^{-16}cm^2
20	22.00	300	6.096
30	17.42	400	5.234
40	14.92	500	4.497
50	13.64	1000	3.050
100	10.71	2000	1.912
150	8.684	3000	1.517
200	7.298		

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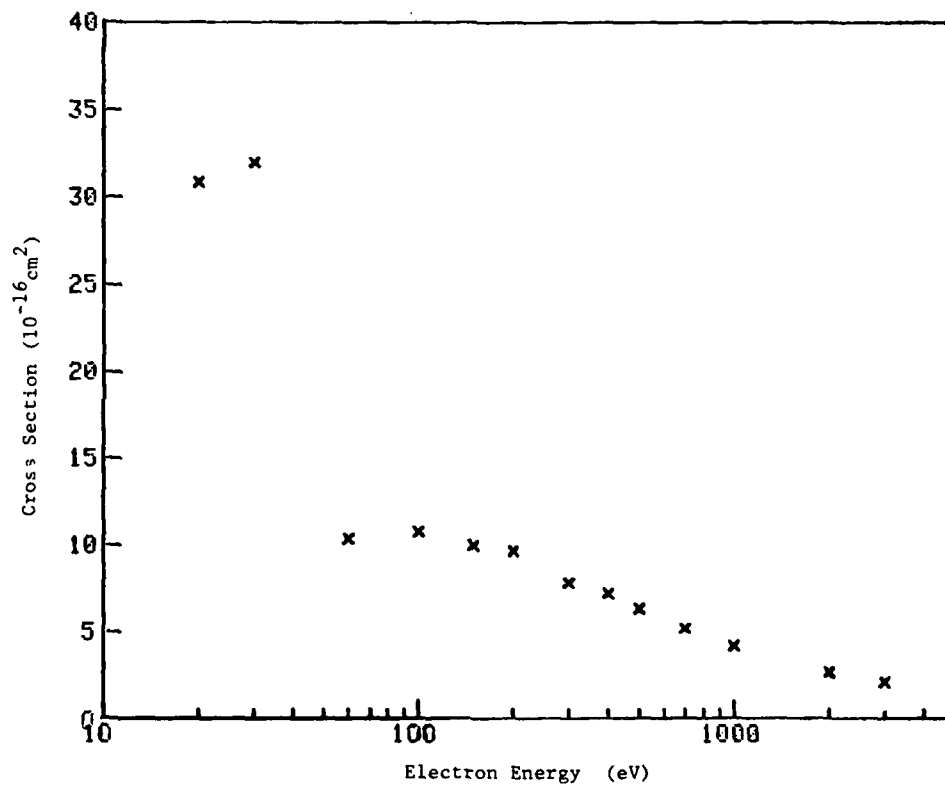


Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B. 12, 979 (1979).

Tabular and Graphical Data C-1.9. Semi-empirical total scattering cross
sections for electrons incident on Xe.

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16}cm^2	eV	10^{-16}cm^2
20	30.86	400	7.158
30	32.01	500	6.261
60	10.32	700	5.153
100	10.67	1000	4.167
150	9.944	2000	2.686
200	9.588	3000	2.132
300	7.785		

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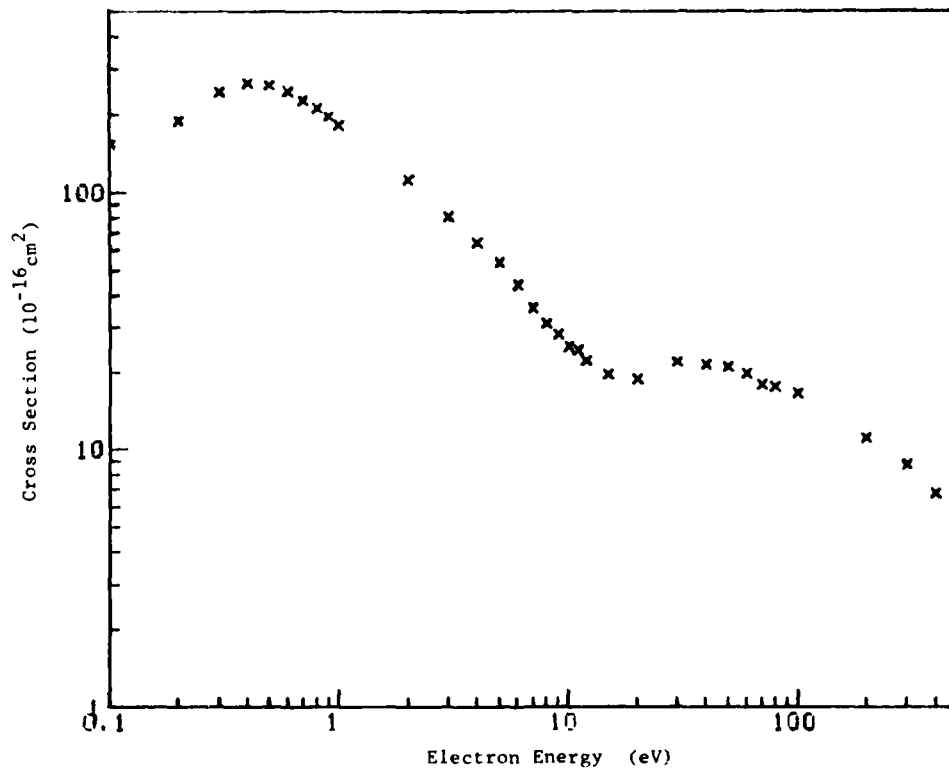


Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12,
979 (1979).

Tabular and Graphical Data C-1.10. Cross sections for total scattering of
electrons incident on Hg.

Electron Energy	Cross Section	Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2	eV	10^{-16} cm^2
0.10	153	4.0	63.6	40	21.5
0.20	189	5.0	53.6	50	21.1
0.30	246	6.0	43.9	60	19.8
0.40	265	7.0	35.6	70	17.9
0.50	261	8.0	30.9	80	17.5
0.60	248	9.0	27.9	100	16.5
0.70	228	10.0	25.2	200	11.1
0.80	213	11	24.3	300	8.76
0.90	197	12	22.2	400	6.75
1.00	184	15	19.5	500	5.80
2.0	113	20	13.8		
3.0	81.1	30	21.9		

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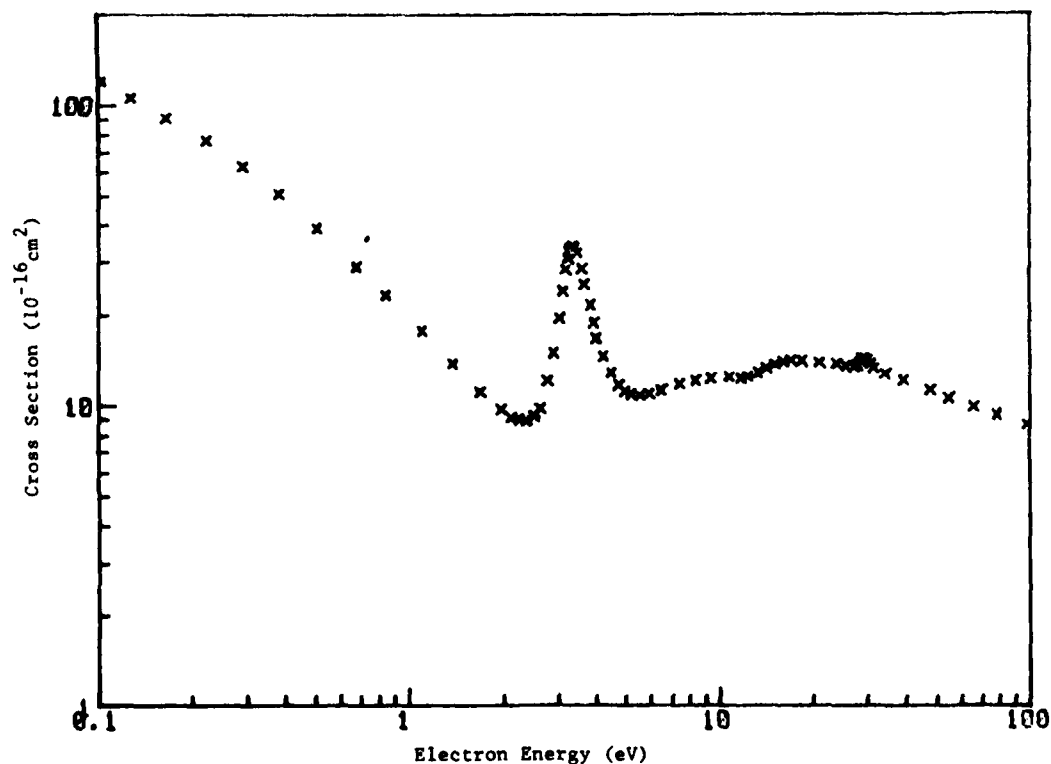


Reference: K. Jost and B. Ohnemus, Phys. Rev. A 19 611 (1979).

Tabular and Graphical Data C-1.11. Calculated cross sections for total scattering of electrons incident on CO₂.

Electron Energy eV	Cross Section 10 ⁻¹⁶ cm ²	Electron Energy eV	Cross Section 10 ⁻¹⁶ cm ²	Electron Energy eV	Cross Section 10 ⁻¹⁶ cm ²
0.101	120	3.37	33.6	17.0	14.2
0.127	105	3.42	33.4	18.5	14.1
0.164	91.2	3.47	32.2	21.0	13.9
0.222	76.8	3.60	28.5	23.9	13.7
0.291	62.7	3.66	25.2	25.5	13.5
0.380	50.6	3.83	21.5	27.0	13.4
0.505	38.8	3.94	18.8	27.8	13.6
0.675	29.0	4.02	16.7	28.3	13.9
0.841	23.3	4.22	14.5	28.7	14.2
1.10	17.8	4.47	12.8	29.3	14.4
1.38	13.8	4.73	11.7	29.9	14.1
1.69	11.1	4.95	11.1	30.6	13.7
1.97	9.70	5.19	10.9	31.4	13.3
2.13	9.17	5.52	10.8	34.2	12.7
2.26	8.97	5.93	10.9	39.2	12.1
2.36	8.88	6.45	11.2	47.6	11.2
2.51	9.23	7.47	11.8	54.6	10.6
2.64	9.83	8.34	12.1	65.7	9.82
2.79	12.1	9.37	12.3	78.2	9.24
2.92	14.9	10.8	12.4	98.1	8.58
3.05	19.4	11.8	12.3		
3.13	24.0	12.3	12.5		
3.19	28.3	13.3	12.9		
3.24	30.9	14.2	13.3		
3.26	32.3	15.1	13.7		
3.31	33.1	16.0	14.0		

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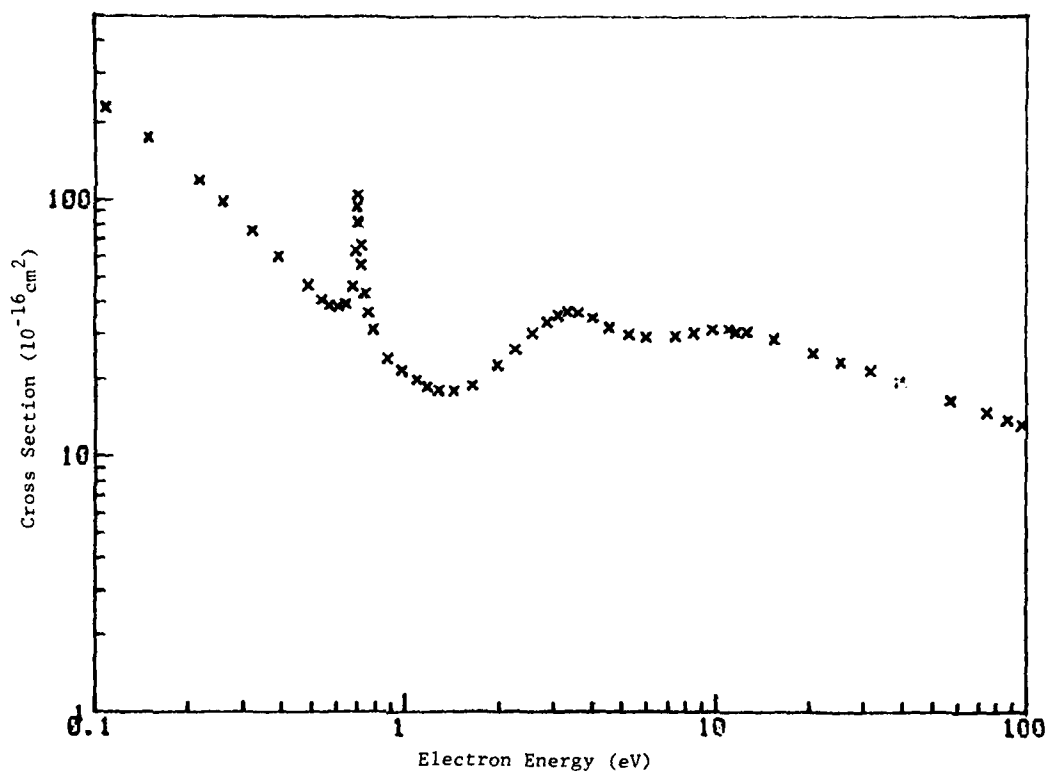
Reference: M. G. Lynch, D. Dill, J. Siegel, and J. L. Dehmer, J. Chem. Phys. 71, 4249 (1979).

Tabular and Graphical Data C-1.12. Calculated cross sections for total scattering of electrons incident on OCS.

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
0.0739	304	0.718	55.4	4.56	31.8
0.108	228	0.743	43.3	5.25	29.9
0.148	174	0.758	36.4	6.00	29.0
0.215	119	0.788	31.3	7.47	29.4
0.259	97.5	0.881	24.2	8.55	30.3
0.321	75.6	0.976	21.7	9.76	31.3
0.388	60.1	1.09	19.9	11.1	31.3
0.482	46.3	1.18	18.7	11.7	30.5
0.537	40.6	1.28	18.1	12.7	30.6
0.569	38.7	1.43	17.8	15.5	28.4
0.605	38.4	1.64	18.9	20.7	25.2
0.640	39.3	1.98	22.7	25.4	23.5
0.674	46.0	2.26	26.3	31.7	21.7
0.691	62.9	2.58	30.3	39.4	19.4
0.701	82.2	2.87	33.4	56.9	16.4
0.701	94.2	3.10	35.6	74.9	14.8
0.702	104	3.33	36.7	87.3	13.9
0.708	81.0	3.61	36.4	96.6	13.3
0.718	66.5	4.01	34.6		

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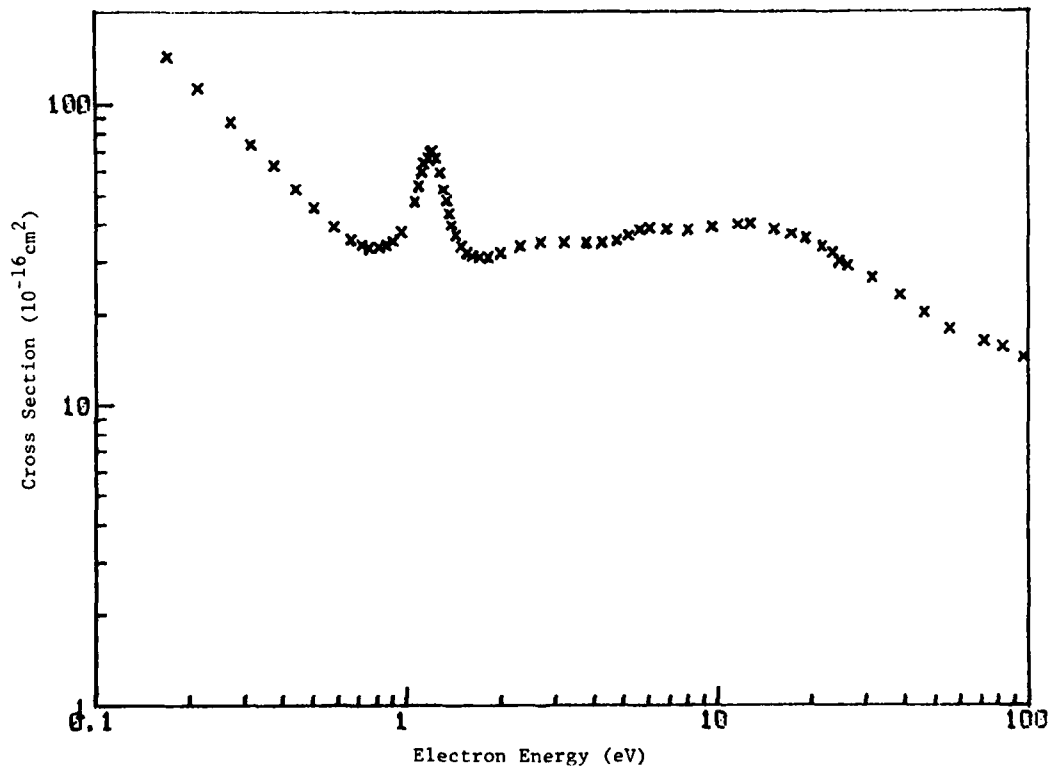


Reference: M. G. Lynch, D. Dill, J. Siegel, and J. L. Dehmer, J. Chem. Phys. 71, 4249 (1979).

Tabular and Graphical Data C-1.13. Calculated cross sections for total
scattering of electrons incident on CS₂.

Electron Energy	Cross Section	Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10 ⁻¹⁶ cm ²	eV	10 ⁻¹⁶ cm ²	eV	10 ⁻¹⁶ cm ²
0.170	143	1.24	66.7	6.87	38.2
0.213	112	1.28	59.3	8.04	38.2
0.272	87.1	1.31	52.3	9.59	38.9
0.316	73.7	1.34	48.3	11.7	39.8
0.373	62.8	1.36	43.6	12.7	40.0
0.441	52.4	1.39	39.8	15.2	38.2
0.502	45.5	1.44	36.7	17.3	36.9
0.581	39.4	1.50	33.7	19.3	35.8
0.662	35.5	1.57	32.1	21.8	33.4
0.717	34.0	1.63	31.4	23.5	32.0
0.758	33.4	1.71	31.0	24.6	29.9
0.817	33.5	1.83	31.0	26.2	28.9
0.858	34.0	2.01	32.1	31.5	26.3
0.907	35.0	2.32	33.8	38.6	23.1
0.958	37.7	2.70	34.4	46.3	20.2
1.06	47.7	3.23	34.5	55.7	17.7
1.09	53.9	3.80	34.6	72.4	16.2
1.12	59.8	4.28	34.6	82.8	15.5
1.14	64.0	4.76	35.1	96.6	14.3
1.17	66.3	5.17	36.6		
1.19	69.6	5.60	38.1		
1.21	69.9	6.05	38.6		

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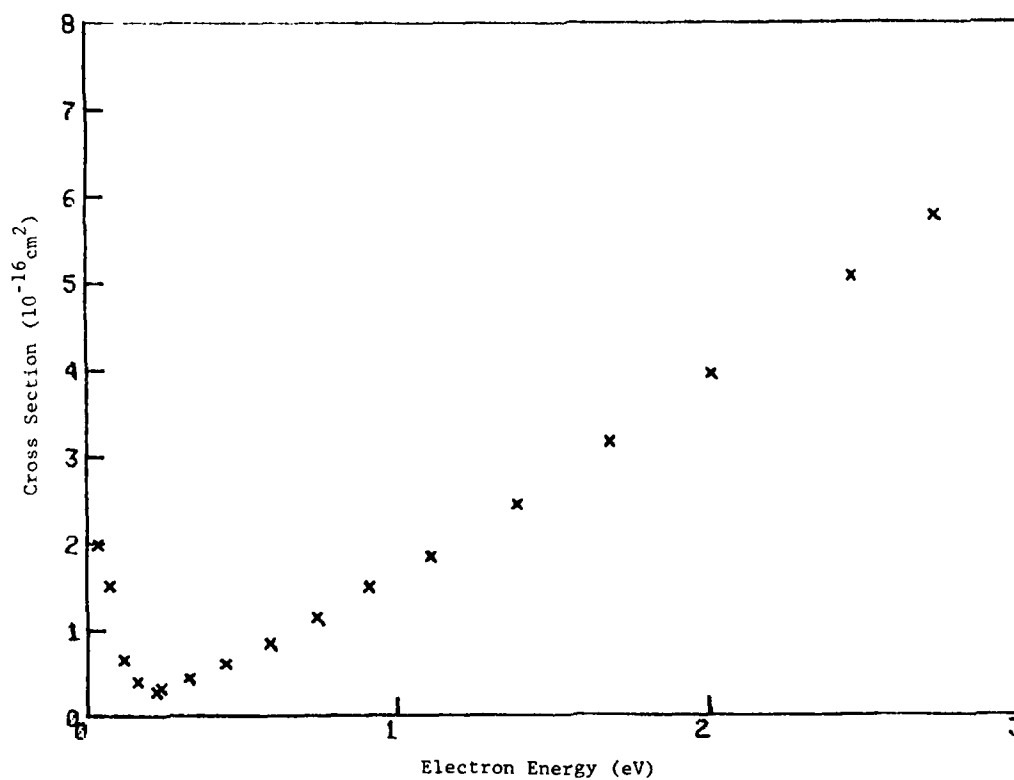


Reference: M. G. Lynch, D. Dill, J. Siegel, and J. L. Dehmer, J. Chem. Phys. 71,
4249 (1979).

Tabular and Graphical Data C-1.14. Cross sections for total scattering of
electrons incident on CH₄.
(0 - 2 eV)

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10 ⁻¹⁶ cm ²	eV	10 ⁻¹⁶ cm ²
0.036	1.98	0.91	1.49
0.073	1.52	1.1	1.85
0.12	0.651	1.4	2.45
0.16	0.397	1.7	3.16
0.22	0.282	2.0	3.95
0.24	0.318	2.5	5.08
0.33	0.443	2.7	5.77
0.44	0.610		
0.59	0.842		
0.74	1.14		

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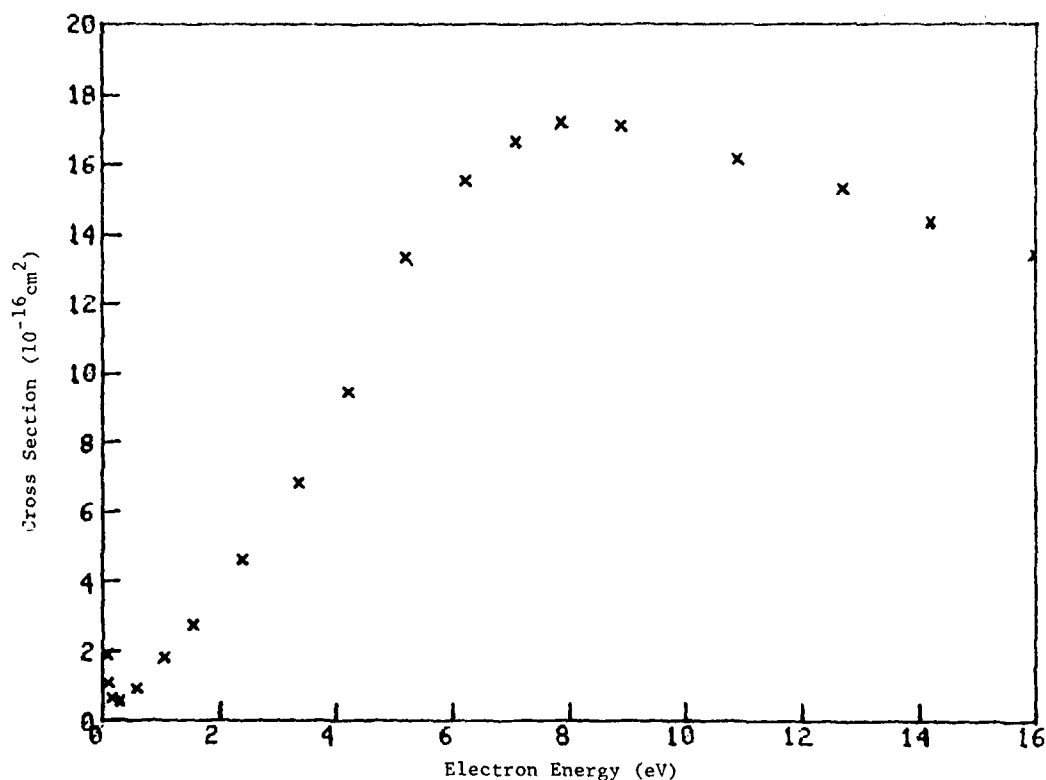


Reference: E. Barbarito, M. Basta, M. Calicchio, and G. Tessari, J. Chem. Phys. 71, 54 (1979).

Tabular and Graphical Data C-1.15. Cross sections for total scattering of
electrons incident on CH₄.
(0 - 16 eV)

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10 ⁻¹⁶ cm ²	eV	10 ⁻¹⁶ cm ²
0.081	1.9	5.2	13
0.090	1.1	6.2	16
0.15	0.64	7.1	17
0.29	0.55	7.9	17
0.58	0.92	8.9	17
1.1	1.8	11	16
1.5	2.7	13	15
2.4	4.6	14	14
3.4	6.8	16	13
4.2	9.5		

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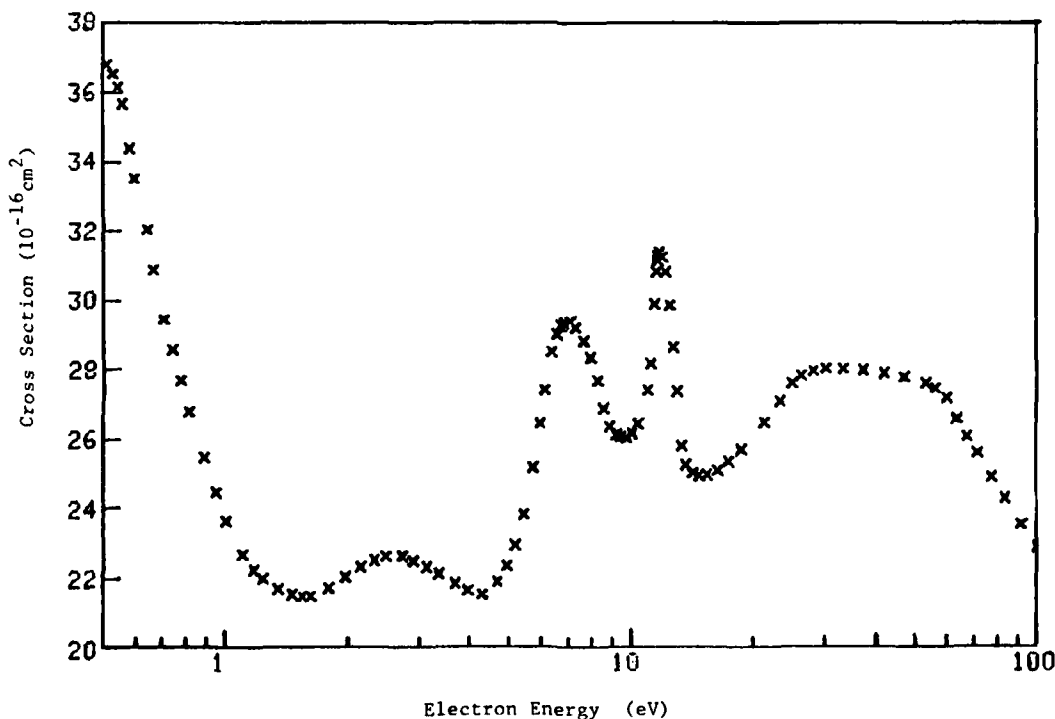
Reference: E. Barbarito, M. Basta, M. Calicchio, G. Tessari, J. Chem. Phys 71, 54 (1979).

Tabular and Graphical Data C-1.16. Cross sections for total scattering of electrons incident on SF₆.

Electron Energy eV	Cross Section 10 ⁻¹⁶ cm ²	Electron Energy eV	Cross Section 10 ⁻¹⁶ cm ²	Electron Energy eV	Cross Section 10 ⁻¹⁶ cm ²
0.508	36.8	4.31	21.5	12.5	29.8
0.528	36.5	4.67	21.9	12.8	28.6
0.545	36.1	4.93	22.3	13.0	27.3
0.558	35.6	5.18	22.9	13.4	25.8
0.580	34.4	5.46	23.8	13.7	25.2
0.597	33.5	5.77	25.2	14.2	25.0
0.640	32.0	5.99	26.4	14.8	24.9
0.665	30.9	6.12	27.4	15.4	24.9
0.708	29.4	6.39	28.5	16.4	25.1
0.743	28.5	6.59	29.0	17.5	25.3
0.777	27.7	6.76	29.2	18.8	25.7
0.815	26.7	6.86	29.3	21.4	26.4
0.886	25.4	7.10	29.3	23.3	27.0
0.948	24.5	7.32	29.2	25.0	27.6
1.01	23.6	7.66	28.8	26.3	27.8
1.10	22.6	7.96	28.3	28.1	27.9
1.17	22.2	8.26	27.7	30.2	28.0
1.24	22.0	8.58	26.9	33.2	28.0
1.35	21.7	8.91	26.3	37.3	27.9
1.46	21.5	9.27	26.1	42.0	27.9
1.54	21.5	9.49	26.0	47.0	27.7
1.62	21.5	9.78	26.0	53.2	27.6
1.81	21.7	10.1	26.1	56.1	27.4
1.98	22.0	10.4	26.4	59.7	27.1
2.16	22.3	11.0	27.4	63.8	26.5
2.33	22.5	11.2	28.1	67.5	26.0
2.49	22.6	11.5	29.9	71.4	25.6
2.73	22.6	11.6	30.8	77.7	24.9
2.90	22.5	11.6	31.1	83.7	24.2
3.14	22.3	11.7	31.2	91.4	23.5
3.36	22.1	11.8	31.3	100	22.8
3.70	21.8	12.0	31.2		
3.98	21.6	12.2	30.8		

Cont. Next Column

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Reference: R. E. Kennerly, R. A. Bonham, and M. McMillan, J. Chem. Phys. 70, 2039 (1979).

C-2. EXCITATION BY ELECTRON IMPACT

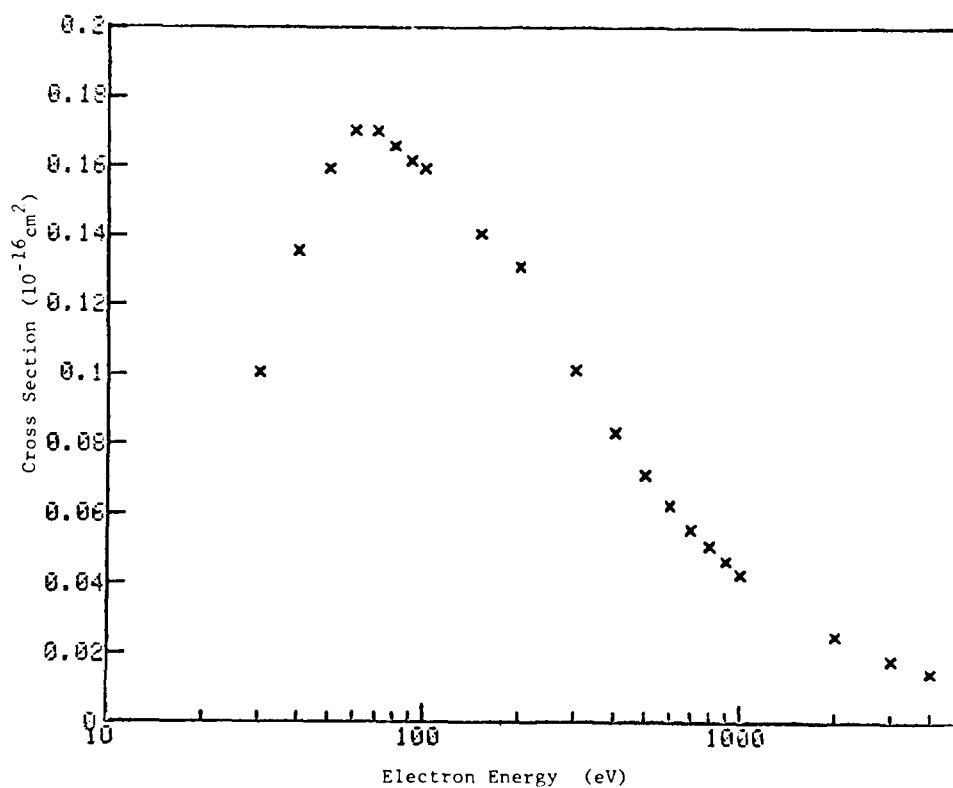
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Tabular and Graphical Data C-2.1. Semi-empirical cross sections for electron impact excitation of Ne.

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
30	0.1005	400	0.08311
40	0.1354	500	0.07110
50	0.1594	600	0.06236
60	0.1704	700	0.05545
70	0.1704	800	0.05049
80	0.1660	900	0.04618
90	0.1616	1000	0.04262
100	0.1594	2000	0.02465
150	0.1407	3000	0.01796
200	0.1309	4000	0.01421
300	0.1011		

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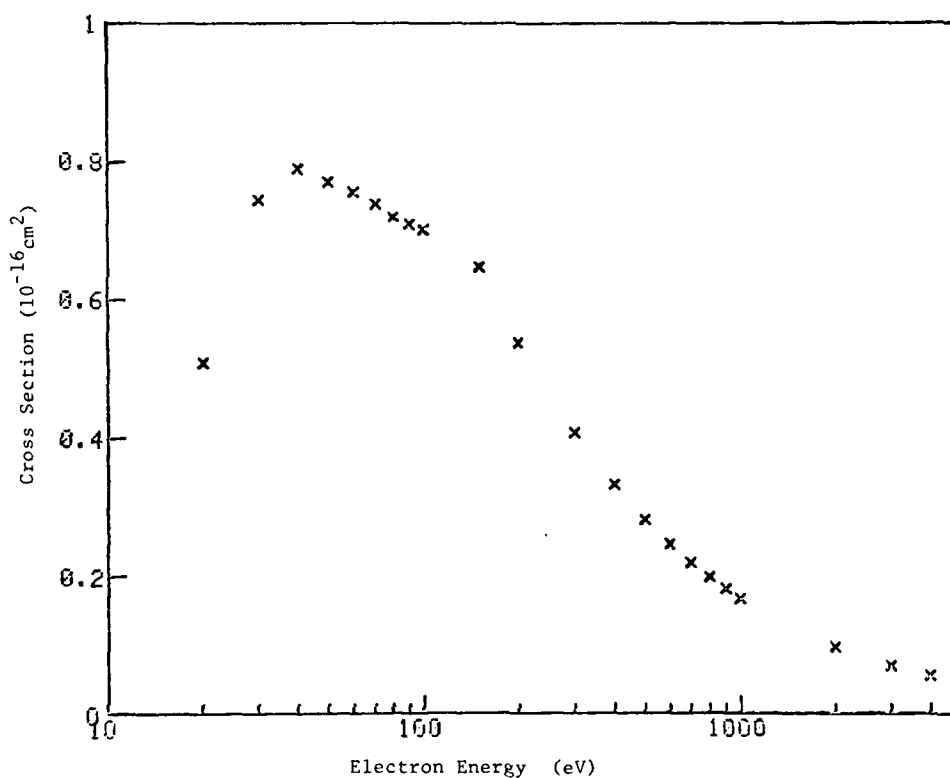


Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12, 979 (1979).

Tabular and Graphical Data C-2.2. Semi-empirical cross sections for electron impact excitation of Ar.

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
20	0.5035	400	0.3324
30	0.7446	500	0.2826
40	0.7383	600	0.2457
50	0.7701	700	0.2193
60	0.7555	800	0.1963
70	0.7373	900	0.1809
80	0.7191	1000	0.1669
90	0.7002	2000	0.09533
100	0.7009	3000	0.06917
150	0.6469	4000	0.05461
200	0.5377		
300	0.4063		

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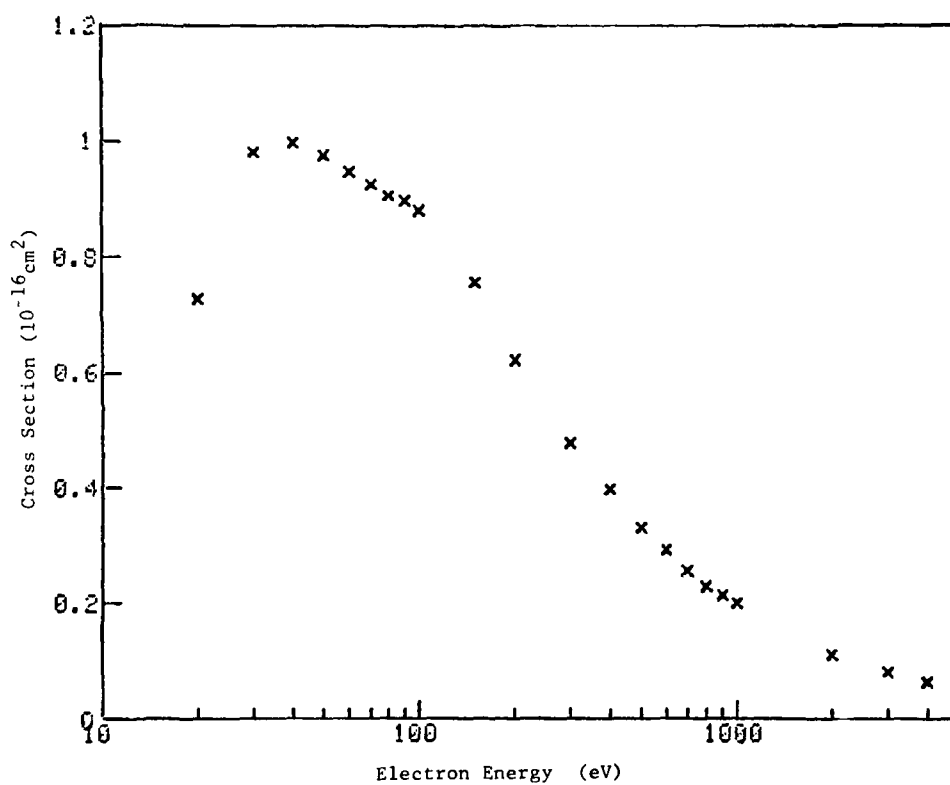


Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12, 979 (1979).

Tabular and Graphical Data C-2.3. Semi-empirical cross sections for electron impact excitation of Kr.

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
20	0.720	400	0.398
30	0.903	500	0.330
40	1.000	600	0.291
50	0.977	700	0.255
60	0.949	800	0.230
70	0.927	900	0.213
80	0.907	1000	0.199
90	0.893	2000	0.112
100	0.882	3000	0.0812
150	0.756	4000	0.0638
200	0.622		
300	0.479		

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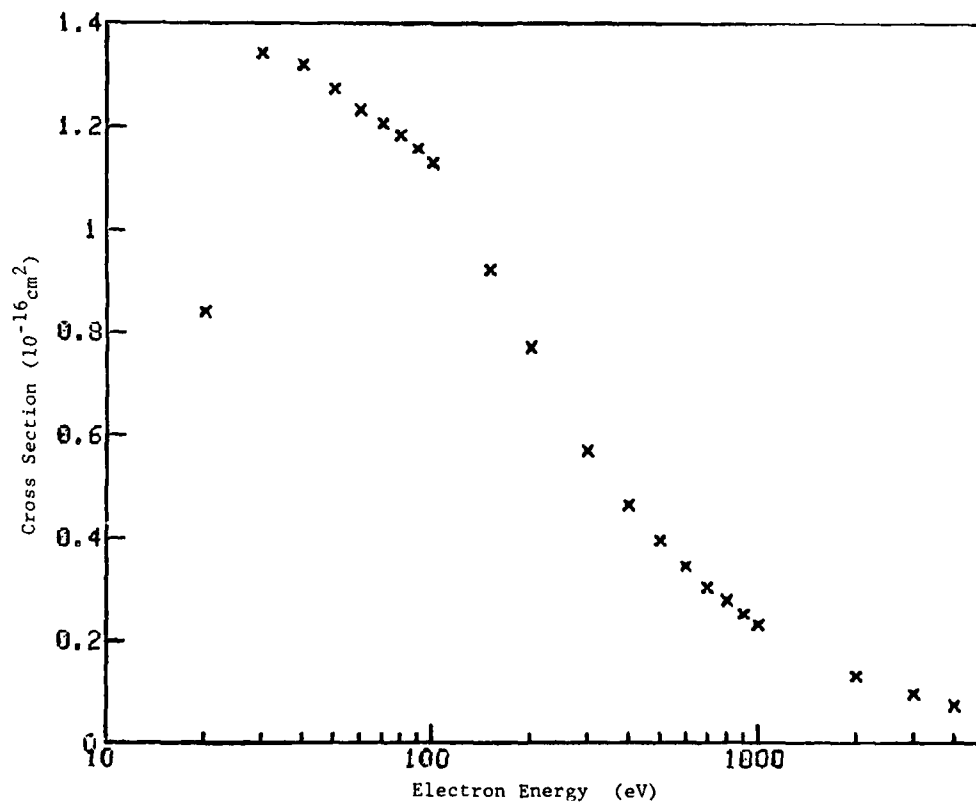


Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12, 979 (1979).

Tabular and Graphical Data C-2.4. Semi-empirical cross sections for electron impact excitation of Xe.

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
20	0.840	400	0.465
30	1.34	500	0.395
40	1.32	600	0.344
50	1.27	700	0.302
60	1.23	800	0.277
70	1.21	900	0.252
80	1.18	1000	0.231
90	1.16	2000	0.132
100	1.13	3000	0.0952
150	0.924	4000	0.0745
200	0.770		
300	0.568		

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Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12, 979 (1979).

Tabular Data C-2.5. Cross sections for electron-
impact excitation of Hg

Electron Energy	Cross Section
eV	10^{-16} cm^2
300	1.32
400	1.06
500	.90

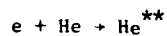
Reference: K. Jost and B. Ohnemus, Phys. Rev A 19,
611, (1979).

Tabular Data C-2.6. Cross Section for electron-
impact excitation of Hg atoms to the 6^3P_1 state.

Electron Energy	Cross Section
eV	10^{-16} cm^2
50	$.434 \pm .04$

Reference: R. D. Kaul, J. Opt. Soc. Am. 69, 150 (1979).

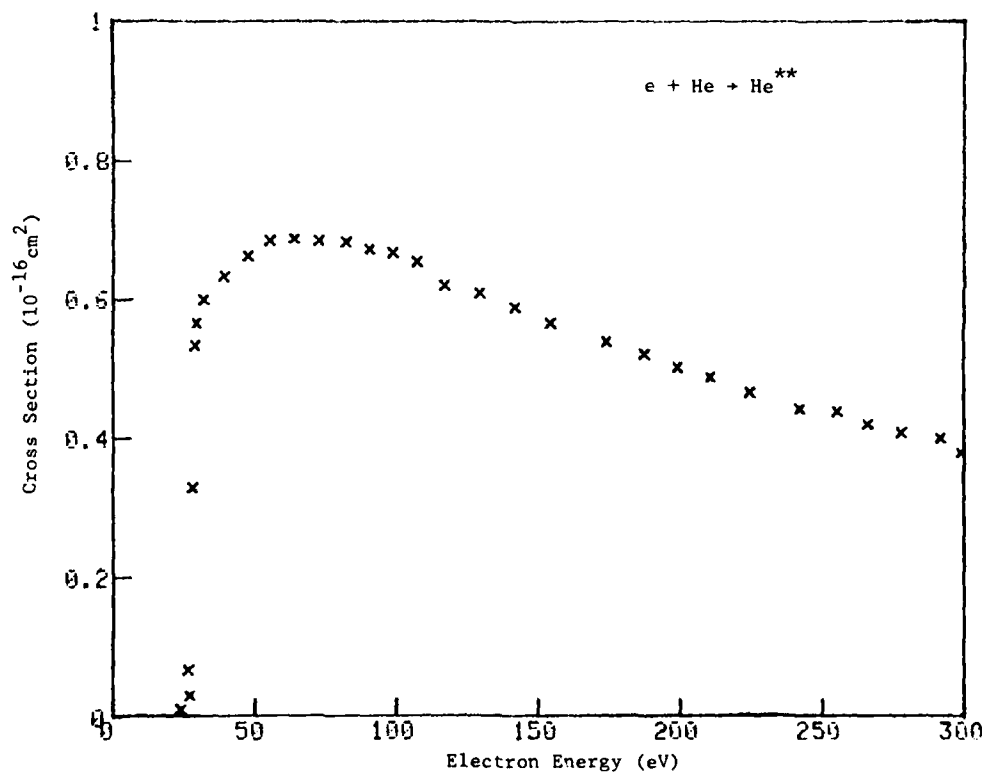
Tabular and Graphical Data C-2.7. Cross sections for electron-impact excitation of He atoms to high-Rydberg states.



Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
24	0.010	73	0.68	200	0.50
27	0.030	82	0.68	210	0.49
27	0.067	90	0.67	220	0.46
28	0.33	99	0.67	240	0.44
29	0.53	110	0.65	260	0.44
29	0.57	120	0.62	270	0.42
32	0.60	130	0.61	280	0.41
39	0.63	140	0.59	290	0.40
48	0.66	150	0.57	300	0.38
55	0.68	170	0.54		
64	0.69	190	0.52		

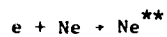
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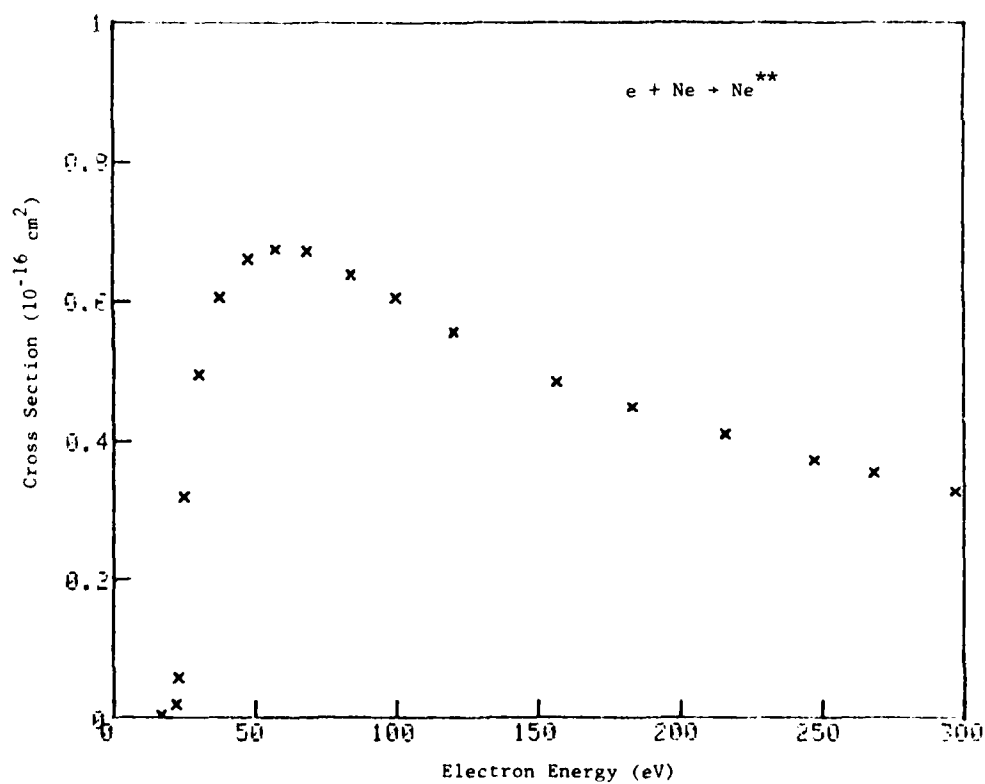
Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, Phys. Rev. A 20, 71 (1979).

Tabular and Graphical Data C-2.8. Cross sections for electron-impact excitation
of Ne atoms to high-Rydberg states



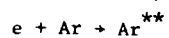
Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
17	0.0031	160	0.49
22	0.020	180	0.45
23	0.058	220	0.41
25	0.32	250	0.37
30	0.50	270	0.35
38	0.61	300	0.33
48	0.66		
57	0.67		
68	0.67		
84	0.64		
100	0.61		
120	0.56		

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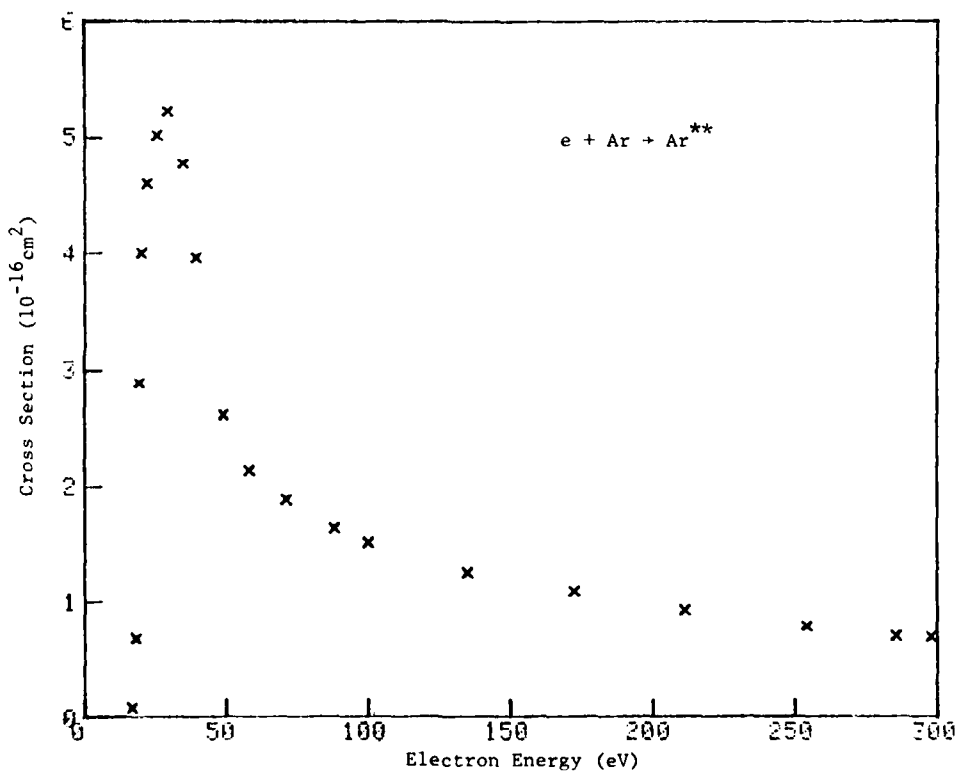
Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, Phys. Rev. A 20, 71 (1979).

Tabular and Graphical Data C-2.9. Cross sections for electron-impact excitation of Ar atoms to high-Rydberg states.



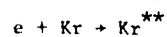
Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
17	0.077	88	1.6
18	0.68	100	1.5
19	2.9	130	1.3
20	4.0	170	1.1
22	4.6	210	0.92
25	5.0	250	0.78
29	5.2	290	0.70
35	4.8	300	0.69
39	4.0		
49	2.6		
58	2.1		
71	1.9		

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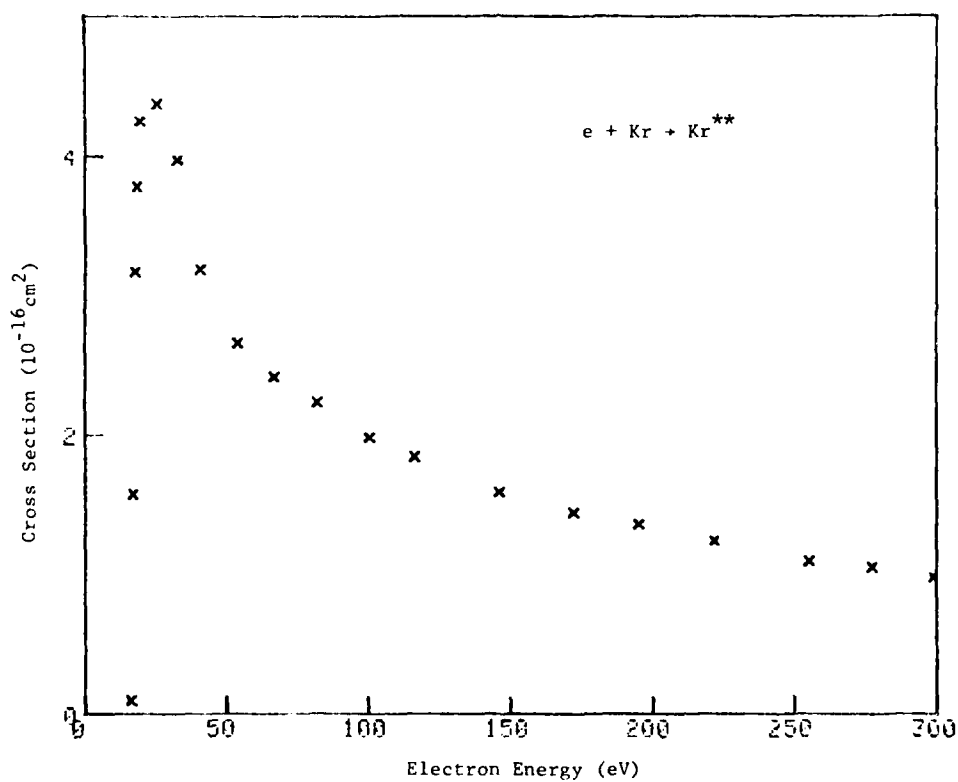
Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, Phys. Rev A 20, 71 (1979).

Tabular and Graphical Data C-2.10. Cross sections for electron-impact excitation
of Kr atoms to high-Rydberg states



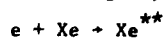
Electron Energy eV	Cross Section 10^{-16}cm^2	Electron Energy eV	Cross Section 10^{-16}cm^2
16	0.096	120	1.8
17	1.6	150	1.6
17	3.2	170	1.4
18	3.8	190	1.4
19	4.3	220	1.2
25	4.4	260	1.1
32	4.0	280	1.1
40	3.2	300	0.98
54	2.7		
66	2.4		
82	2.2		
100	2.0		

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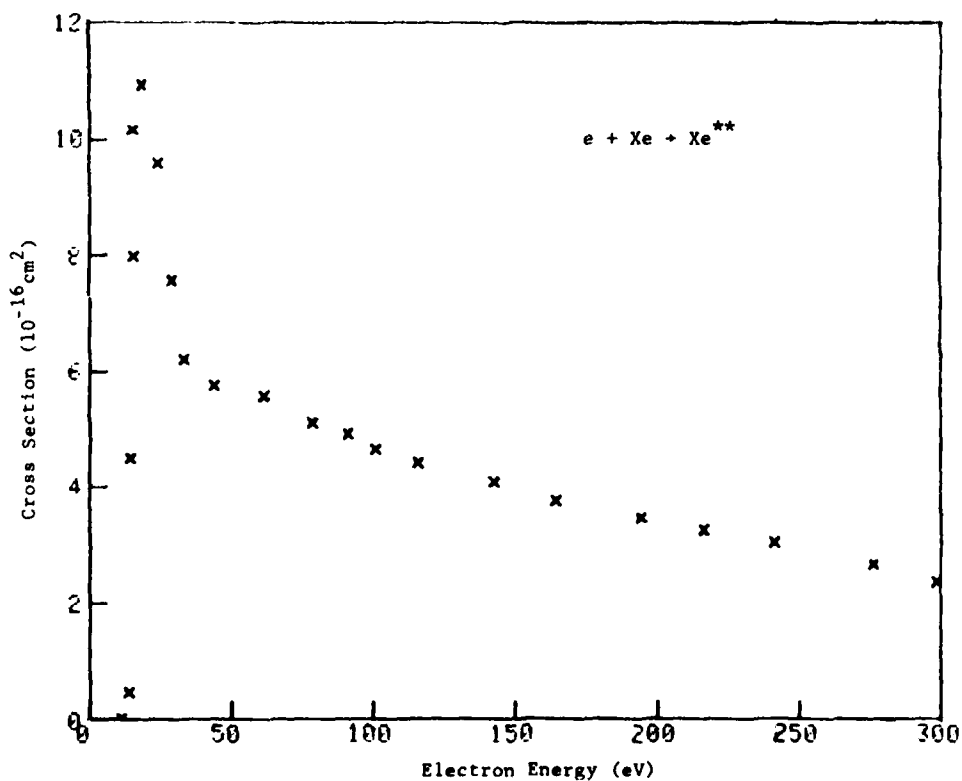
Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, Phys. Rev. A 20, 71 (1979).

Tabular and Graphical Data C-2.11. Cross sections for electron-impact excitation
of Xe atoms to high-Rydberg states.



Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
11	0	91	4.9
14	0.46	100	4.7
14	4.5	120	4.4
15	8.0	140	4.1
15	10	160	3.8
18	11	190	3.5
24	9.6	220	3.3
29	7.6	240	3.0
33	6.2	280	2.7
44	5.7	300	2.4
61	5.6		
78	5.1		

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Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, Phys. Rev. A 20, 71 (1979).

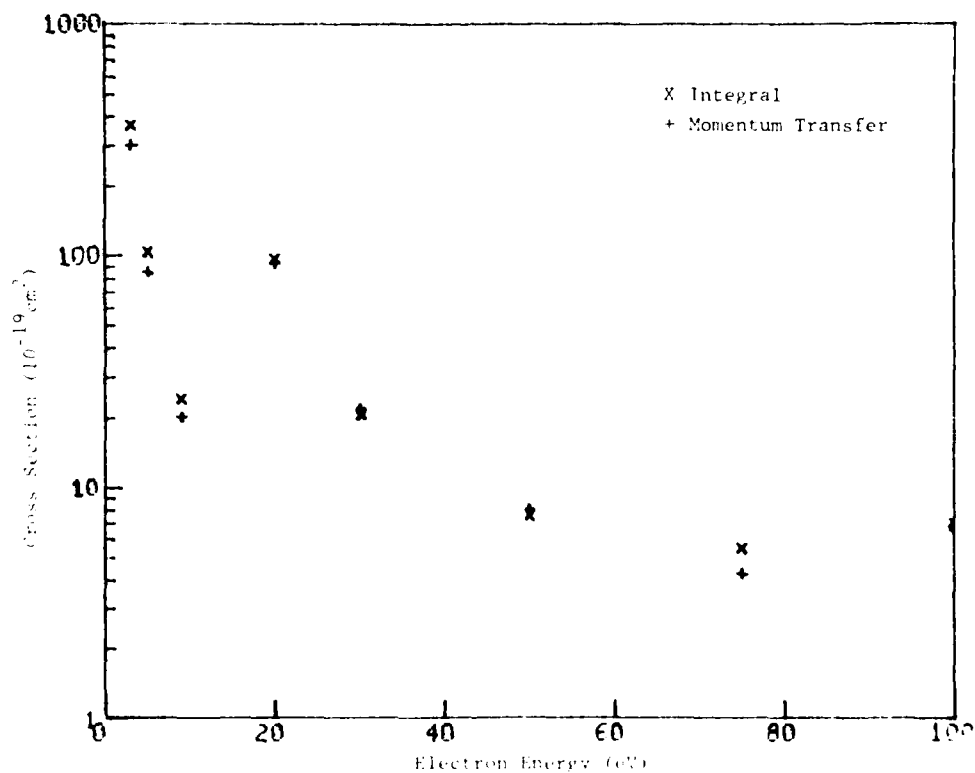
Tabular and Graphical Data C-2.12. Cross sections for $v = 0 \rightarrow 1$
vibrational excitation in CO by electron impact.

Integral cross section

Electron Energy	Cross Section
eV	10^{-19} cm^2
3.0	365
5.0	104
9.0	24.1
20	96.7
30	20.6
50	7.57
75	5.42
100	7.03

Momentum-transfer cross section

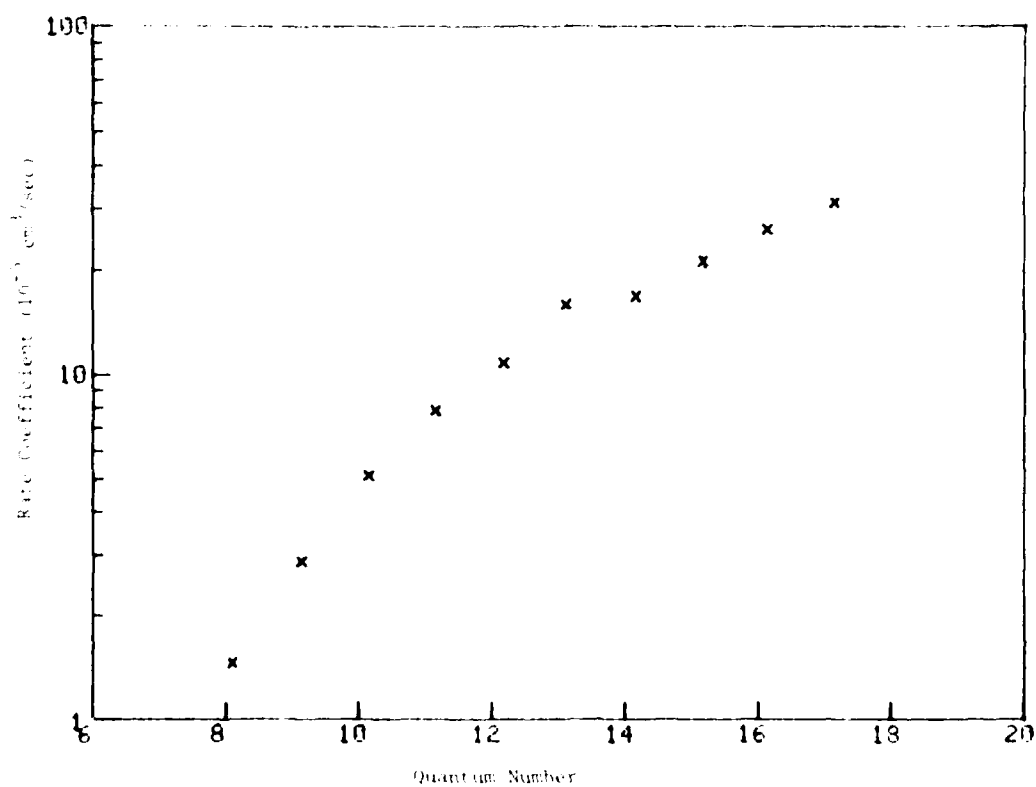
Electron Energy	Cross Section
eV	10^{-19} cm^2
3.0	300
5.0	85.0
9.0	20.3
20	93.4
30	21.3
50	7.97
75	4.71
100	6.48



Reference A Chutjian and H Tanaka, J. Phys. B 13, 1901 (1980)

Tabular and Graphical Data C-2 13. Rate Coefficients
for electron impact depopulation of excited states of
He as a function of principal quantum number.

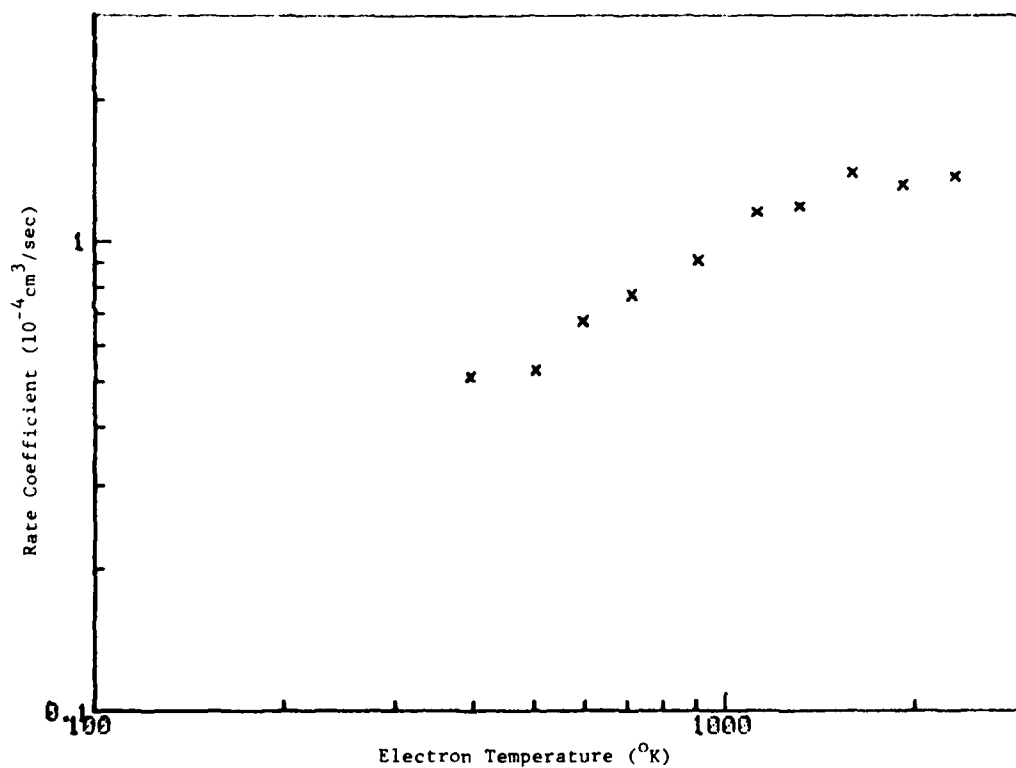
Quantum Number	Rate Coef
	$10^{-5} \text{ cm}^3/\text{sec}$
8	1.46
9	2.87
10	5.10
11	7.86
12	10.8
13	15.9
14	16.9
15	21.1
16	26.2
17	31.2



Reference: E. Devos, J. Boulmer, J.-F. Delpech, J. Physique (Paris) 40, 215 (1979).

Tabular and Graphical Data C-2.14. Rate coefficients
for electron-impact depopulation of He(n=10) as a
function of electron temperature.

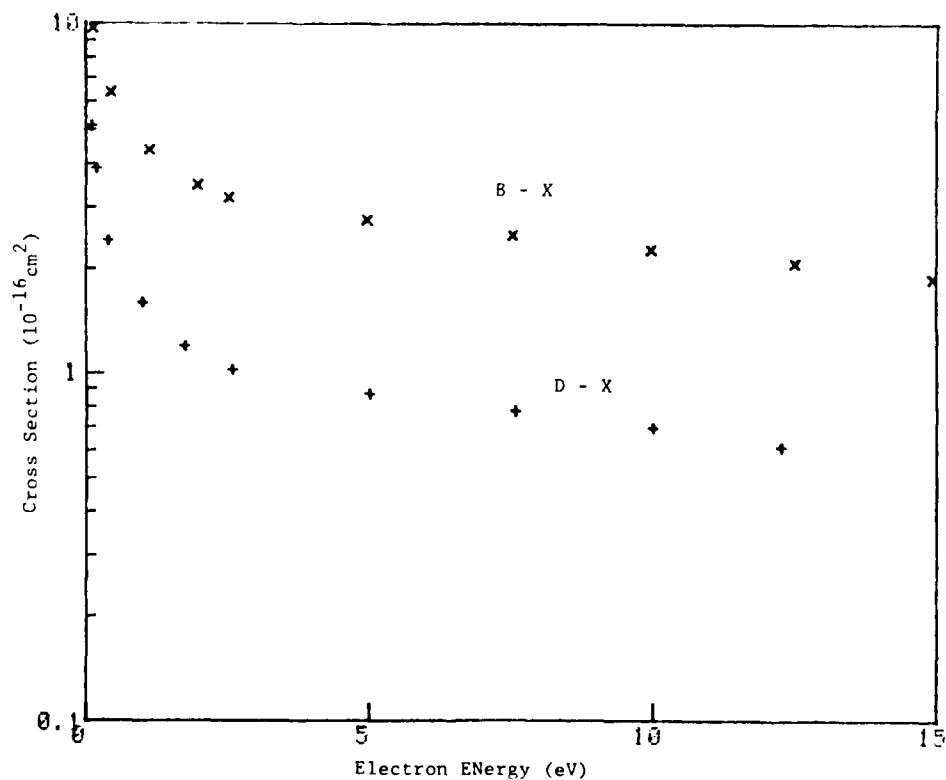
Temperature	Rate Coef
$^{\circ}\text{K}$	$10^{-4} \text{ cm}^3/\text{sec}$
394	0.516
500	0.532
596	0.677
714	0.770
906	0.913
1120	1.15
1310	1.19
1590	1.40
1920	1.32
2320	1.38



Reference: F. Devos, J. Boulmer, and J. -F. Delpech, J. Physique (Paris) 40,
215 (1979).

Tabular and Graphical Data C-2.15. Calculated cross sections for electron-
impact deexcitation of excimer states of KrF

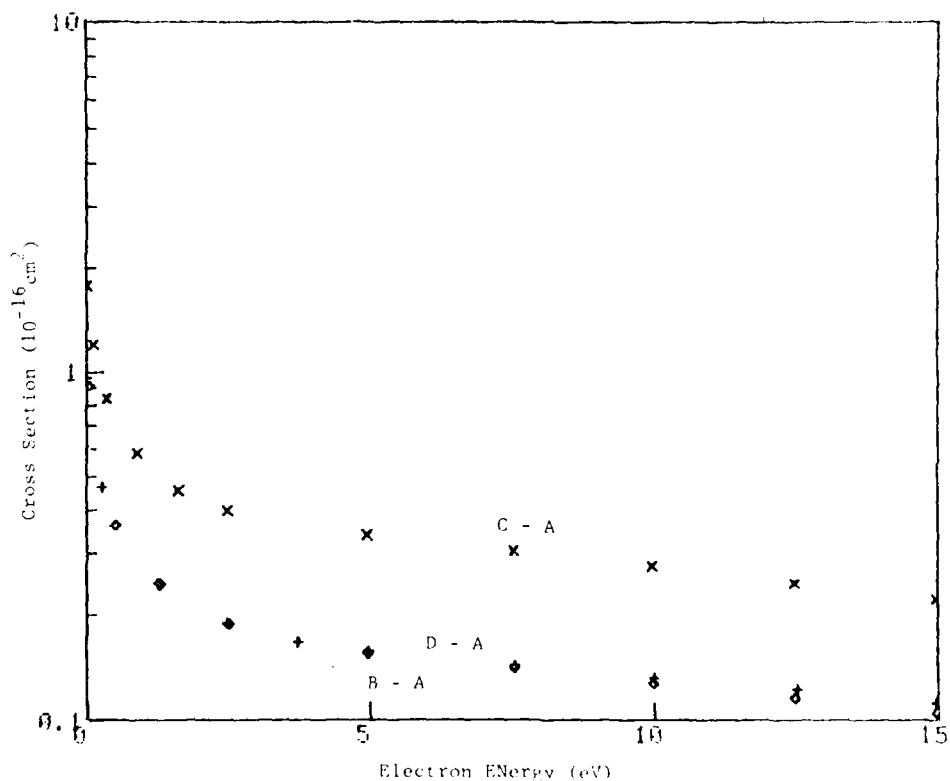
e + KrF(B) + KrF(X)		e + KrF(D) + KrF(X)	
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
0.12	9.70	0.10	5.11
0.42	6.37	0.18	3.87
1.1	4.37	0.39	2.40
1.9	3.47	1.0	1.59
2.5	3.19	1.7	1.19
5.0	2.76	2.6	1.02
7.5	2.49	5.0	0.865
10.0	2.26	7.6	0.778
13	2.07	10.0	0.696
15	1.86	12	0.614



Reference: A. U. Hazi, T.N. Rescigno, and A. E. Orel, Appl. Phys. Lett. 35, 477 (1979).

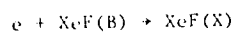
Tabular and Graphical Data C-2.16. Calculated cross section for electron-impact deexcitation of excimer states of KrF.

e + KrF(C) → KrF(A)		e + KrF(D) → KrF(A)		e + KrF(B) → KrF(A)	
Electron Energy	Cross Section	Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2	eV	10^{-16} cm^2
0.0064	1.77	0.00010	0.961	0.041	0.910
0.12	1.50	0.27	0.463	0.50	0.359
0.34	0.853	1.3	0.245	1.3	0.244
0.55	0.565	2.5	0.189	2.5	0.187
1.0	0.453	3.7	0.166	5.0	0.155
1.25	0.395	5.0	0.156	7.5	0.146
4.4	0.337	7.5	0.143	10.0	0.126
7.5	0.305	10	0.132	12	0.115
10.0	0.274	13	0.122	15	0.105
15	0.245	15	0.112		
15	0.221				

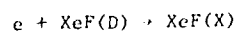


Reference A. U. Hazi, T. N. Rescigno, and A. E. Orei, Appl. Phys. Lett. 35, 477 (1979).

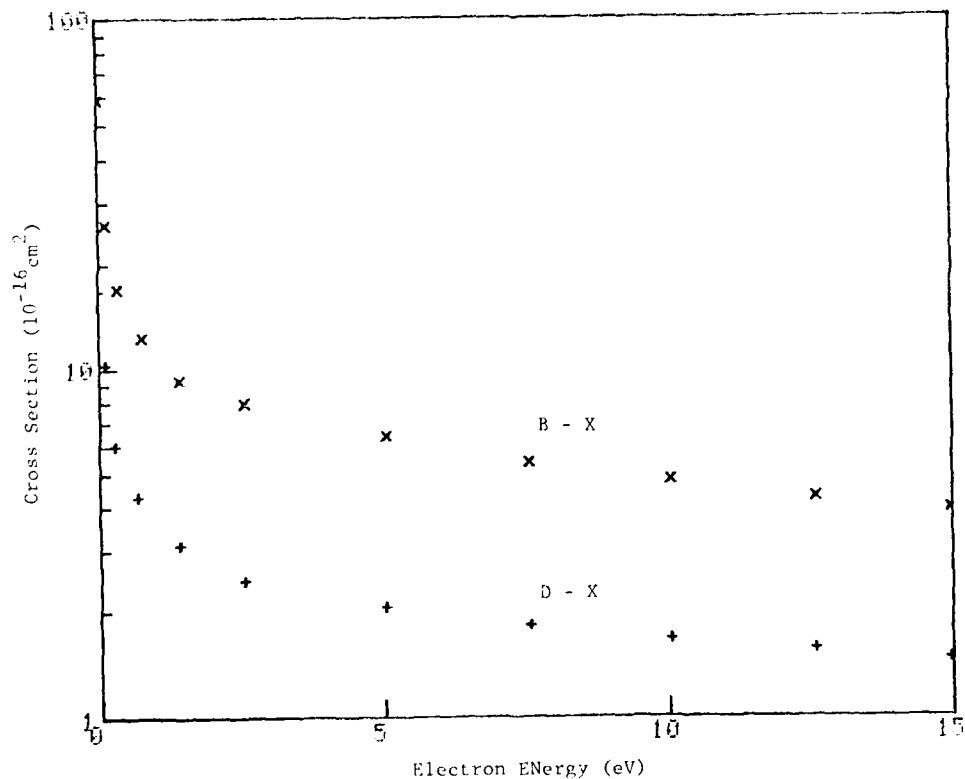
Tabular and Graphical Data C-2.17. Calculated cross sections for electron-impact deexcitation of excimer states of XeF



Electron Energy eV	Cross Section 10^{-16} cm^2
0.00068	59.0
0.11	25.9
0.33	16.9
0.75	12.2
1.4	9.30
2.5	7.95
5.0	6.39
7.6	5.37
10	4.78
13	4.25
15	3.92



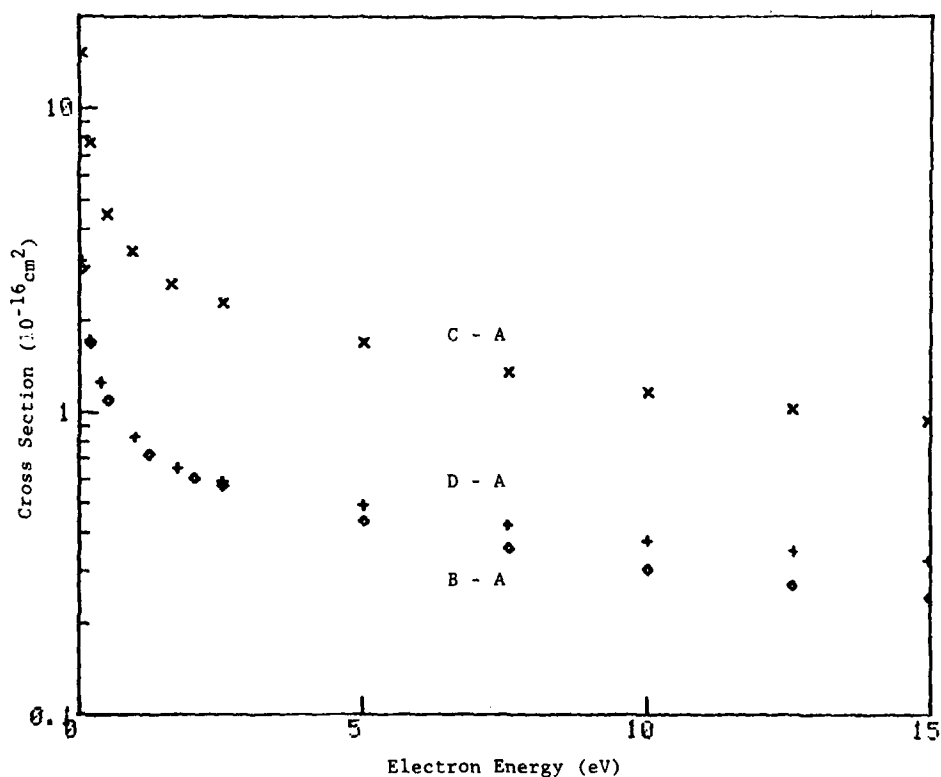
Electron Energy eV	Cross Section 10^{-16} cm^2
0.00010	16.7
0.095	10.3
0.28	6.01
0.66	4.26
1.4	3.10
2.5	2.46
5.0	2.06
7.6	1.83
10	1.67
13	1.55
15	1.46



Reference: A. U. Hazi, T. N. Rescigno, and A. E. Orel, Appl. Phys. Lett. 35, 477 (1979).

Tabular and Graphical Data C-2.18. Calculated cross sections for electron-impact deexcitation of excimer states of XeF

e + XeF(C) → XeF(A)		e + XeF(D) → XeF(A)		e + XeF(B) → XeF(A)	
Electron Energy	Cross Section	Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10 ⁻¹⁶ cm ²	eV	10 ⁻¹⁶ cm ²	eV	10 ⁻¹⁶ cm ²
0.033	15.2	0.027	3.14	0.043	2.95
0.17	7.72	0.17	1.72	0.17	1.69
0.49	4.50	0.36	1.25	0.49	1.10
0.92	3.37	0.98	0.823	1.2	0.722
1.6	2.63	1.7	0.652	2.0	0.604
2.5	2.29	2.5	0.590	2.5	0.568
5.0	1.70	5.0	0.492	5.0	0.436
7.6	1.35	7.6	0.421	7.6	0.355
10	1.16	10	0.377	10	0.301
13	1.03	13	0.349	13	0.267
15	0.931	15	0.321	15	0.242



Reference: A. U. Hazi, T. N. Rescigno, and A. E. Orel, Appl. Phys. Lett. 35, 477 (1979).

Tabular Data C-2.19. Calculated rate coefficients for electron impact deexcitation of excimer states of KrF in units of $10^{-9} \text{ cm}^3/\text{sec}$ (taken from Fig. 4 of the reference).

Transition	Electron Energy (eV)	1	2	4
B - X		27	30	31
D - X		9.0	9.7	10
C - A		3.4	3.7	3.8
D - A		1.6	1.8	1.9
B - A		1.6	1.8	1.9

Tabular Data C-2.20. Calculated rate coefficients for electron-impact deexcitation of excimer states of XeF in units of $10^{-9} \text{ cm}^3/\text{sec}$ (taken from Fig. 5 of the reference).

Transition	Electron Energy (eV)	1	2	4
B - X		66	71	74
D - X		21	23	24
C - A		18	19	20
D - A		4.8	5.1	5.5
B - A		4.5	4.9	4.9

Reference: A. U. Hazi, T.N. Rescigno, and A. E. Orel, Appl. Phys. Lett. 35 477 (1979).

C-3. DISSOCIATION BY ELECTRON IMPACT

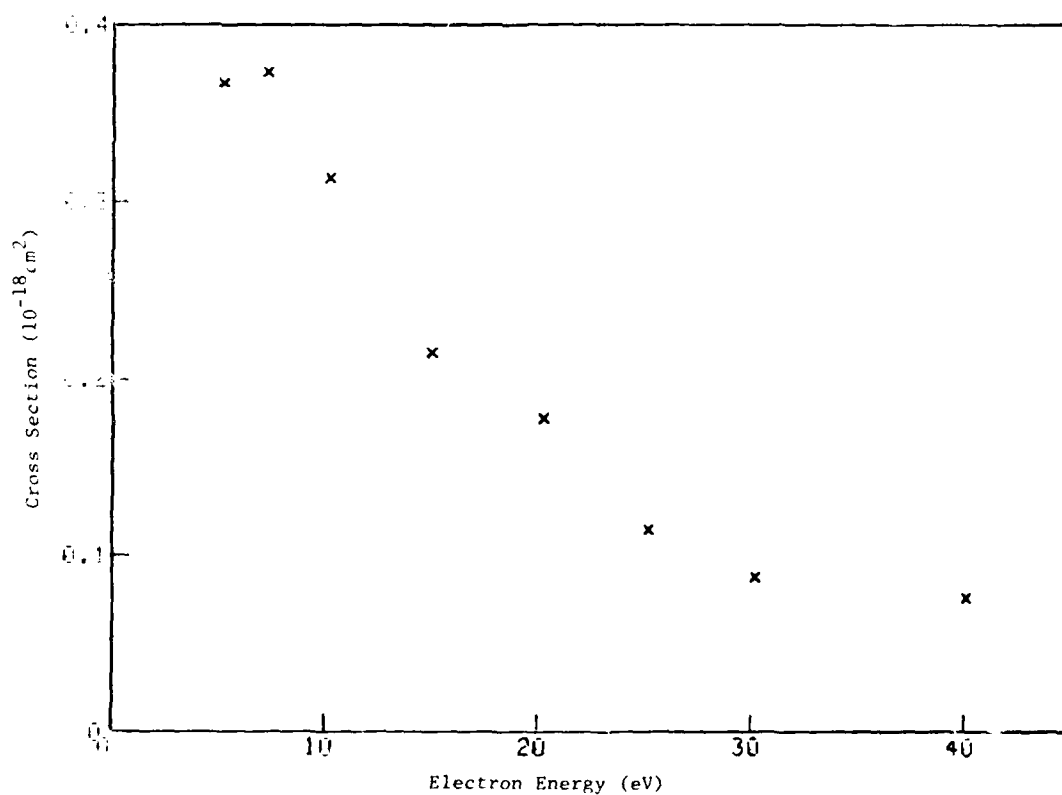
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Tabular and Graphical Data C-3.1. Calculate
cross sections for electron impact dissociation of F_2
(Distorted wave model with static exchange)

electron Energy	Cross Section
ev	10^{-16} cm^2
5.2	0.368
7.2	0.374
10	0.314
15	0.219
20	0.173
25	0.119
30	0.0880
40	0.0752



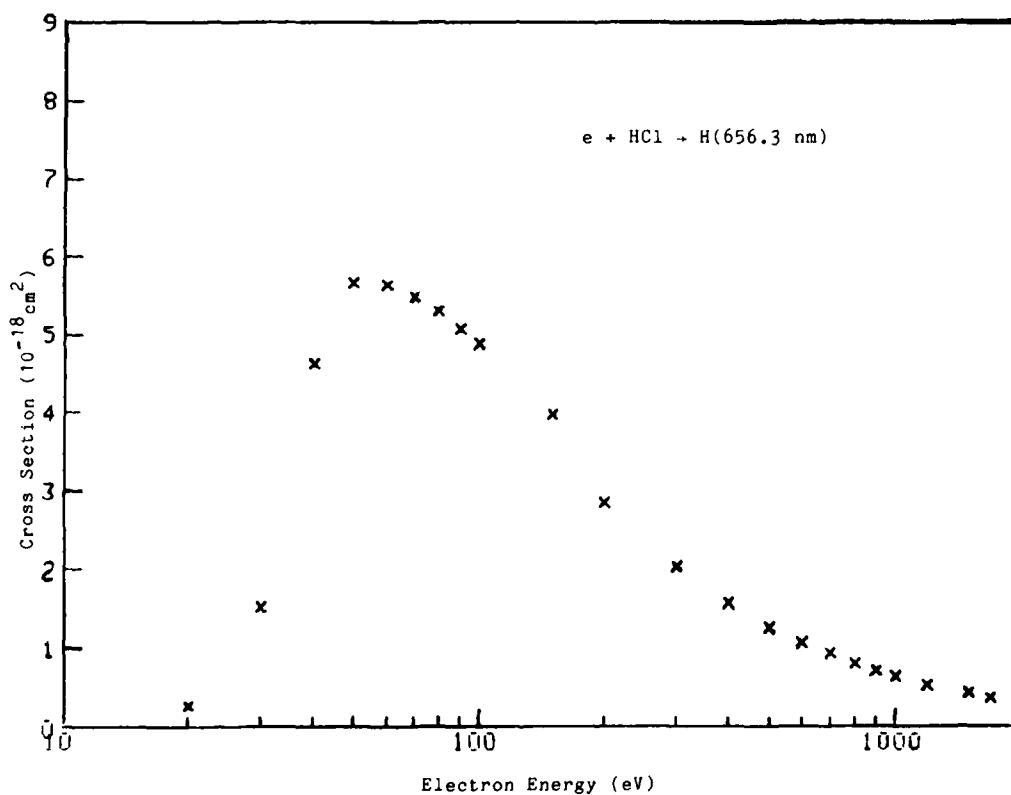
Reference: A. W. Fliflet, V. McKoy, and T. N. Rescigno, Phys. Rev. A 21, 788 (1980).

Tabular and Graphical Data C-3.2. Cross sections for electron-impact
dissociation of HCl to form excited fragments.



Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-18} cm^2	eV	10^{-18} cm^2
20	0.265	400	1.57
30	1.51	500	1.27
40	4.62	600	1.07
50	5.66	700	0.932
60	5.64	800	0.803
70	5.49	900	0.714
80	5.30	1000	0.638
90	5.07	1200	0.531
100	4.88	1500	0.432
150	3.98	1700	0.381
200	2.85	2000	0.324
300	2.03		

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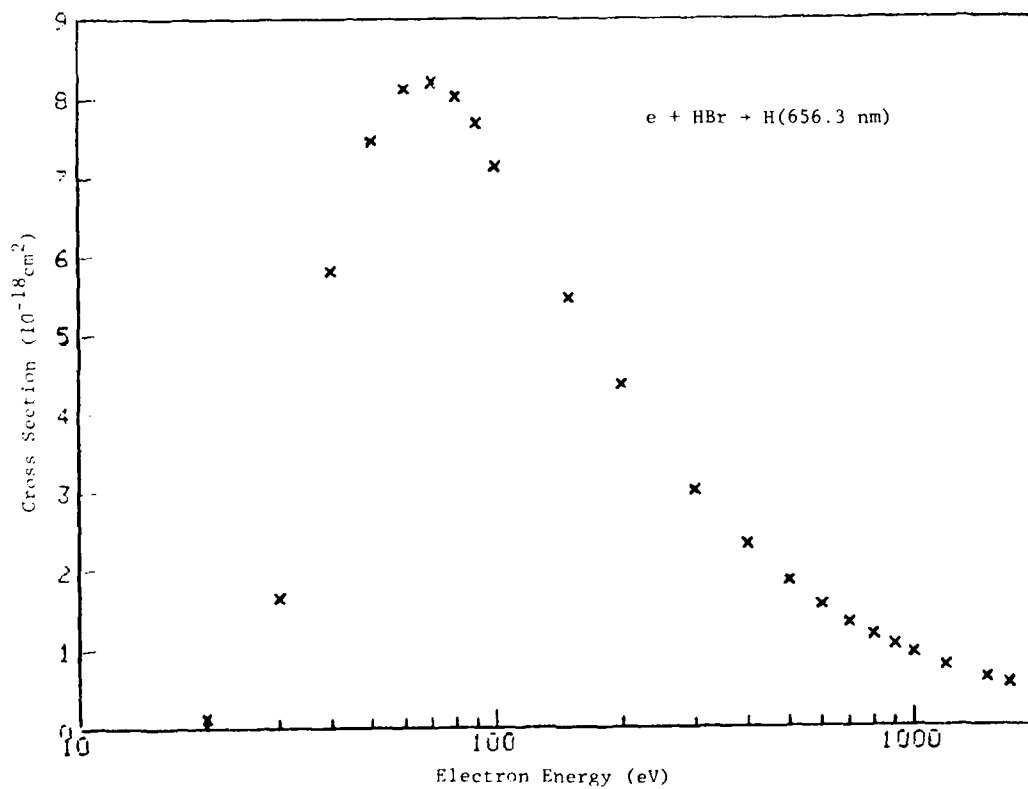


Reference: G. R. Mohlmann and F. J. de Heer, Chem. Phys. 40, 157 (1979)

Tabular and Graphical Data C-3.3. Cross sections for electron-impact
dissociation of HBr to form excited fragments
 $e + \text{HBr} \rightarrow \text{H} (656.3 \text{ nm})$

Electron Energy eV	Cross Section 10^{-18} cm^2	Electron Energy eV	Cross Section 10^{-18} cm^2
20	0.110	400	2.32
30	1.65	500	1.85
40	5.82	600	1.55
50	7.47	700	1.33
60	6.13	800	1.17
70	8.21	900	1.03
80	8.02	1000	0.941
90	7.69	1200	0.780
100	7.14	1500	0.624
150	5.47	1700	0.545
200	4.37	2000	0.464
300	3.01		

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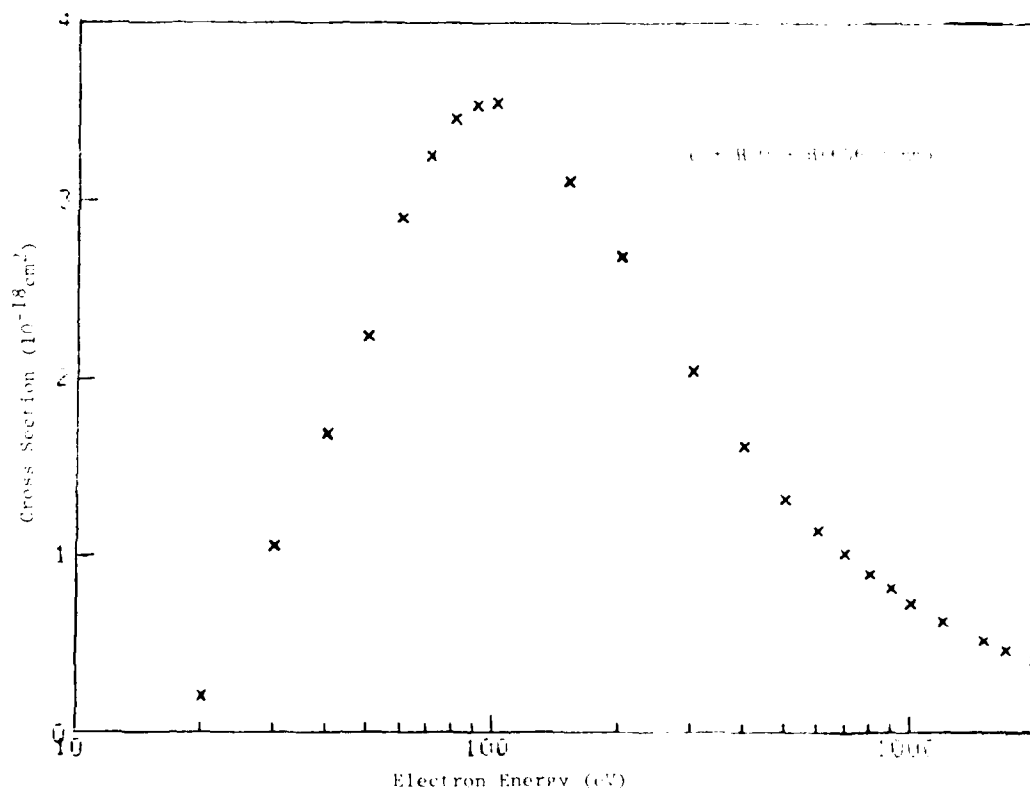


Reference: G. R. Mohlmann and F. J. de Heer, Chem. Phys. 40, 157 (1979).

Tabular and Graphical Data (Fig. 1) Cross Sections for Electron Impact
Dissociation of H_2O to Form Excited Fragment
 $e + H_2O \rightarrow H + O(^1D, ^3P, \dots)$

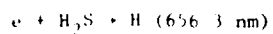
Electron Energy (eV)	Cross Section (10^{-18} cm^2)	Electron Energy (eV)	Cross Section (10^{-18} cm^2)
40	0.000	40	0.000
50	0.000	50	0.000
60	0.000	60	0.000
70	0.000	70	0.000
80	0.000	80	0.000
90	0.000	90	0.000
100	0.000	100	0.000
110	0.000	110	0.000
120	0.000	120	0.000
130	0.000	130	0.000
140	0.000	140	0.000
150	0.000	150	0.000
160	0.000	160	0.000
170	0.000	170	0.000
180	0.000	180	0.000
190	0.000	190	0.000
200	0.000	200	0.000

Fig. 1. Excited states.



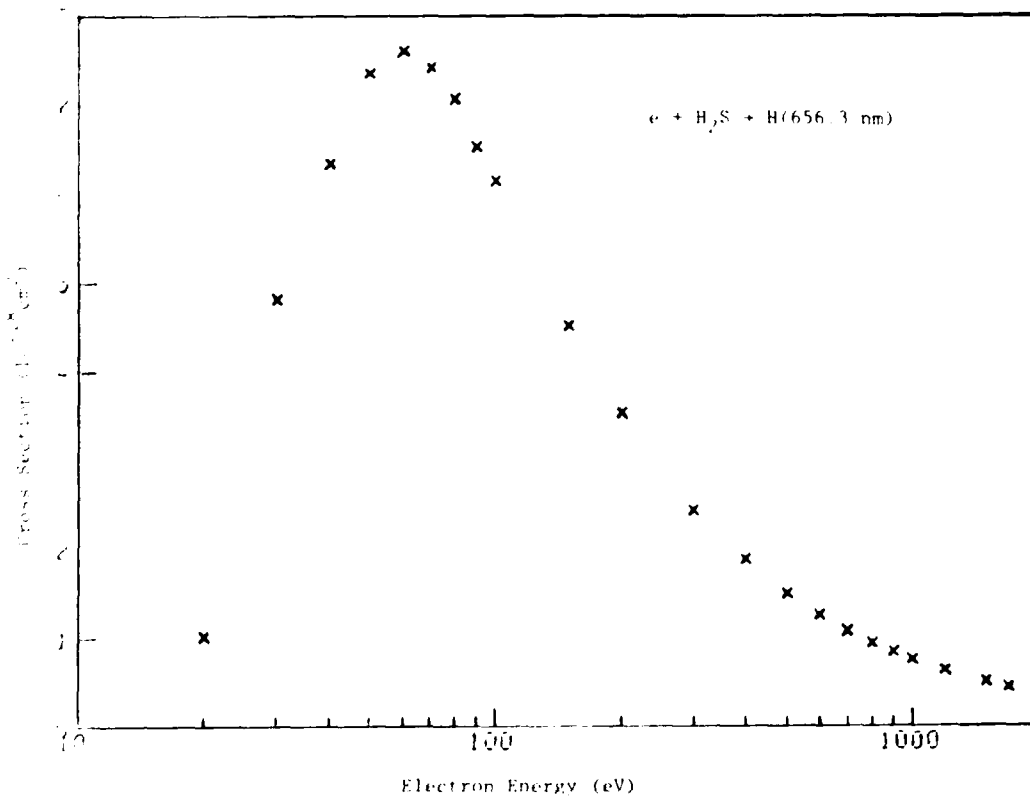
Reference: G. R. Mohlmann and F. J. de Heer, Chem. Phys. 40, 157 (1979).

Tabular and Graphical Data C-3.5 Cross sections for electron-impact
dissociation of H₂S to form excited fragments.



Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10 ⁻¹⁸ cm ²	eV	10 ⁻¹⁸ cm ²
400	1.50	400	1.50
500	4.531	500	1.50
600	6.576	600	1.50
700	7.527	700	1.50
800	7.527	800	2.948
900	7.40	900	3.543
1000	7.25	1000	3.754
1200	6.204	1200	3.632
1500	5.16	1500	3.506
1800	4.13	1700	3.446
2000	3.225	2000	3.372
3000	2.43		

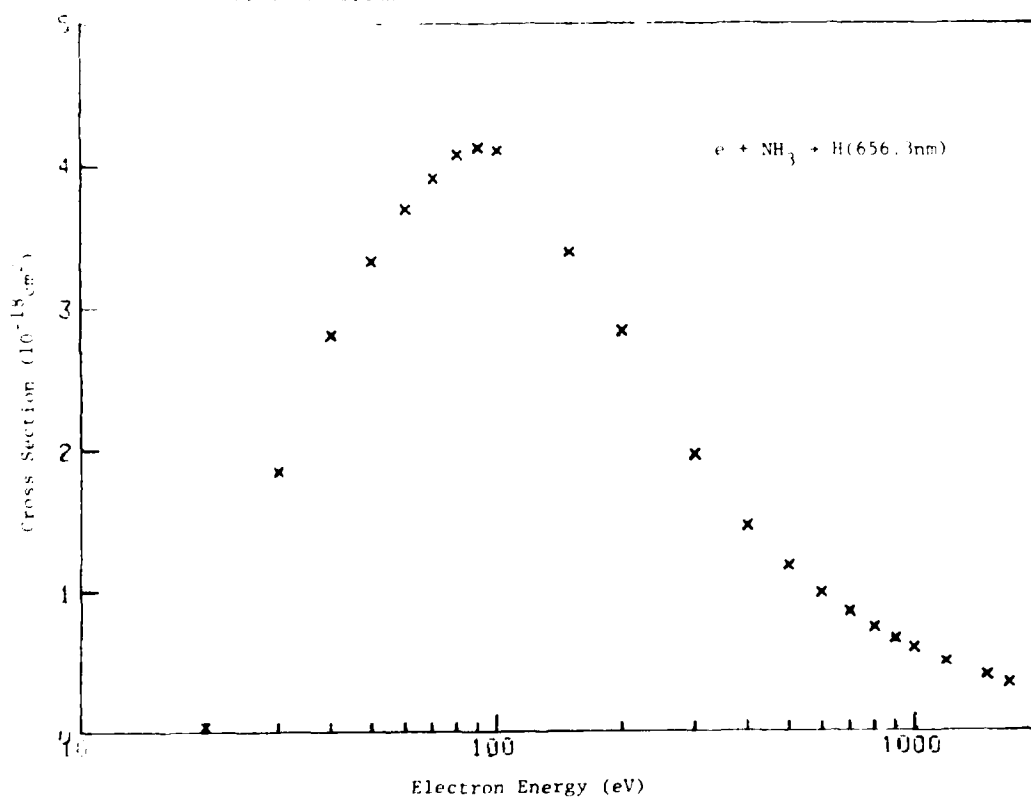
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Reference G. R. Mohlmann and F. J. de Heer, Chem. Phys. 40, 157 (1979).

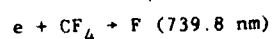
Tabular and Graphical Data C-3.6. Cross sections for electron-impact
dissociation of NH_3 to form excited fragments
 $e + \text{NH}_3 \rightarrow \text{H} (656.3 \text{ nm})$

Electron Energy (eV)	Electron Energy (eV)	Electron Energy (eV)	Cross Section (10^{-18} cm^2)
E_1	E_2	E_3	$\sigma_{\text{exc}}(E)$
10	10.0	10	1.45
20	20.0	20	1.15
30	30.0	30	1.4
40	40.0	40	1.5
50	50.0	50	1.25
60	60.0	60	1.515
70	70.0	70	1.555
80	80.0	80	1.45
90	90.0	90	1.45
100	100.0	100	1.45
150	150.0	150	1.45
200	200.0	200	1.45
300	300.0	300	1.45
400	400.0	400	1.45
500	500.0	500	1.45
600	600.0	600	1.45
700	700.0	700	1.45
800	800.0	800	1.45
900	900.0	900	1.45
1000	1000.0	1000	1.45



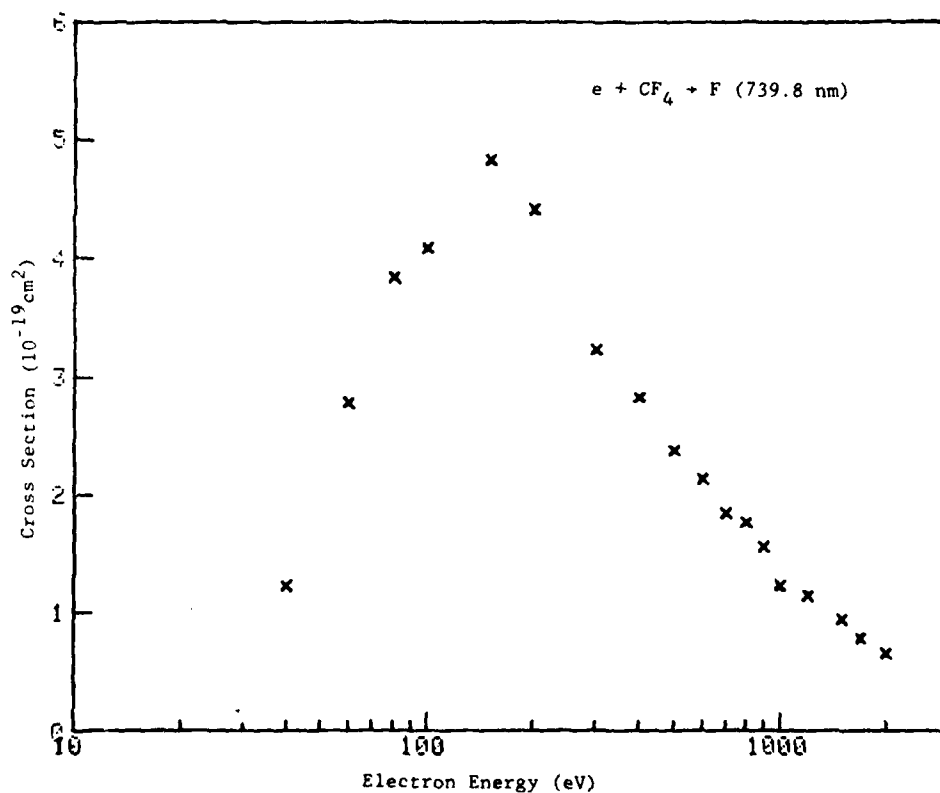
Reference: G. R. Mohlmann and F. J. de Heer, Chem. Phys. 40, 157 (1979).

Tabular and Graphical Data C-3.7. Cross sections for electron-impact
dissociation of CF_4 to form excited fragments.



Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-19}cm^2	eV	10^{-19}cm^2
40	1.23	700	1.85
60	2.79	800	1.76
80	3.85	900	1.56
100	4.10	1000	1.23
150	4.84	1200	1.15
200	4.43	1500	0.943
300	3.24	1700	0.779
400	2.83	2000	0.656
500	2.38		
600	2.13		

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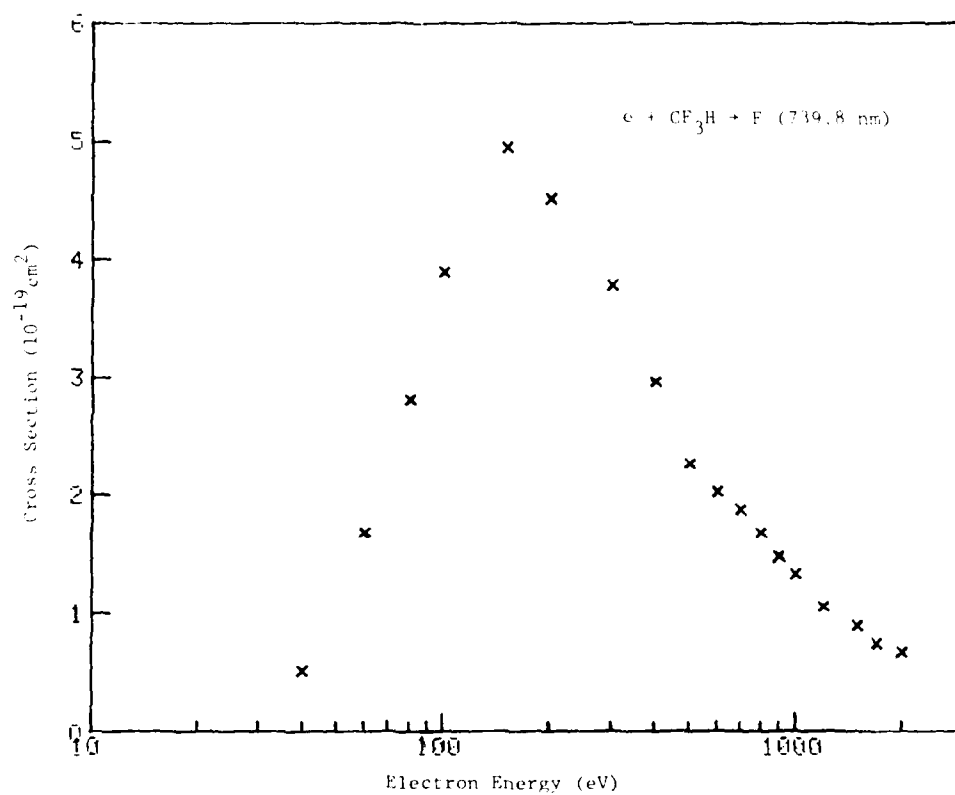


Reference: H. A. van Sprang, H. H. Brongersma, and F. J. de Heer, Chem. Phys. 35, 51 (1978).

Tabular and Graphical Data C-3.8. Cross sections for electron-impact
dissociation of CF_3H to form excited fragments
 $\text{e} + \text{CF}_3\text{H} \rightarrow \text{F} (739.8 \text{ nm})$

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-19} cm^2	eV	10^{-19} cm^2
40	0.51	700	1.9
50	1.7	800	1.7
60	1.5	900	1.5
100	3.9	1000	1.5
150	5.5	1200	1.1
200	4.5	1500	0.90
300	3.3	1700	0.74
400	3.0	2000	0.66
500	2.3		
600	2.0		

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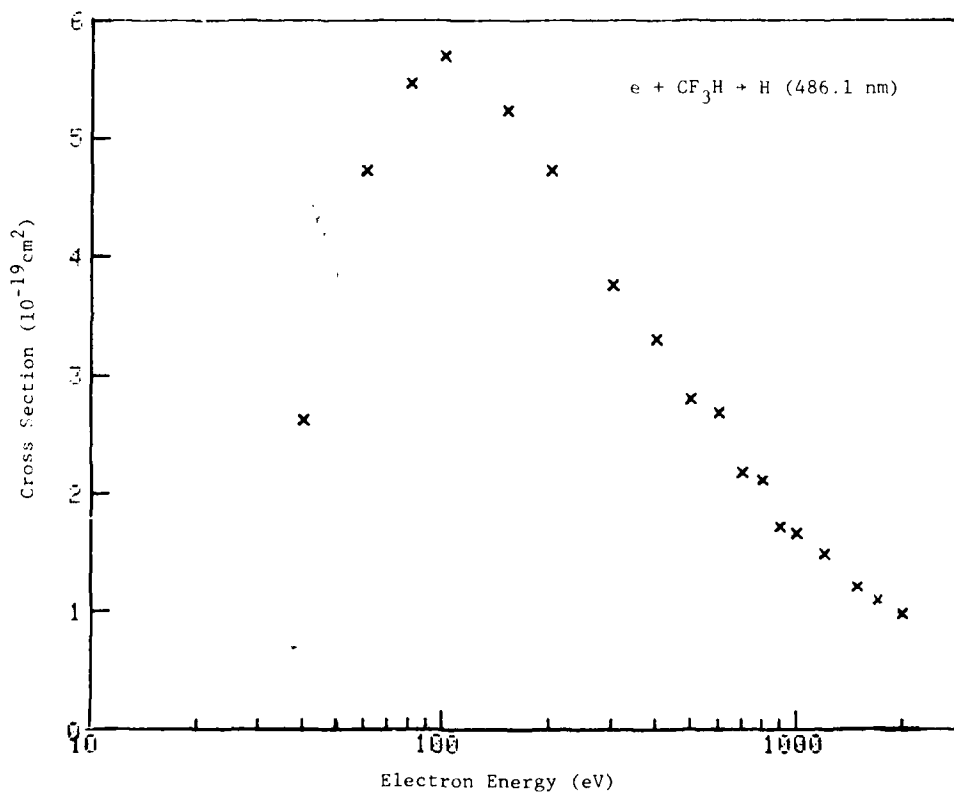


Reference: H. A. van Sprang, H. H. Brongersma, and F. J. de Heer, Chem. Phys. 35, 51 (1978).

Tabular and Graphical Data C-3.9. Cross sections for electron-impact
dissociation of CF_3H to form excited fragments.
 $\text{e} + \text{CF}_3\text{H} \rightarrow \text{H} (486.1 \text{ nm})$

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-19} cm^2	eV	10^{-19} cm^2
40	2.6	700	2.4
60	4.7	800	2.1
80	5.5	900	1.7
100	5.7	1000	1.7
150	5.2	1200	1.5
200	4.7	1500	1.2
300	3.8	1700	1.1
400	3.3	2000	0.97
500	2.8		
600	2.7		

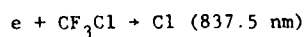
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Reference: H. A. van Sprang, H. H. Brongersma, and F. J. de Heer, Chem. Phys. 35
51 (1978).

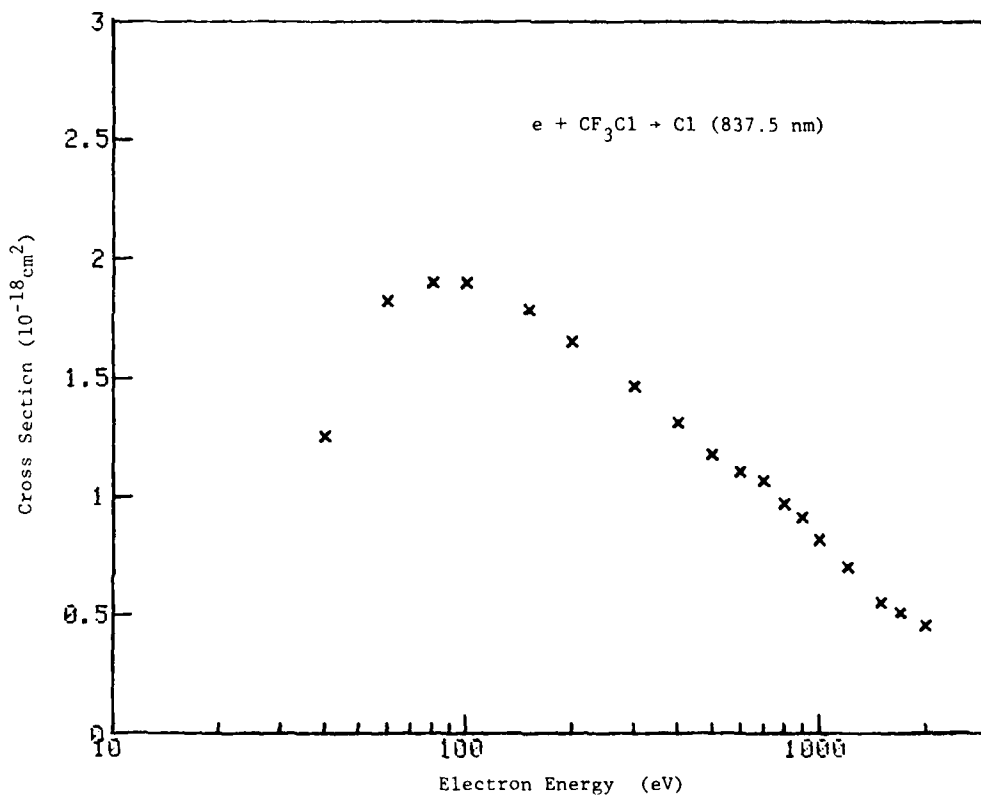
Tabular and Graphical Data C-3.10. Cross Sections for electron-impact

dissociation of CF_3Cl to form excited fragments.



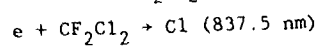
Electron Energy eV	Cross Section 10^{-18} cm^2	Electron Energy eV	Cross Section 10^{-18} cm^2
40	1.3	700	1.1
60	1.8	800	0.97
80	1.9	900	0.91
100	1.9	1000	0.82
150	1.8	1200	0.70
200	1.7	1500	0.55
300	1.5	1700	0.51
400	1.3	2000	0.46
500	1.2		
600	1.1		

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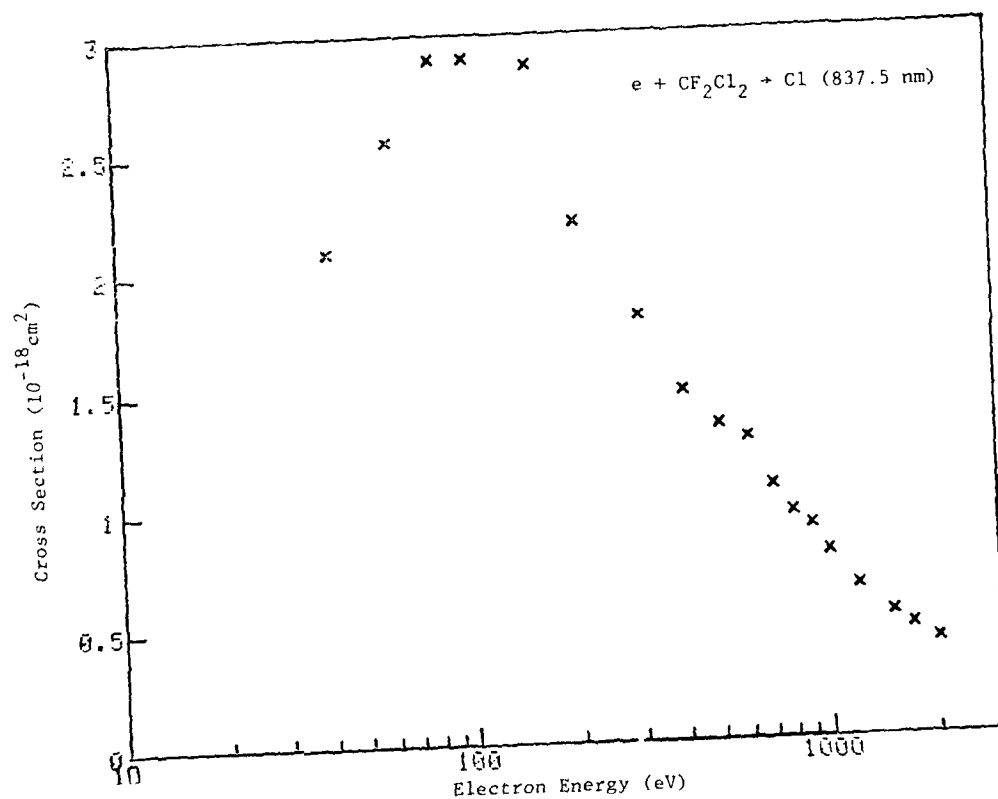
Reference: H. A. van Sprang, H. H. Brongersma, and F. J. de Heer, Chem. Phys. 35, 51 (1978).

Tabular and Graphical Data C-3.11. Cross sections for electron-impact
dissociation of CF_2Cl_2 to form excited fragments.



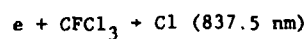
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-18} cm^2	eV	10^{-18} cm^2
40	2.1	700	1.1
60	2.6	800	0.96
80	2.9	900	0.90
100	2.9	1000	0.78
150	2.9	1200	0.64
200	2.2	1500	0.52
300	1.6	1700	0.46
400	1.5	2000	0.41
500	1.3		
600	1.3		

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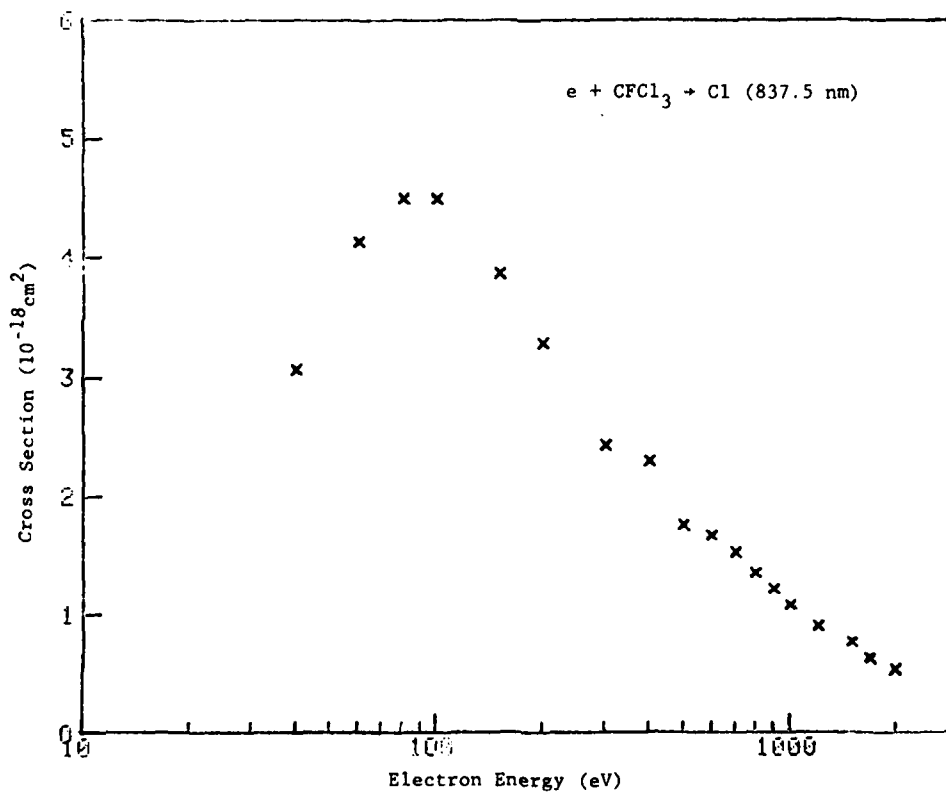
Reference: H. A. van Sprang, H. H. Brongersma, and F. J. de Heer, Chem. Phys. 35, 51 (1978).

Tabular and Graphical Data C-3.12. Cross sections for electron-impact
dissociation of CFCl_3 to form excited fragments.



Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-18}cm^2	eV	10^{-18}cm^2
40	3.1	700	1.5
60	4.1	800	1.4
80	4.5	900	1.2
100	4.5	1000	1.1
150	3.9	1200	0.90
200	3.3	1500	0.77
300	2.4	1700	0.63
400	2.3	2000	0.54
500	1.8		
600	1.7		

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Reference: H. A. van Sprang, H. H. Brongersma, and F. J. de Heer, Chem. Phys. 35, 51 (1978).

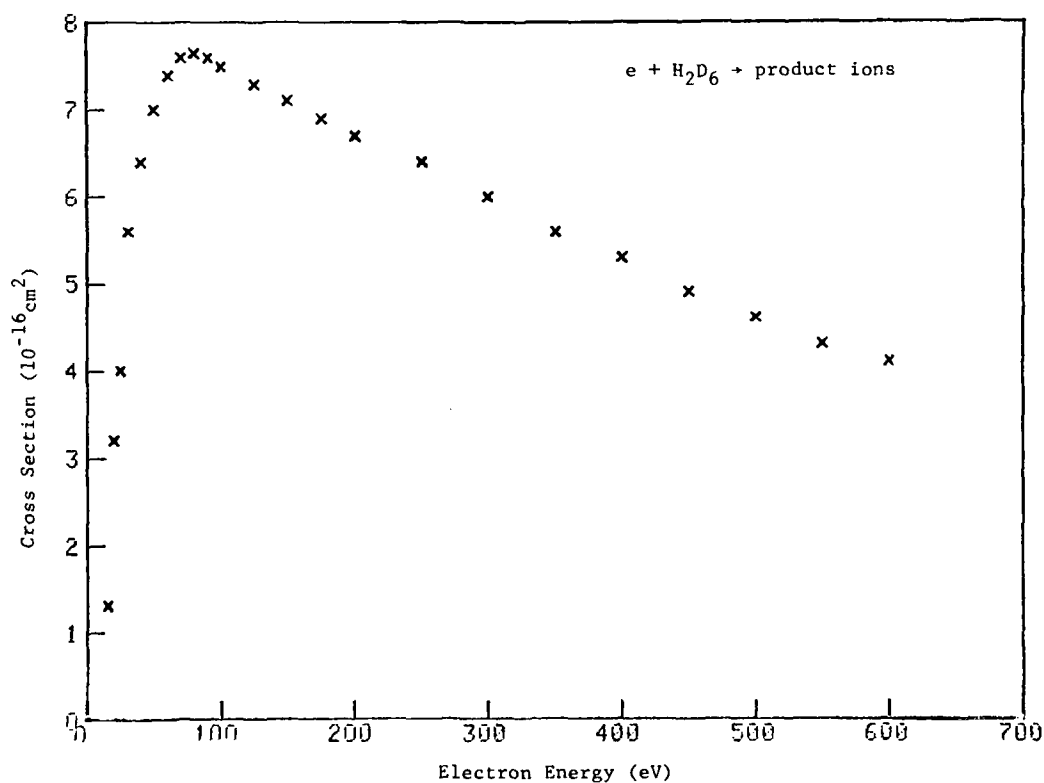
Tabular and Graphical Data C-3.13. Total dissociation cross sections for electrons

incident on C_2D_6 .

$e + C_2D_6 \rightarrow$ product ions

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	$10^{-16} cm^2$	eV	$10^{-16} cm^2$
15	1.30	150	7.10
20	3.20	180	6.90
25	4.00	200	6.70
30	5.60	250	6.40
40	6.40	300	6.00
50	7.00	350	5.60
60	7.40	400	5.30
70	7.60	450	4.90
80	7.65	500	4.60
90	7.60	550	4.30
100	7.50	600	4.10
130	7.30		

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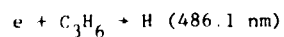


Tabular Data C-3.14. Ratio of total dissociation cross sections for C_2H_6 to that of C_2D_6 .

Electron Energy	Ratio	Electron Energy	Ratio
eV		eV	
50	1.14	300	1.06
100	1.07	400	1.05
150	1.07	500	1.14
200	1.03	600	1.11

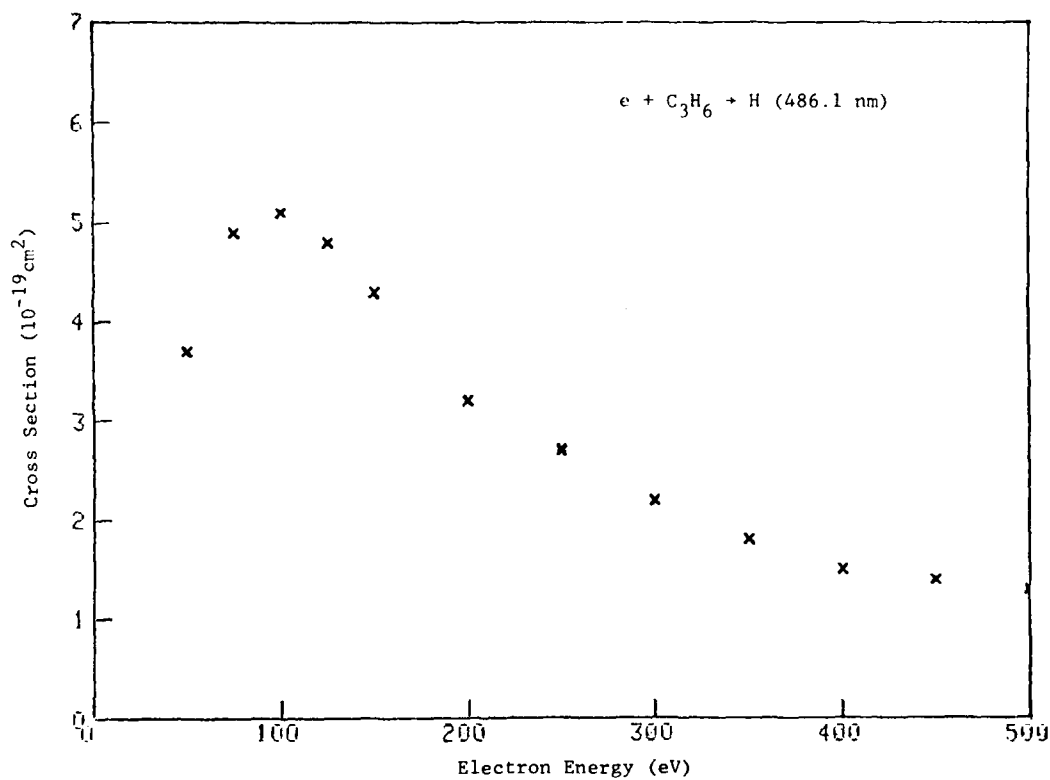
Reference: H. F. Winters, Chem. Phys. 36, 364 (1979).

Tabular and Graphical Data C-3.15. Cross sections for electron-impact
dissociation of propylene to form excited fragments.



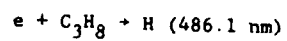
electron energy	Cross Section	Electron Energy	Cross Section
eV	10^{-19} cm^2	eV	10^{-19} cm^2
50.0	3.7	300	2.2
75.	4.9	350	1.5
100	5.1	400	1.5
125	4.3	450	1.4
150	4.3	500	1.3
200	3.2		
250	2.7		

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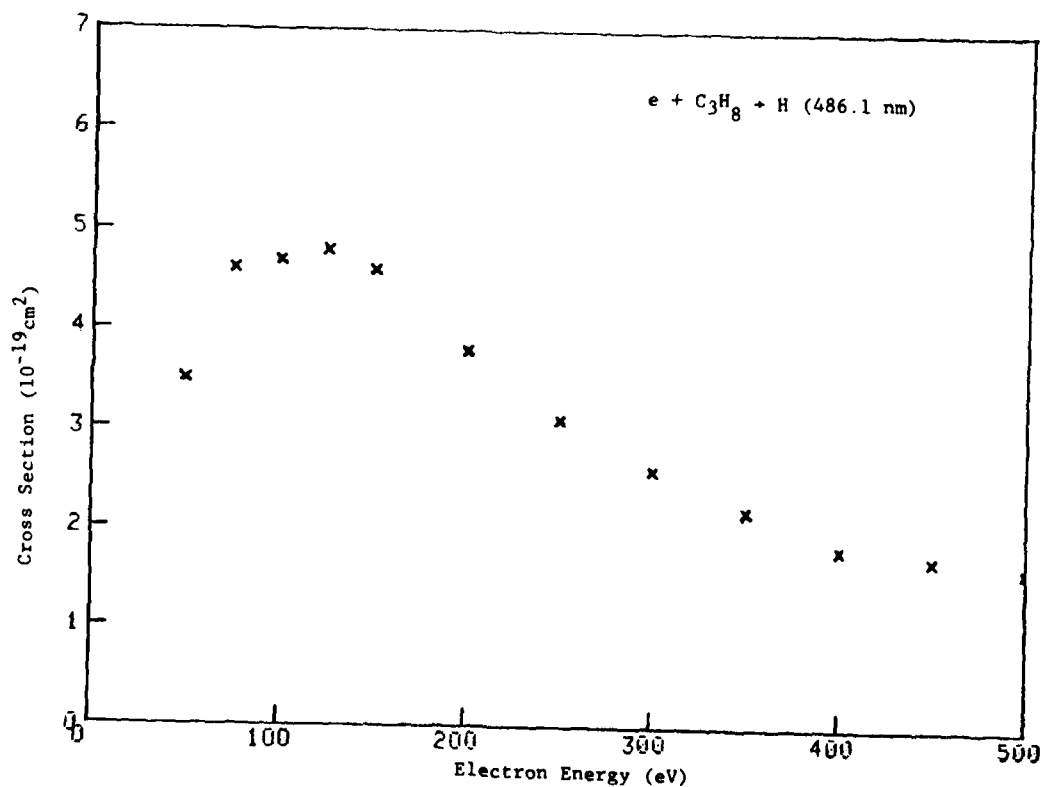
Reference: J. M. Kurepa and M. D Tasic, Chem. Phys. 38, 361 (1979).

Tabular and Graphical Data C-3.16. Cross sections for electron-impact
dissociation of propane to form excited fragments.



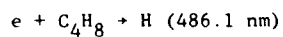
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-19}cm^2	eV	10^{-19}cm^2
50.0	3.5	300	2.6
75.0	4.6	350	2.2
100	4.7	400	1.8
125	4.8	450	1.7
150	4.6	500	1.6
200	3.8		
250	3.1		

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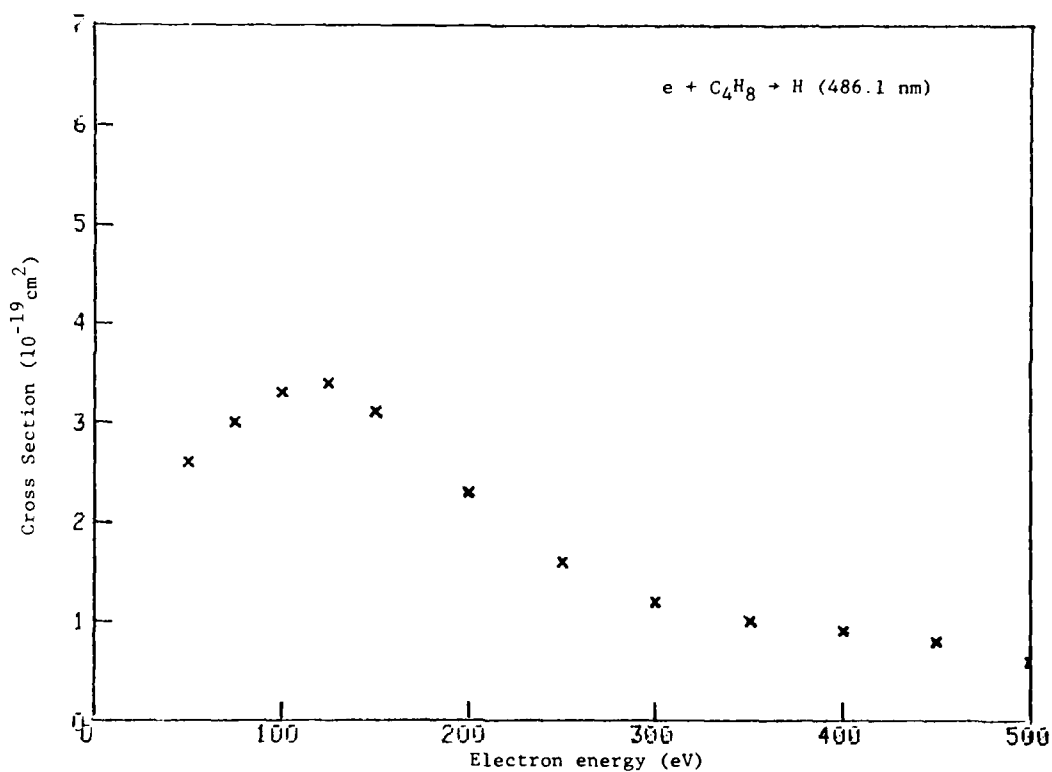
Reference: J. M. Kurepa and M. D. Tasic, Chem. Phys. 38, 361 (1979).

Tabular and Graphical Data C-3.17. Cross sections for electron-impact
dissociation of 1-butene to form excited fragments.



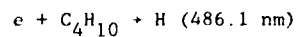
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-19} cm^2	eV	10^{-19} cm^2
50.0	2.6	300	1.2
75.0	3.0	350	1.00
100	3.3	400	0.90
125	3.4	450	0.80
150	3.1	500	0.60
200	2.3		
250	1.6		

Cont. Next Column



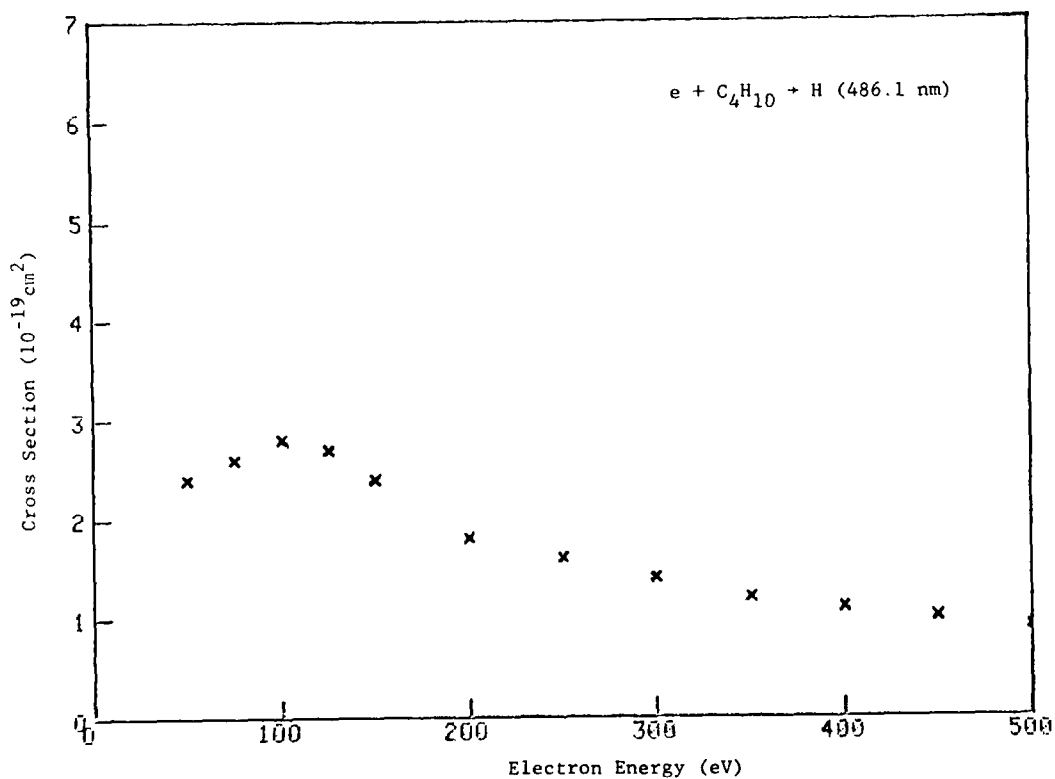
Reference: J. M. Kurepa and M. D. Tasic, Chem. Phys. 38, 361 (1979).

Tabular and Graphical Data C-3.18. Cross sections for electron-impact
dissociation of n-butane to form excited fragments.



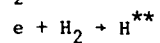
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-19} cm^2	eV	10^{-19} cm^2
50.0	2.4	300	1.4
75.0	2.6	350	1.2
100	2.6	400	1.1
125	2.7	450	1.00
150	2.4	500	0.90
200	1.8		
250	1.6		

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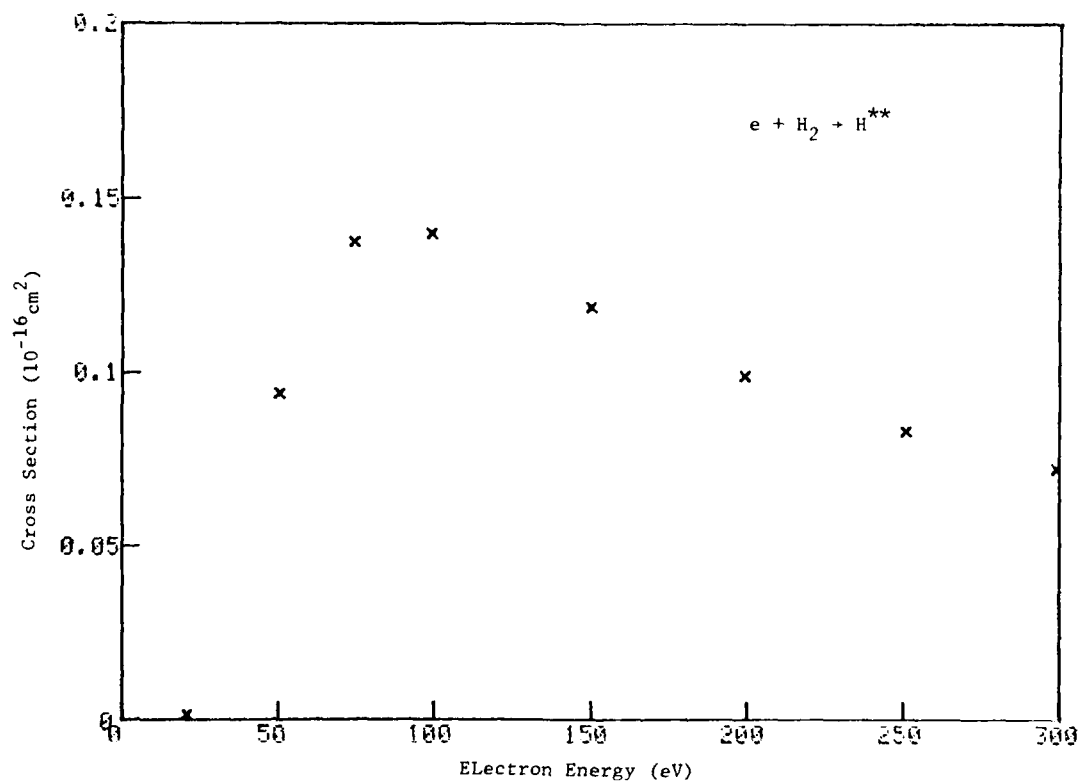


Reference: J. M. Kurepa and M. D. Tasic, Chem. Phys. 38, 361 (1979).

Tabular and Graphical Data C-3.19. Cross sections for electron-
impact dissociation of H_2 to form high-Rydberg fragments.

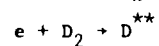


Electron Energy	Cross Section
eV	10^{-16} cm^2
21	0.0014
51	0.094
74	0.14
99	0.14
150	0.12
200	0.099
250	0.083
300	0.072

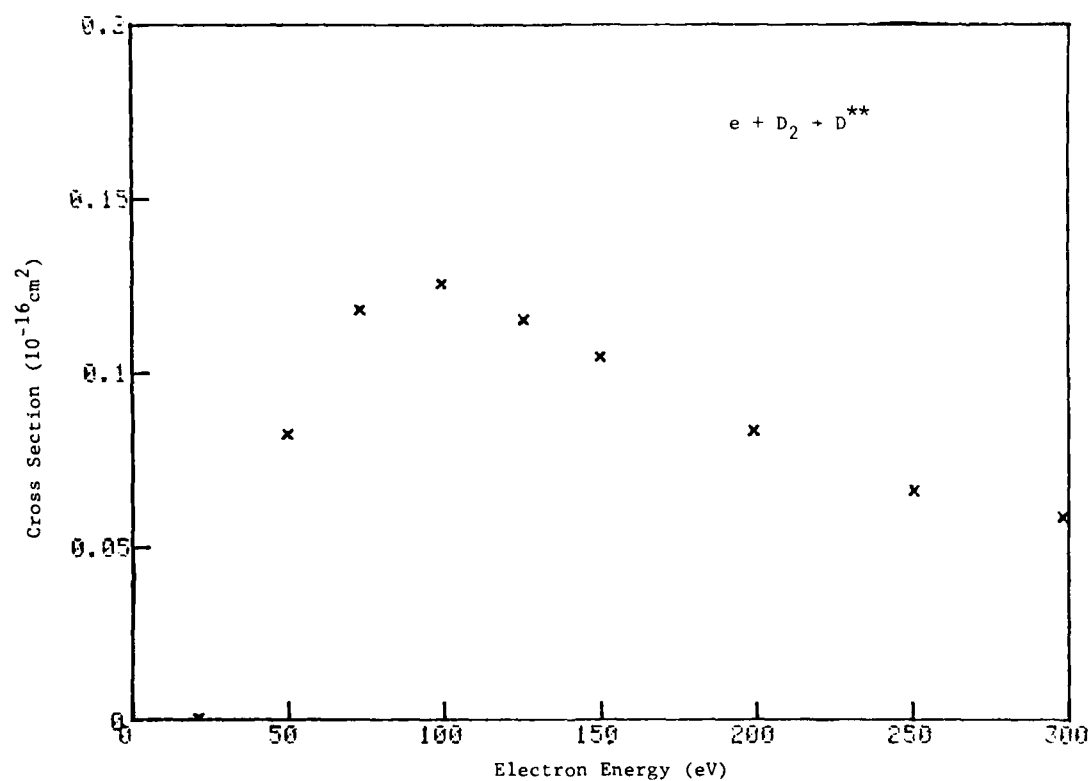


Reference: J. A. Schiavone, S. M. Tarr and R. S. Freund, J. Chem. Phys. 70,
4468 (1979).

Tabular and Graphical Data C-3.20. Cross sections for electron-impact dissociation of D_2 to form high-Rydberg fragments.

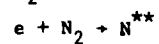


Electron Energy	Cross Section
eV	10^{-16} cm^2
21	0.00046
50	0.083
73	0.12
100	0.13
125	0.12
150	0.10
200	0.083
250	0.066
300	0.059

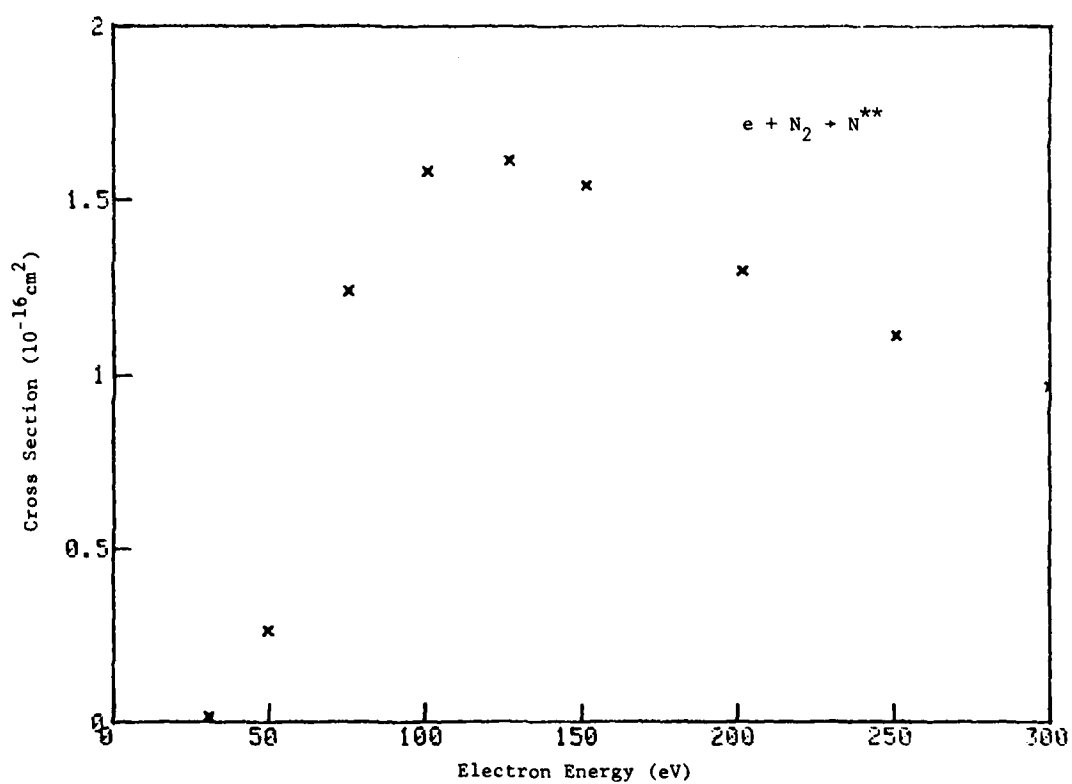


Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund. J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.21. Cross sections for electron-
impact dissociation of N_2 to form high-Rydberg fragments.

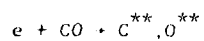


Electron Energy	Cross Section
eV	10^{-16} cm^2
31	0.012
50	0.26
76	1.2
100	1.6
125	1.6
150	1.5
200	1.3
250	1.1
300	0.96

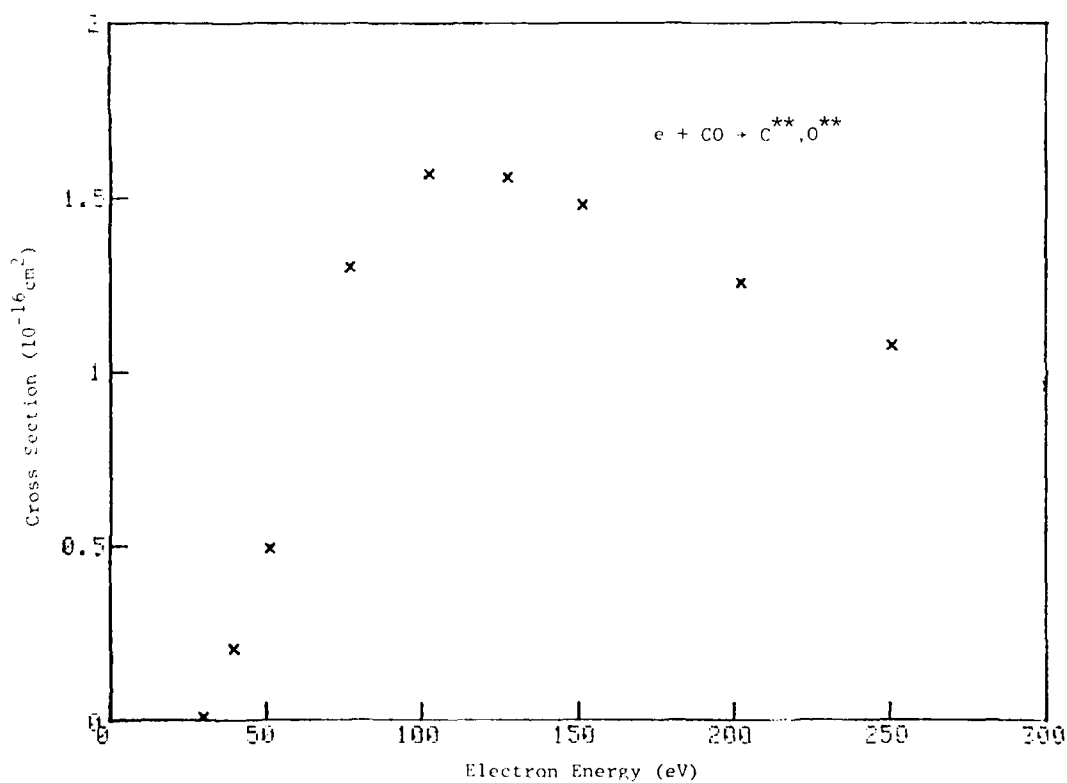


Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.22. Cross sections for electron-
impact dissociation of CO to form high-Rydberg fragments.

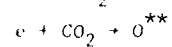


Electron Energy	Cross Section
eV	10^{-16} cm^2
30	0.0080
40	0.20
51	0.49
77	1.3
100	1.6
125	1.6
150	1.5
200	1.3
250	1.1
300	0.96

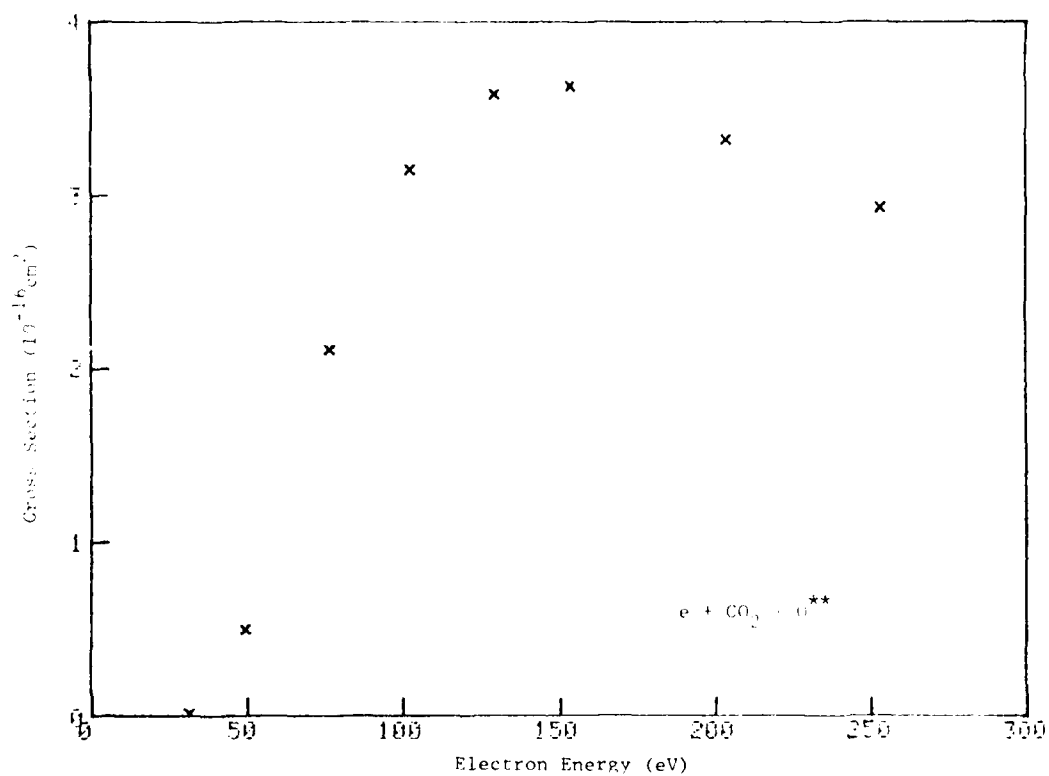


Reference: J. A. Schiavone, S.M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.23. Cross sections for electron-
impact dissociation of CO₂ to form high-Rydberg fragments.

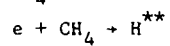


Electron Energy	Cross Section
eV	10 ⁻¹⁶ cm ²
31	0.016
49	0.50
77	2.1
100	3.2
125	3.6
150	3.6
200	3.3
250	2.9
300	2.6

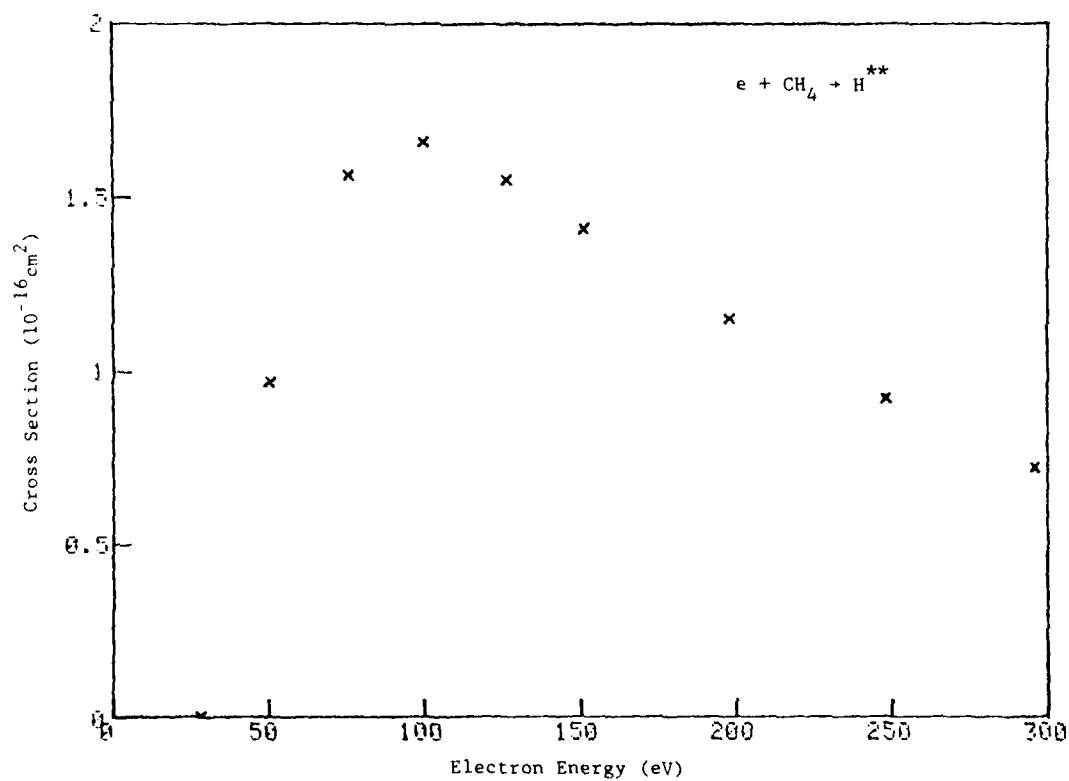


Reference: J. A. Schiavone, S.M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979)

Tabular and Graphical Data C-3.24. Cross sections for electron-
impact dissociation of CH_4 to form high-Rydberg fragments.

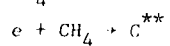


Electron Energy eV	Cross Section 10^{-16} cm^2
28	0.00045
50	0.97
76	1.6
100	1.7
125	1.6
150	1.4
200	1.2
250	0.92
300	0.72



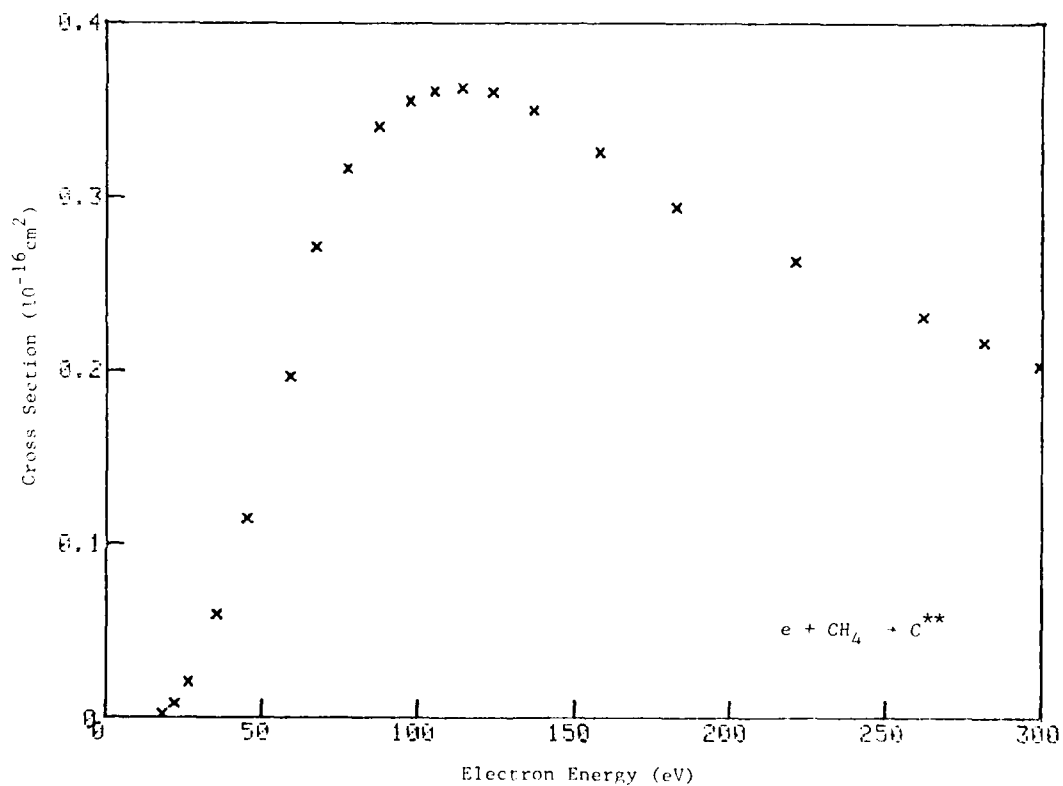
Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.25. Cross sections for electron-
impact dissociation of CH_4 to form high-Rydberg fragments.



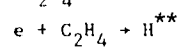
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
18	0.0018	110	0.36
22	0.0082	120	0.36
26	0.021	140	0.35
35	0.060	160	0.33
45	0.11	180	0.29
59	0.20	220	0.26
67	0.27	260	0.23
77	0.32	280	0.22
87	0.34	300	0.20
97	0.36		
100	0.36		

Cont. Next Column

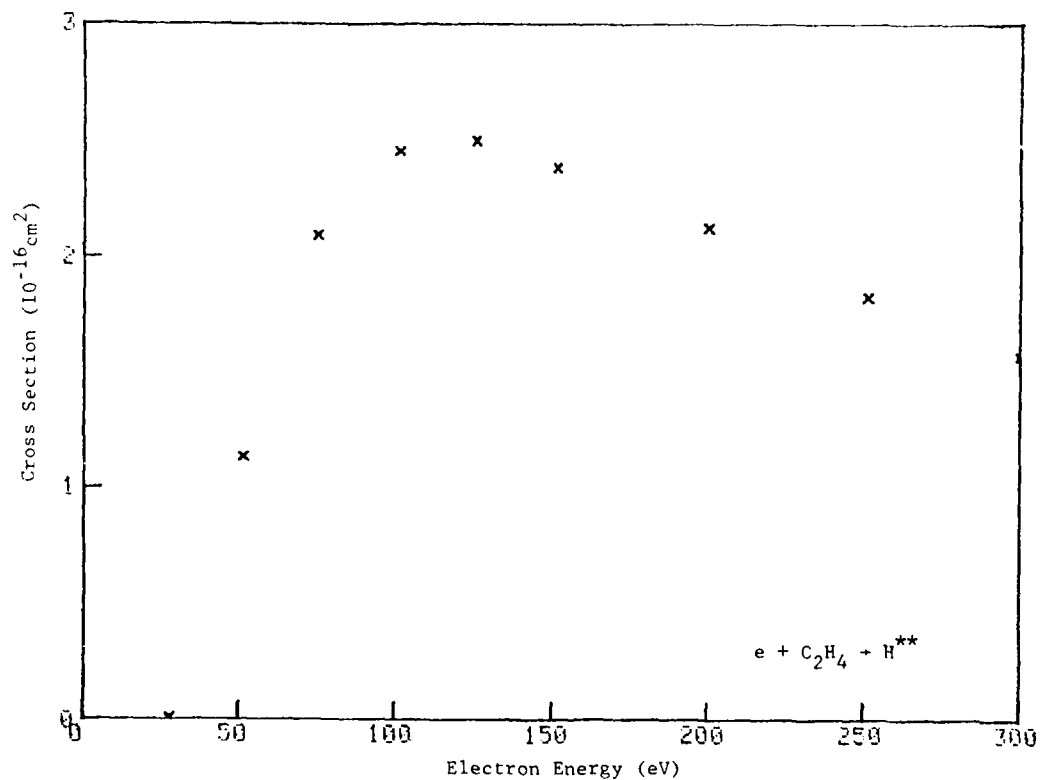


Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.26. Cross section for electron-
impact dissociation of C_2H_4 to form high-Rydberg fragments.

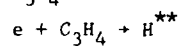


Electron Energy	Cross Section
eV	$10^{-16} cm^2$
28	0.0065
51	1.1
75	2.1
100	2.5
125	2.5
150	2.4
200	2.1
250	1.8
300	1.6

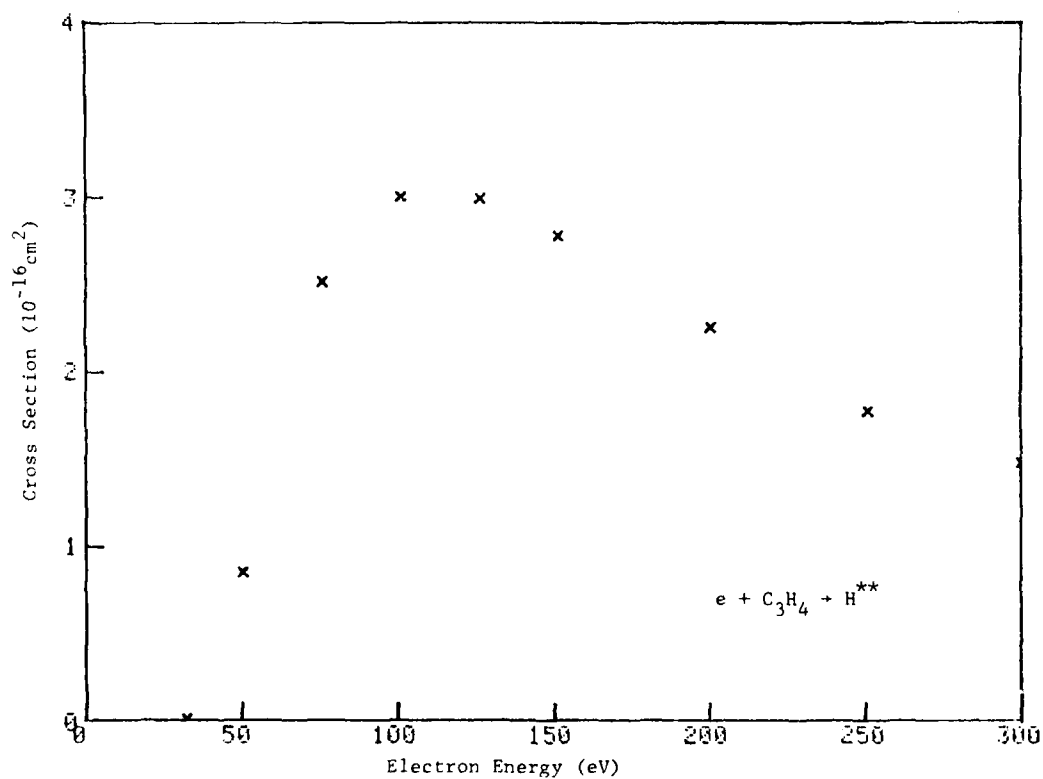


Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.27. Cross sections for electron-
impact dissociation of C_3H_4 to form high-Rydberg fragments.

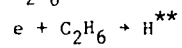


Electron Energy	Cross Section
eV	10^{-16} cm^2
32	0.0050
50	0.85
76	2.5
100	3.0
125	3.0
150	2.8
200	2.3
250	1.8
300	1.5

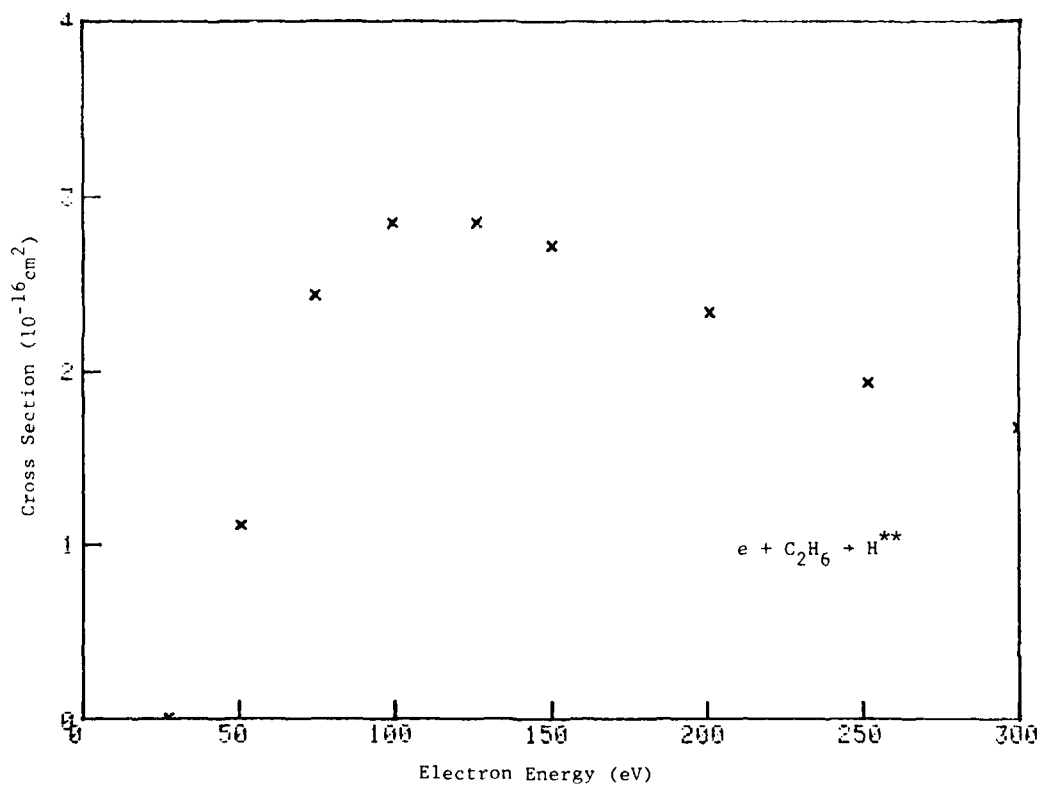


Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.28. Cross sections for electron-
impact dissociation of C_2H_6 to form high-Rydberg fragments.

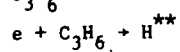


Electron Energy eV	Cross Section $10^{-16} cm^2$
27	0.0035
51	1.1
74	2.4
100	2.9
125	2.9
150	2.7
200	2.3
250	1.9
300	1.7

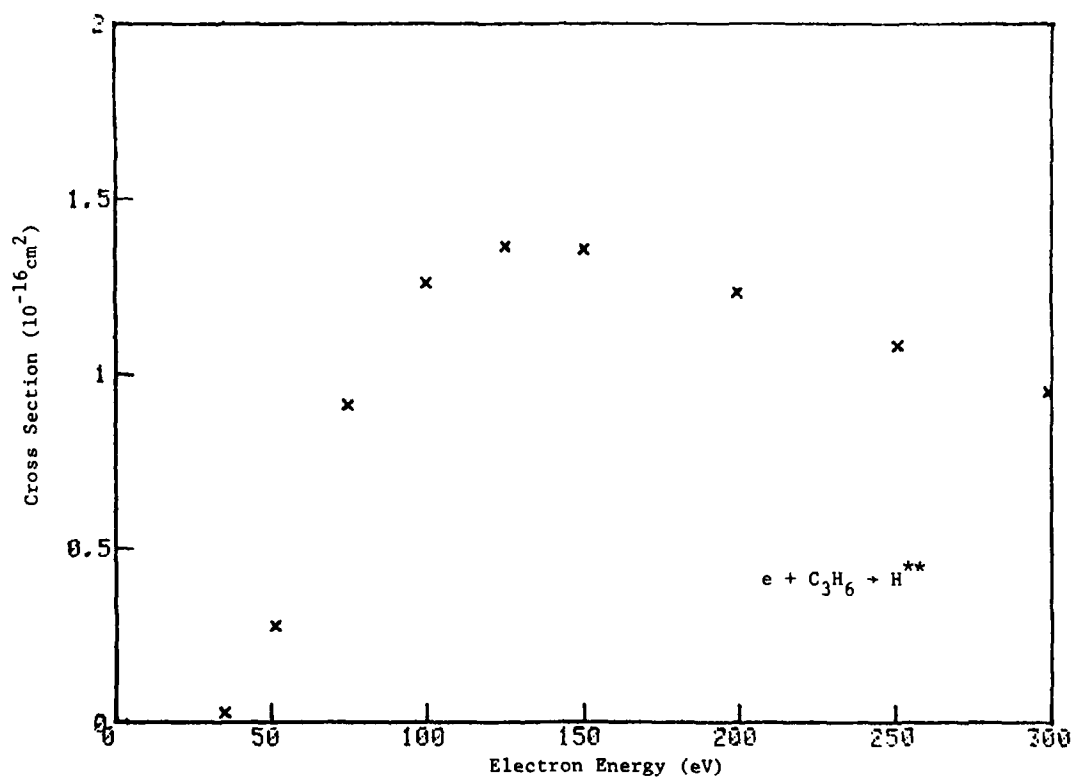


Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.29. Cross sections for electron-
impact dissociation of C_3H_6 to form high-Rydberg fragments.

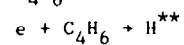


Electron Energy	Cross Section
eV	10^{-16} cm^2
35	0.024
51	0.27
75	0.91
100	1.3
125	1.4
150	1.4
200	1.2
250	1.1
300	0.95

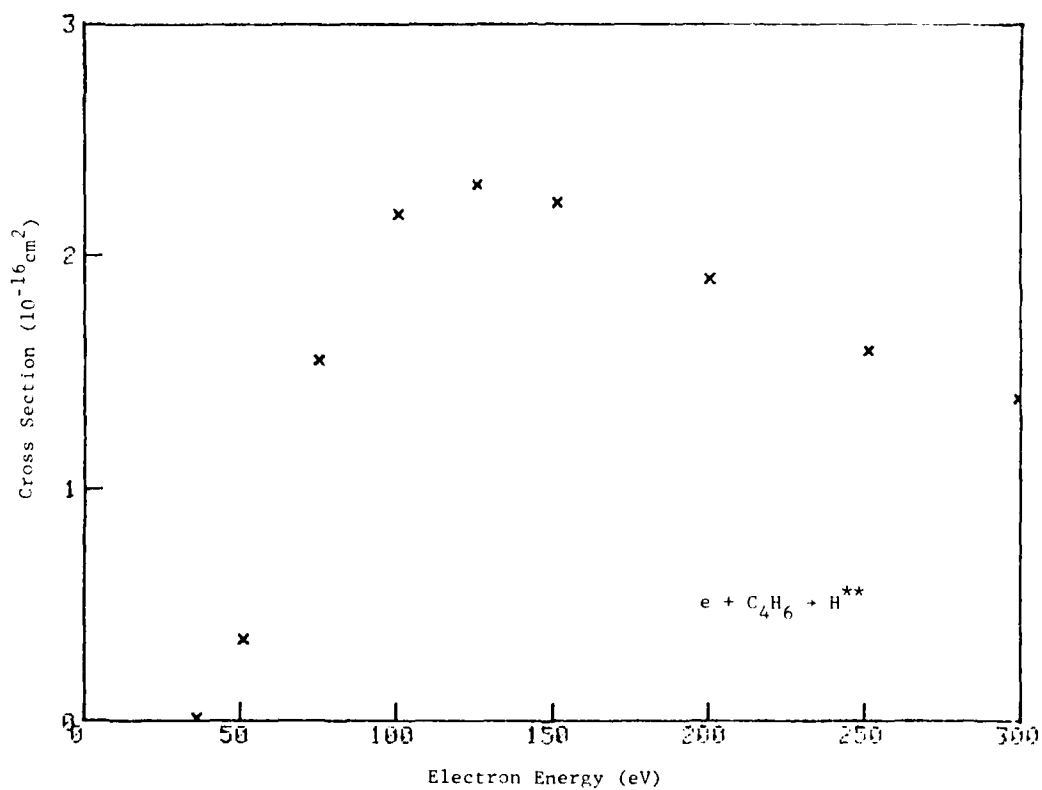


Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.30 Cross sections for electron-
impact dissociation of C_4H_6 to form high-Rydberg fragments.

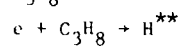


Electron Energy	Cross Section
eV	10^{-16} cm^2
36	0.0095
51	0.35
75	1.6
100	2.2
125	2.3
150	2.2
200	1.9
250	1.6
300	1.4

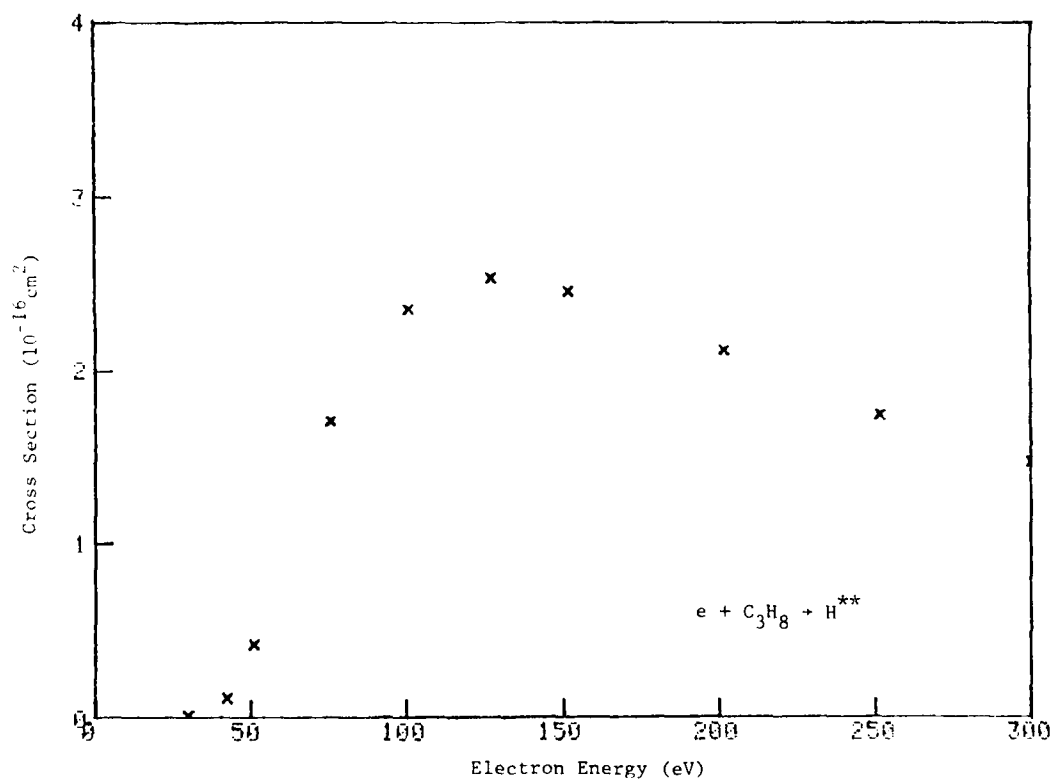


Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.31. Cross sections for electron-
impact dissociation of C_3H_8 to form high-Rydberg fragments.

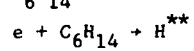


Electron Energy eV	Cross Section 10^{-16} cm^2
30	0.0076
42	0.11
51	0.42
76	1.7
100	2.4
125	2.5
150	2.5
200	2.1
250	1.7
300	1.5

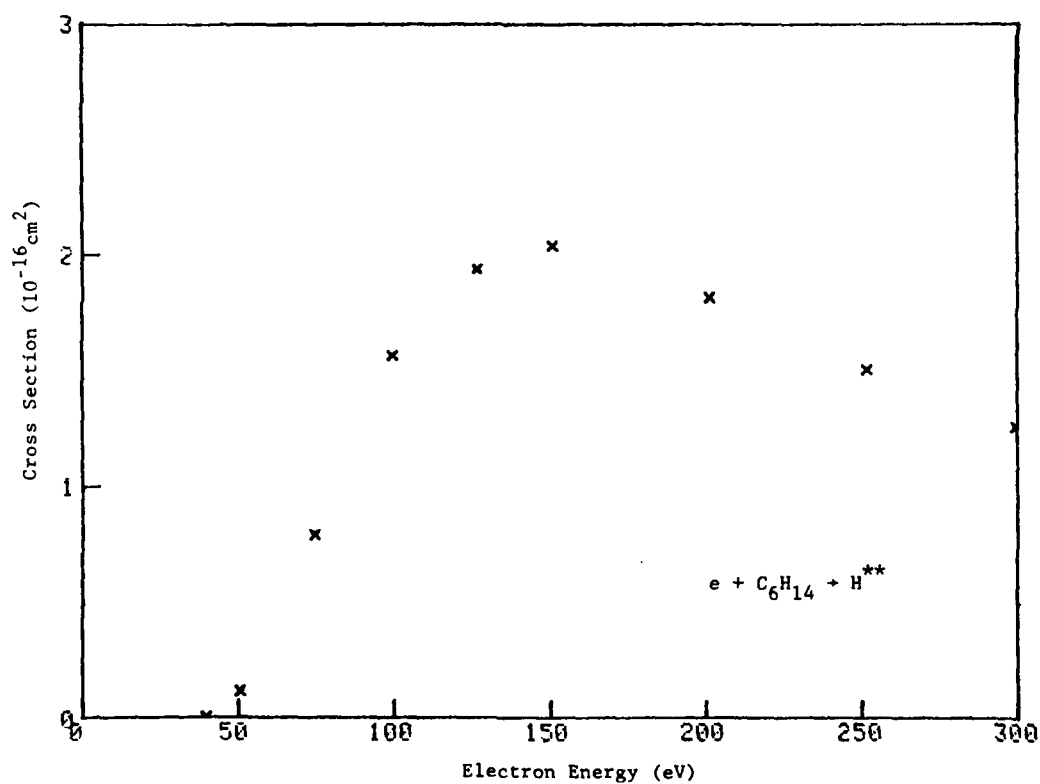


Reference: J. A. Schiavone, S. M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

Tabular and Graphical Data C-3.32. Cross sections for electron-
impact dissociation of C_6H_{14} to form high-Rydberg products.



Electron Energy	Cross Section
eV	10^{-16} cm^2
25	0.00
50	0.12
74	0.79
100	1.6
125	1.9
150	2.0
200	1.8
250	1.5
300	1.3



Reference: J. A. Schiavone, S.M. Tarr, and R. S. Freund, J. Chem. Phys. 70, 4468 (1979).

C-4. IONIZATION BY ELECTRON IMPACT

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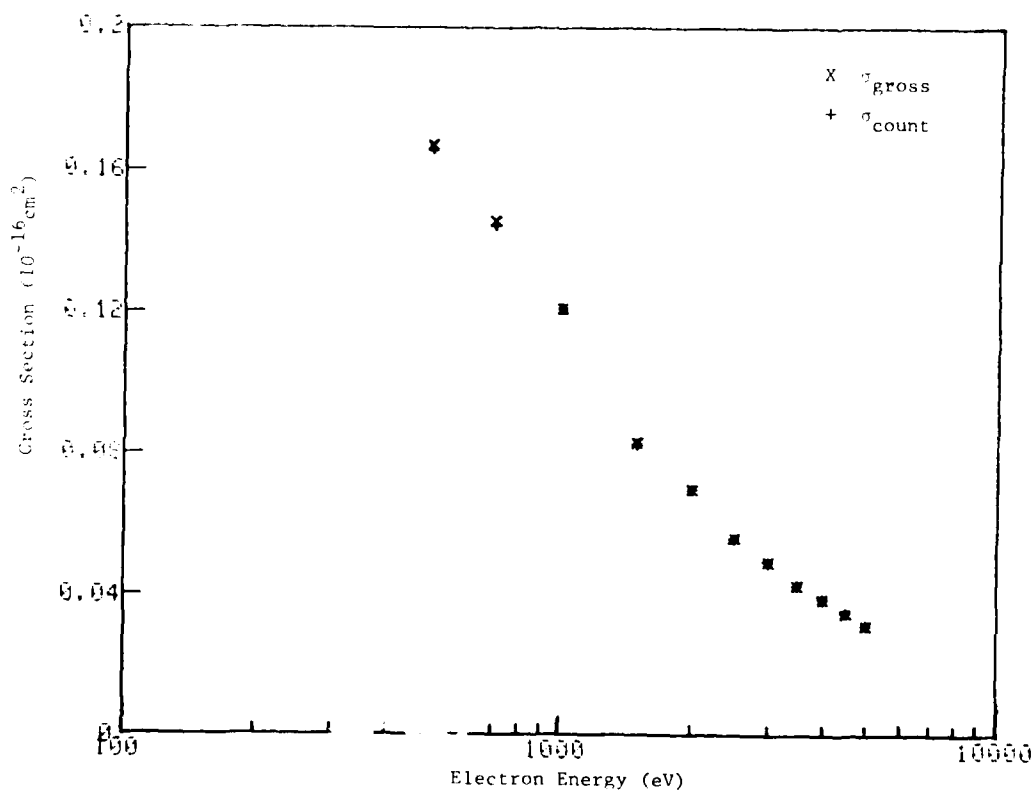
Tabular and Graphical C-4.1. Cross sections for electron impact ionization of He.

$$\sigma_{\text{gross}} = \sum_n \sigma_n$$

$$\sigma_{\text{count}} = \sum_n \sigma_n$$

σ_n = cross section for n-fold ionization

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
500	0.167	500	0.166
700	0.146	700	0.145
1000	0.121	1000	0.121
1500	0.0828	1500	0.0825
2000	0.0699	2000	0.0696
2500	0.0560	2500	0.0557
3000	0.0494	3000	0.0492
3500	0.0427	3500	0.0426
4000	0.0386	4000	0.0386
4500	0.0348	4500	0.0347
5000	0.0314	5000	0.0313



Reference: P. Nagy, A. Skutlartz, and V. Schmidt, J. Phys. B 13, 1249 (1980).

Tabular and Graphical Data C-4.2. Gross ionization cross sections
for electrons incident on Ne.

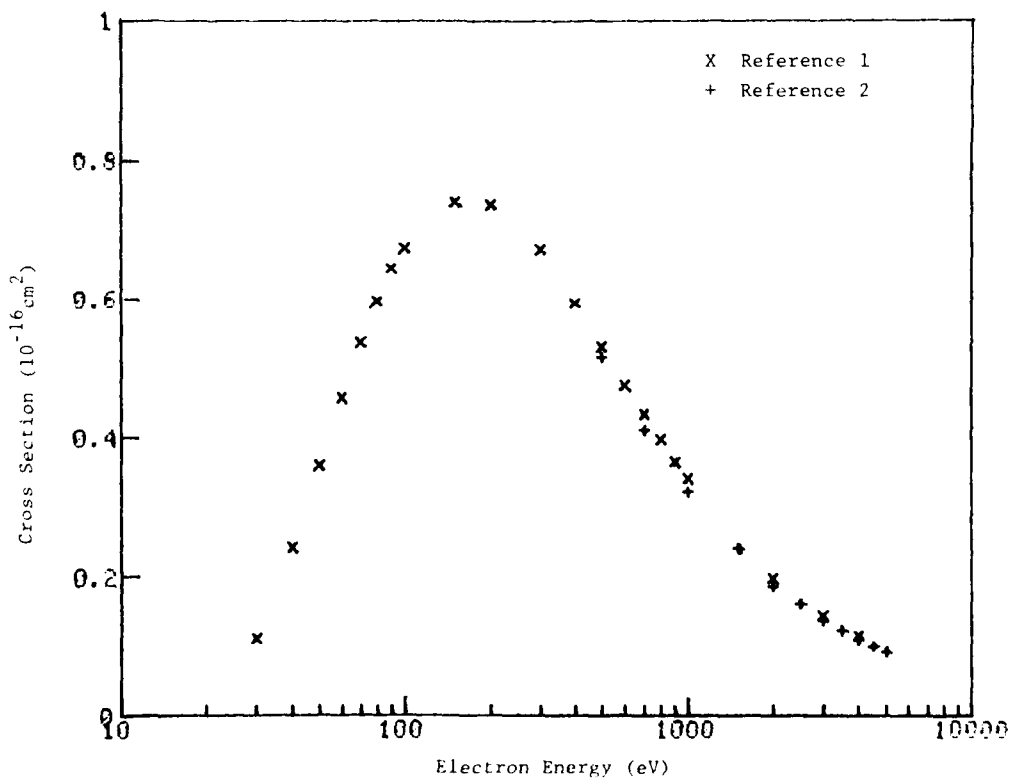
$$\sigma_{\text{gross}} = \sum_n \sigma_n$$

Ref. 1, Experimental average

Electron Energy eV	Cross Section 10^{-16} cm^2
30	0.1106
40	0.2421
50	0.3595
60	0.4575
70	0.5376
80	0.5978
90	0.6460
100	0.6745
150	0.7403
200	0.7381
300	0.6720
400	0.5947
500	0.5323
600	0.4757
700	0.4334
800	0.3973
900	0.3654
1000	0.3405
2000	0.1967
3000	0.1430
4000	0.1131

Ref. 2

Electron Energy eV	Cross Section 10^{-16} cm^2
500	0.515
700	0.409
1000	0.320
1500	0.238
2000	0.185
2500	0.158
3000	0.135
3500	0.120
4000	0.107
4500	0.0973
5000	0.0898



Reference 1: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12, 979 (1979).

Reference 2: P. Nagy, A. Skutlartz, and V. Schmidt, J. Phys. B 13, 1249 (1980).

Tabular and Graphical Data C-4.3. Gross ionization cross sections for
electrons incident on Ar.

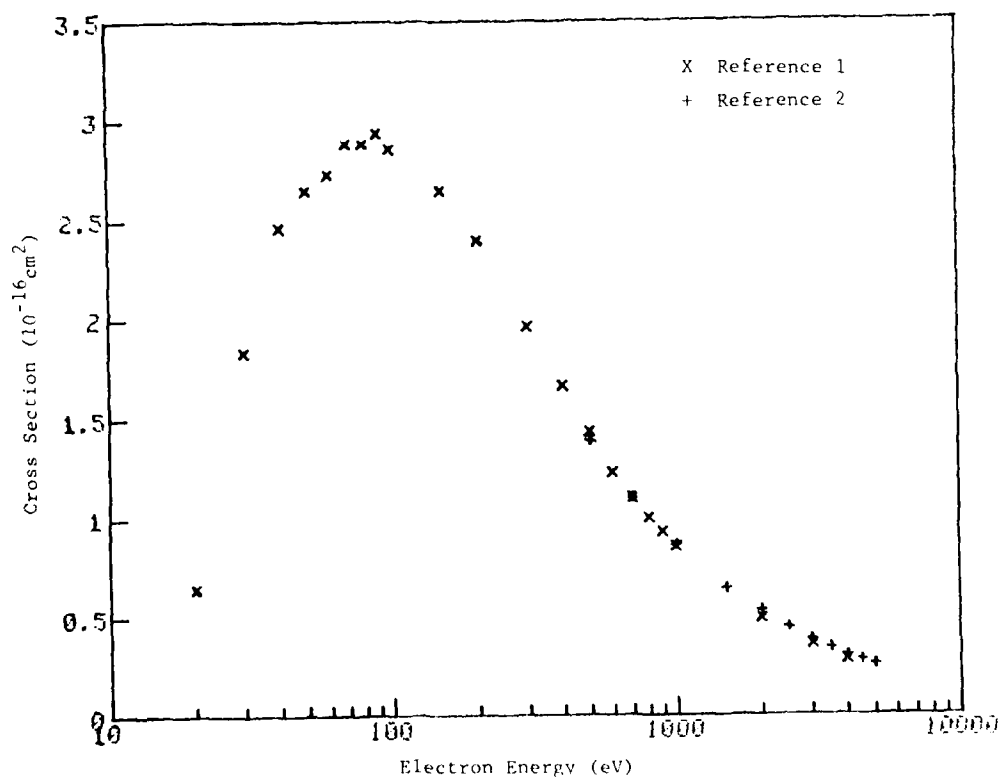
$$\sigma_{\text{gross}} = \sum_n n \sigma_n$$

Ref. 1, Experimental average

Electron Energy	Cross Section
eV	10^{-16} cm^2
20	0.6437
30	1.838
40	2.463
50	2.652
60	2.732
70	2.890
80	2.887
90	2.940
100	2.859
150	2.651
200	2.398
300	1.960
400	1.664
500	1.435
600	1.226
700	1.099
800	1.000
900	0.9232
1000	0.8501
2000	0.4875
3000	0.3548
4000	0.2759

Ref. 2

Electron Energy	Cross Section
eV	10^{-16} cm^2
500	0.515
700	0.409
1000	0.320
1500	0.238
2000	0.185
2500	0.158
3000	0.135
3500	0.120
4000	0.107
4500	0.0973
5000	0.0893



- Reference 1: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12, 979 (1979).
Reference 2: P. Nagy, A. Skutlartz, and V. Schmidt, J. Phys. B 13, 1249 (1980).

Tabular and Graphical Data C-4.4. Gross ionization cross sections for
electrons incident on Kr.

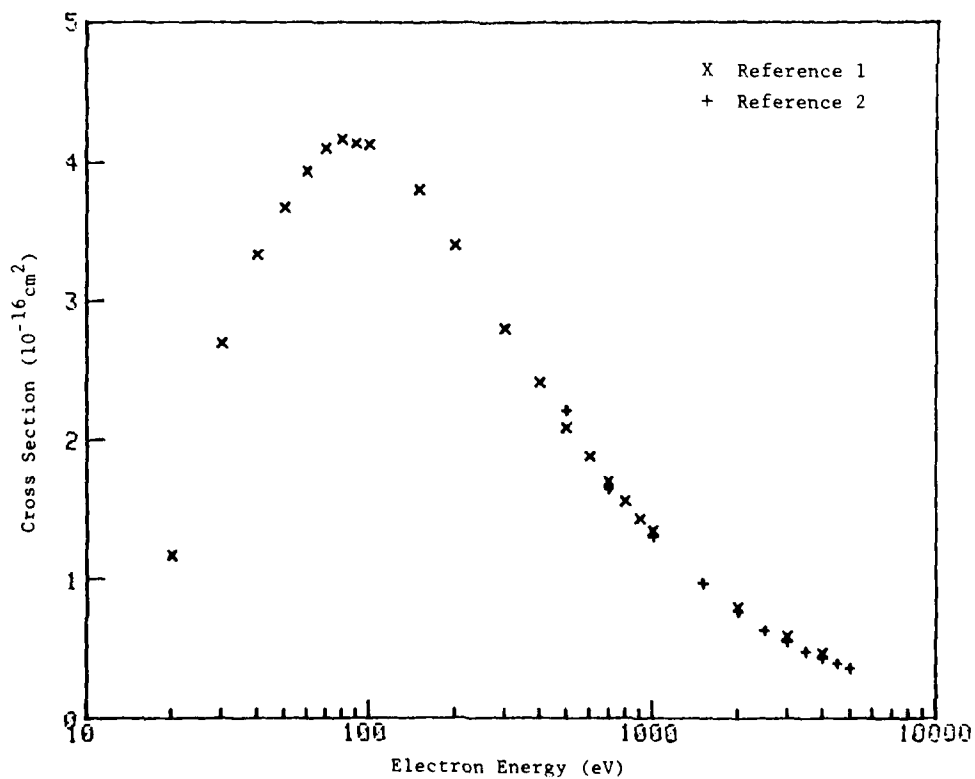
$$\sigma_{\text{gross}} = \sum_n \sigma_n$$

Ref. 1, Experimental average

Ref. 2

Electron Energy eV	Cross Section 10^{-16} cm^2
20	1.169
30	2.696
40	3.338
50	3.679
60	3.934
70	4.105
80	4.172
90	4.141
100	4.117
150	3.805
200	3.408
300	2.798
400	2.417
500	2.087
600	1.886
700	1.706
800	1.557
900	1.434
1000	1.342
2000	0.7913
3000	0.5866
4000	0.4640

Electron Energy eV	Cross Section 10^{-16} cm^2
500	2.21
700	1.65
1000	1.30
1500	0.961
2000	0.763
2500	0.620
3000	0.542
3500	0.474
4000	0.425
4500	0.388
5000	0.353



Reference 1. F. J. deHeer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B, 979 (1979).

Reference 2. P. Nagy, A. Skutlartz, and V. Schmidt, J. Phys. B 13, 1249 (1980).

Tabular and Graphical Data C-4.5. Gross ionization cross sections for
electrons incident on Xe.

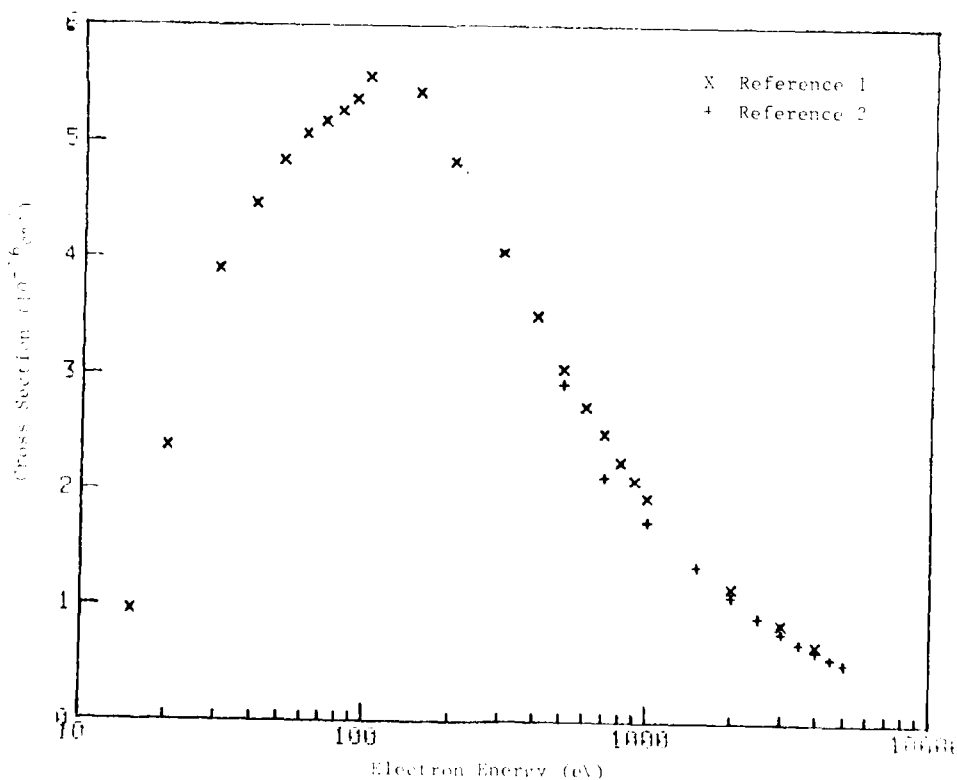
$$\sigma_{\text{gross}} = \sum_n \sigma_n$$

Ref. 1, Experimental average

Electron Energy	Gross Section
eV	10^{-16} cm^2
10	0.9554
20	2.369
30	3.906
40	4.477
50	4.847
60	5.076
70	5.186
80	5.275
90	5.377
100	5.461
120	5.446
140	4.850
160	4.057
180	3.505
200	3.055
250	2.725
300	2.457
350	2.248
400	2.096
450	1.941
500	1.754
520	0.8557
4500	0.6770

Ref. 2

Electron Energy	Gross Section
eV	10^{-16} cm^2
100	2.92
700	2.11
1000	1.73
1500	1.34
2000	1.09
2500	0.911
3000	0.761
3500	0.657
4000	0.600
4500	0.564
5000	0.516



Reference 1. E. J. deHeer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12, 979 (1979).

Reference 2. E. Nagy, A. Skutlartz, and V. Schmidt, J. Phys. B 13, 1749 (1980).

Tabular Data C-4.6. Ratio of count ionization to gross ionization

cross sections for noble gas atoms.

$$\sigma_{\text{gross}} = \sum_n \sigma_n$$

$$\sigma_{\text{count}} = \sum_n \sigma_n$$

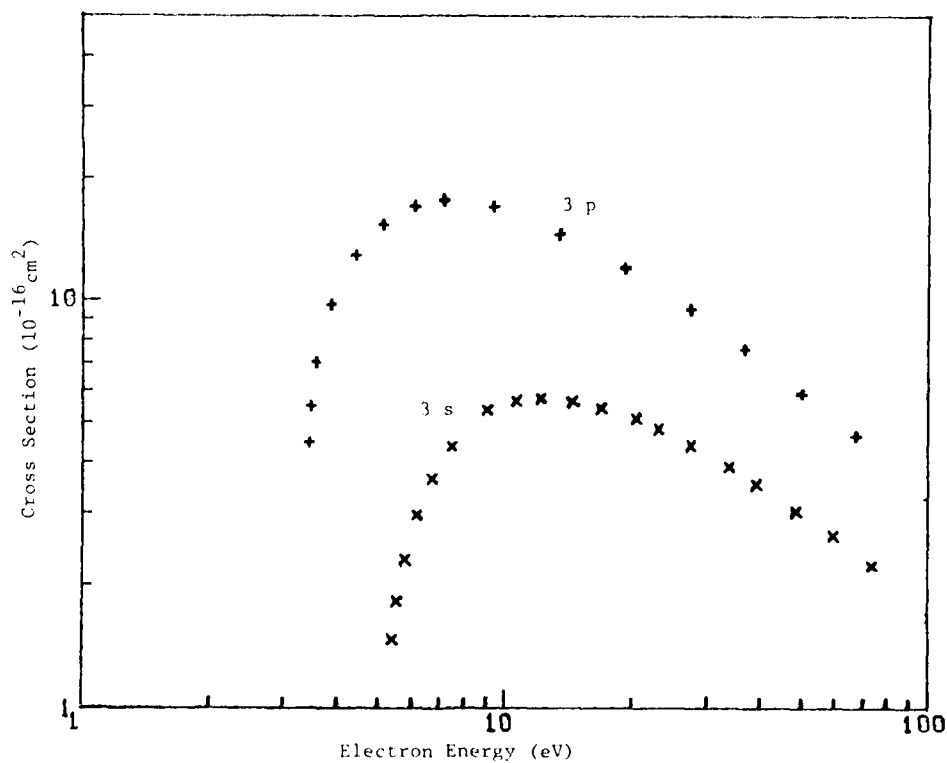
σ_n = cross section for n-fold ionization

Electron Energy (eV)	Ne	Ar	Kr	Xe
30	1	1		1.0
40	1	1	1.0	.97
50	1	1	.972	.93
60	1	.995	.924	.886
70	1	.986	.905	.873
80	.9927	.968	.90	.86
90	.9851	.959	.89	.85
100	.9771	.9498	.8817	.846
150	.9646	.9396	.887	.8510
200	.9548	.9353	.88	.8443
300	.9416	.9342	.869	.8337
400	.9420	.9305	.875	.8395
500	.9450	.9297	.86	.8276
600	.9524	.9280	.8667	.8195
700	.9527	.9278	.8662	.8068
800	.9548	.9275	.8478	.7914
900	.9553	.9277	.8629	.7925
1000	.9556	.9269	.8612	.7937
2000	.9526	.9225	.8519	.7814
3000	.9493	.9216	.8416	.7704
4000	.9586	.9197	.8299	.7582

Reference: F. J. de Heer, R. H. J. Jansen, and W. van der Kaay, J. Phys. B 12, 979 (1979).

Tabular and Graphical Data C-4.7. Calculated cross sections for electron-impact
ionization of excited states of Ne

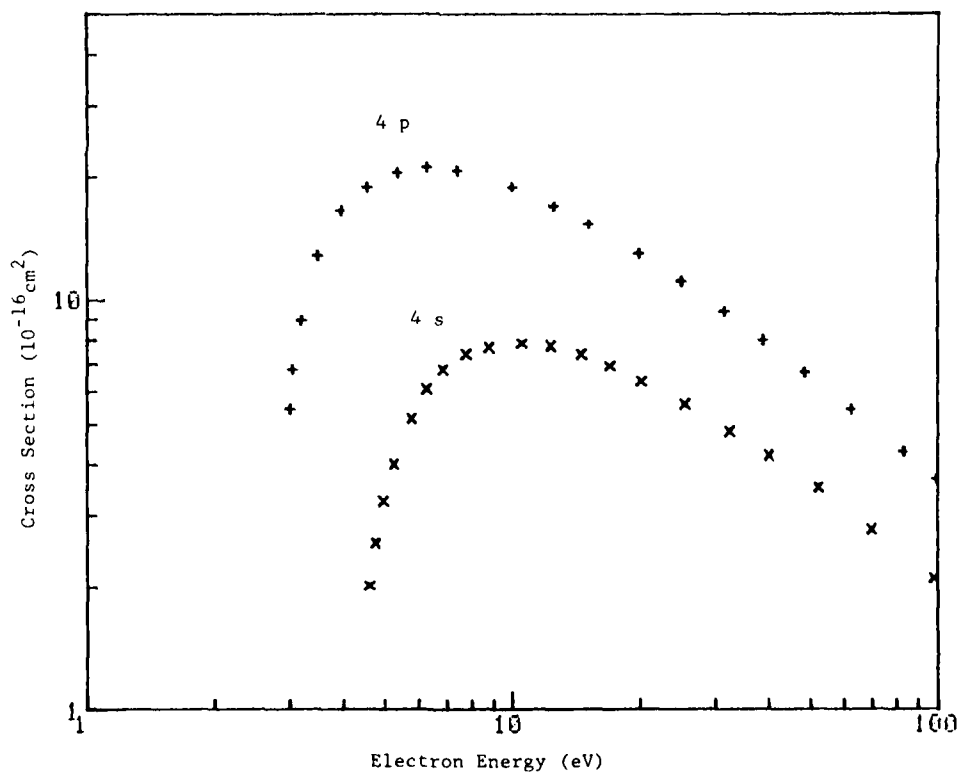
Ne(3s)		Ne(3p)	
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
5.4	1.48	3.4	4.47
5.5	1.82	3.5	5.51
5.8	2.31	3.6	7.06
6.2	2.96	3.9	9.77
6.7	3.65	4.4	12.9
7.5	4.36	5.1	15.4
9.1	5.41	6.1	17.0
11	5.68	7.1	17.6
12	5.76	9.3	17.1
14	5.68	13	14.6
17	5.47	19	12.0
20	5.14	28	9.55
23	4.85	37	7.60
28	4.41	50	5.90
34	3.91	67	4.66
39	3.54	100	3.28
49	3.06		
60	2.64		
74	2.23		
100	1.68		



Reference: H. A. Hyman, Phys. Rev. A 20, 855 (1979).

Tabular and Graphical Data C-4.8. Calculated cross sections for electron-impact ionization of excited states of Ar.

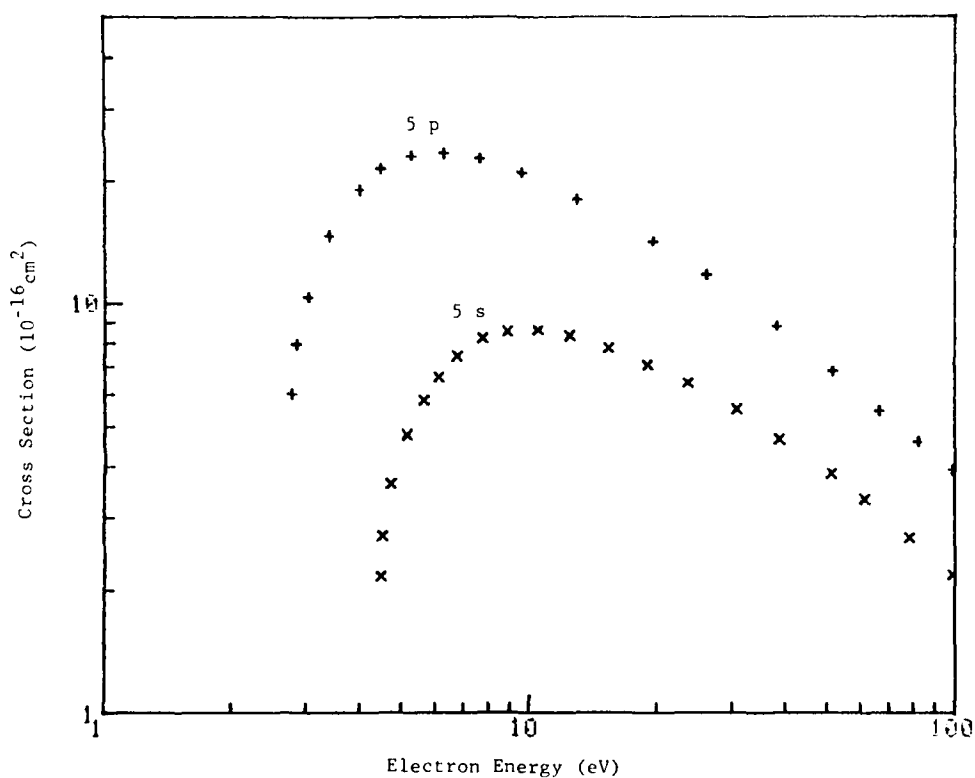
Ar(4s)		Ar(4p)	
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
4.6	2.01	3.0	5.43
4.8	2.55	3.0	6.81
5.0	3.24	3.2	8.93
5.3	3.98	3.5	12.8
5.8	5.16	3.9	16.6
6.3	6.08	4.5	19.0
6.9	6.78	5.4	20.6
7.8	7.39	6.3	21.2
8.8	7.71	7.4	20.8
10	7.88	10.0	19.0
12	7.74	12	17.0
15	7.41	15	15.4
17	6.94	20	13.1
20	6.37	25	11.1
25	5.59	31	9.38
32	4.81	39	8.02
40	4.19	49	6.68
52	3.50	62	5.42
70	2.76	83	4.28
98	2.10	99	3.67



Reference: H. A. Hyman, Phys. Rev. A 20, 855 (1979).

Tabular and Graphical Data C-4.9. Calculated cross sections for electron-impact ionization of excited states of Kr.

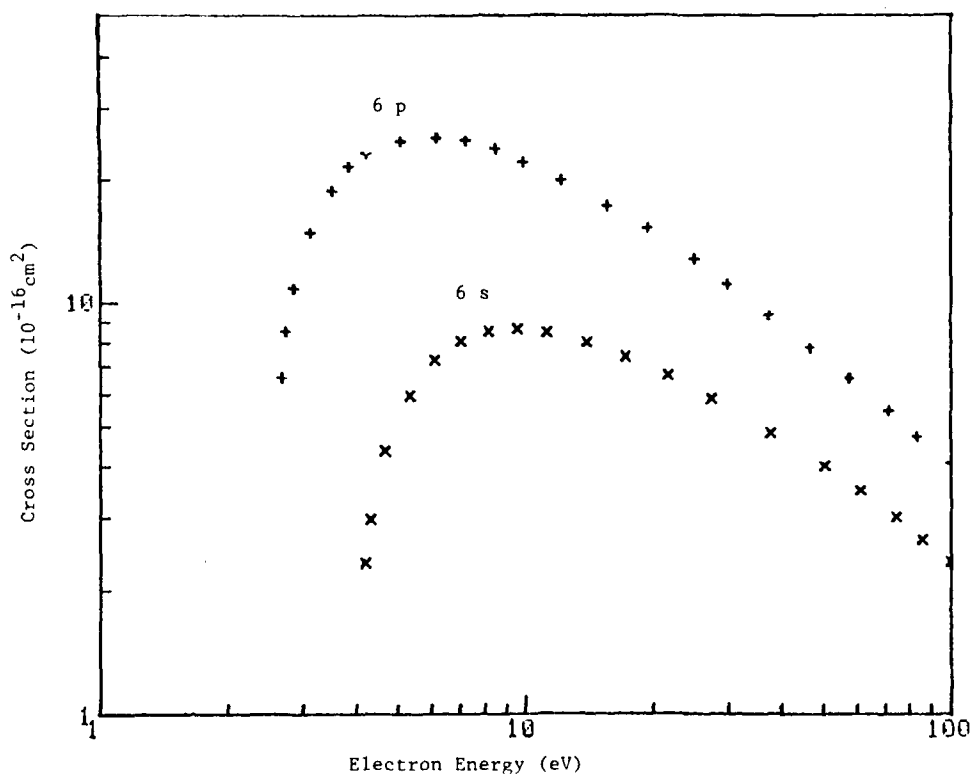
Kr(5s)		Kr(5p)	
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
4.5	2.16	2.8	6.01
4.5	2.70	2.8	7.93
4.7	3.63	3.0	10.4
5.2	4.79	3.4	14.7
5.6	5.81	4.0	19.1
6.1	6.63	4.4	21.5
6.7	7.44	5.2	23.1
7.7	8.26	6.3	23.5
8.9	8.58	7.6	22.8
10	8.65	9.6	21.0
12	8.36	13	18.0
15	7.79	20	14.3
19	7.11	26	11.8
24	6.39	38	8.80
31	5.51	52	6.85
39	4.68	67	5.45
51	3.84	82	4.58
62	3.32	99	3.92
79	2.67		
99	2.17		



Reference: H. A. Hyman, Phys. Rev. A 20, 855 (1979).

Tabular and Graphical Data C-4.10. Calculated cross sections for electron-impact ionization of excited states of Xe.

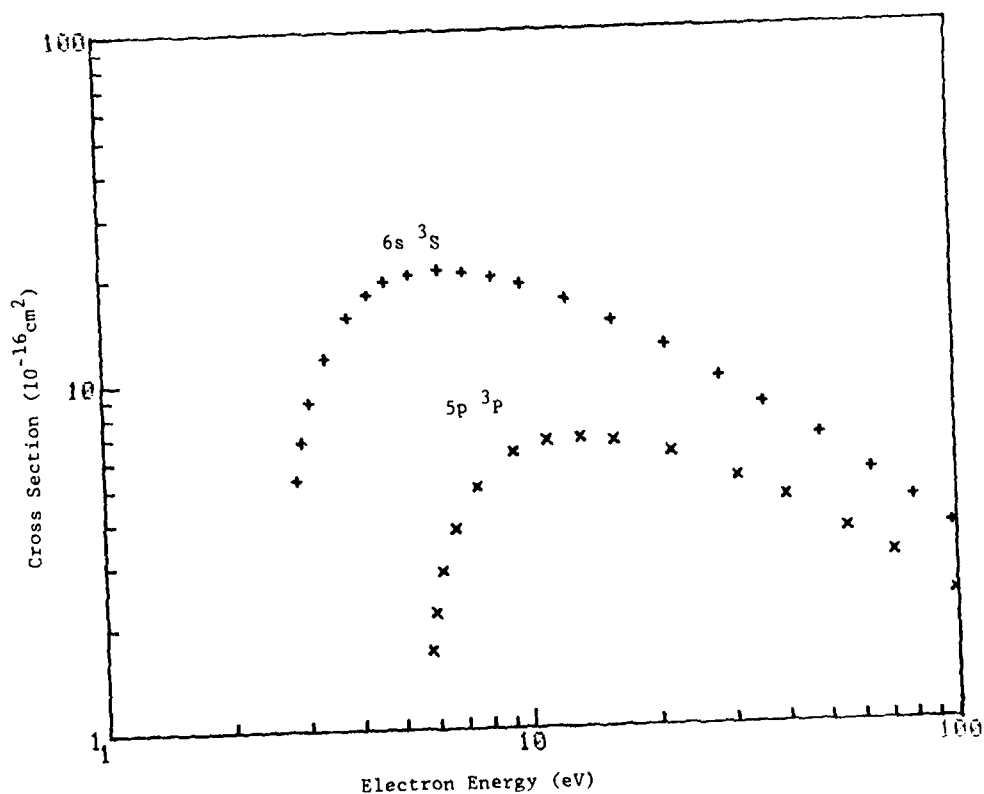
Xe(6p)		Xe(6p)	
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
4.2	2.32	2.7	6.54
4.3	2.97	2.7	8.49
4.7	4.35	2.8	10.9
5.3	5.96	3.1	14.9
6.1	7.27	3.5	18.9
7.0	8.12	3.8	21.7
8.2	8.58	4.2	23.3
9.6	8.68	5.1	25.0
11	8.50	6.2	25.5
14	8.05	7.3	25.0
17	7.42	8.5	24.0
22	6.69	9.9	22.2
27	5.84	12	20.0
38	4.82	16	17.4
50	4.00	19	15.3
61	3.49	25	12.8
74	3.02	30	11.2
86	2.65	37	9.31
100	2.34	47	7.76
		58	6.55
		71	5.46
		83	4.73
		100	4.07



Reference: H. A. Hyman, Phys. Rev. A 20, 855 (1979).

Tabular and Graphical Data C-4.11a. Calculated cross sections for electron-impact ionization of excited states of Cd.

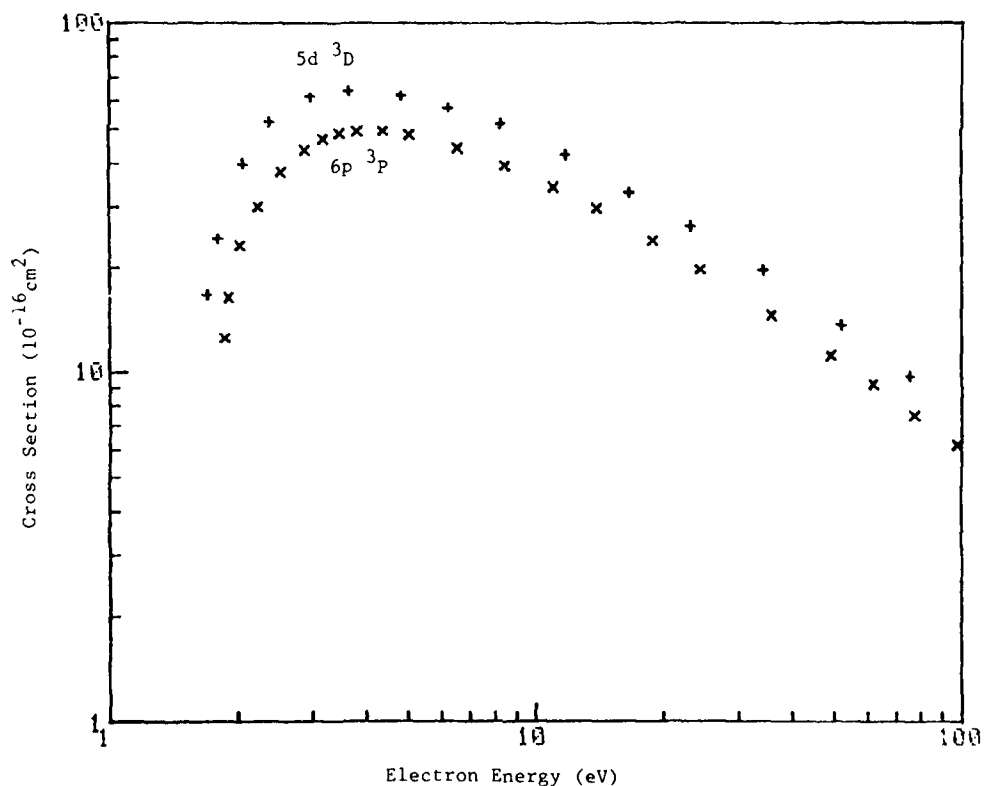
Cd(5p ³ P)		Cd(6s ³ S)	
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10 ⁻¹⁶ cm ²	eV	10 ⁻¹⁶ cm ²
5.8	1.67	2.8	5.21
5.9	2.14	2.9	6.71
6.2	2.84	3.1	8.72
6.7	3.75	3.3	11.6
7.5	4.94	3.8	15.3
9.2	6.20	4.2	17.6
11	6.67	4.7	19.2
13	6.72	5.3	20.2
16	6.59	6.2	20.6
22	6.07	7.2	20.3
31	5.12	8.4	19.6
40	4.47	9.7	18.7
55	3.61	12	16.8
71	3.04	16	14.5
99	2.34	21	12.3
		28	9.90
		36	8.26
		48	6.62
		63	5.29
		80	4.37
		98	3.63



Reference: H. A. Hyman, Phys. Rev. A 20, 855 (1979).

Tabular and Graphical Data C-4.11b. Calculated cross sections for electron-impact ionization of excited states of Cd.

Cd(6p ³ P)		Cd(5d ³ D)	
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10 ⁻¹⁶ cm ²	eV	10 ⁻¹⁶ cm ²
1.9	12.5	1.7	16.6
1.9	16.4	1.8	24.1
2.0	23.1	2.1	39.5
2.2	29.9	2.4	52.1
2.5	37.5	3.0	61.8
2.9	43.6	3.6	64.5
3.2	46.8	4.8	62.3
3.5	48.6	6.2	57.7
3.8	49.3	8.2	52.0
4.4	49.5	12	42.4
5.1	48.2	17	32.8
6.6	44.1	23	26.5
8.4	39.3	34	19.6
11	34.2	52	13.6
14	29.6	76	9.69
19	24.0	100	7.25
24	19.7		
36	14.5		
50	11.2		
62	9.17		
78	7.49		
98	6.18		



Reference: H. A. Hyman, Phys. Rev. A 20, 855 (1979).

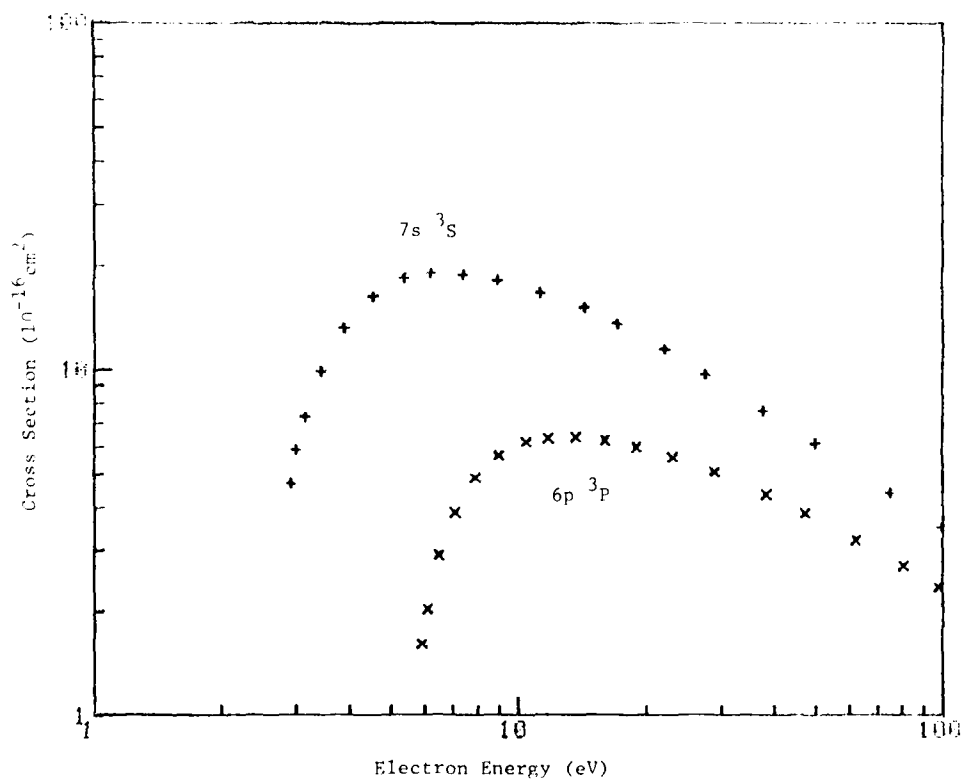
Tabular Data C-4.12. Cross Sections for electron impact
ionization of Hg.

Electron Energy eV	Cross Section 10^{-16} cm^2
300	3.14
400	2.66
500	2.32

Reference: K. Jost and B. Ohnemus, Phys. Rev. A 19
611 (1979).

Tabular and Graphical Data C-4.13a. Calculated cross sections for electron-impact ionization for excited states of Hg.

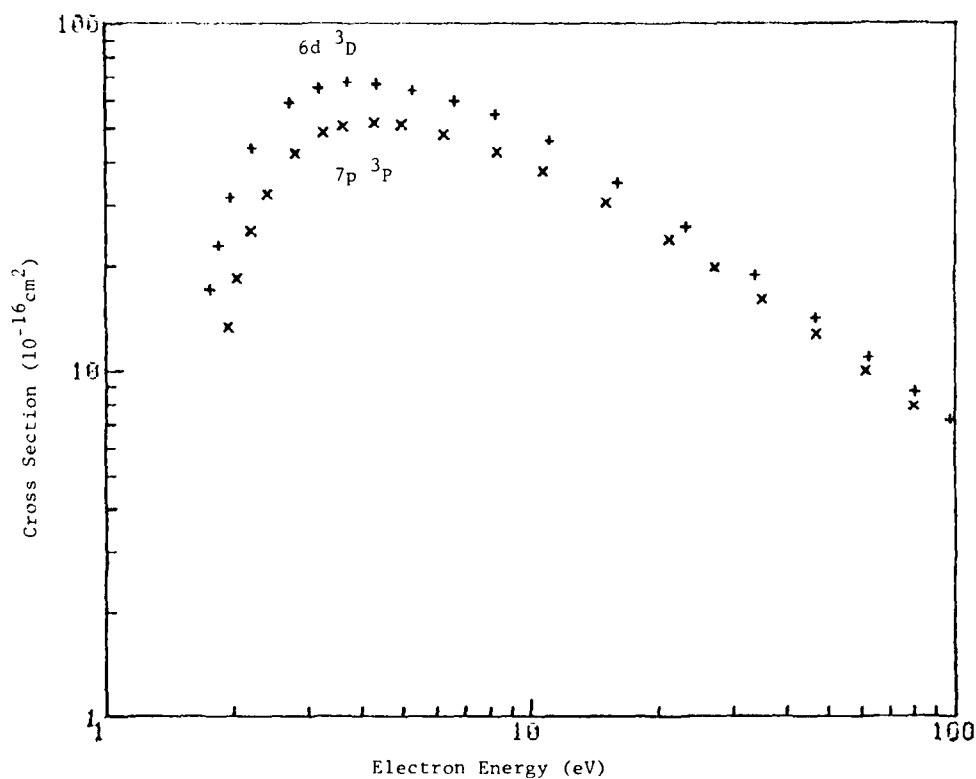
Hg(6p ^3P)		Hg(7s ^3S)	
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
5.9	1.61	2.9	4.72
6.1	2.03	3.0	5.86
6.5	2.92	3.1	7.34
7.1	3.88	3.4	9.87
7.9	4.90	3.9	13.2
9.0	5.68	4.6	16.3
10	6.20	5.4	18.4
12	6.37	6.2	19.0
14	6.39	7.4	18.8
16	6.27	9.0	18.1
19	5.99	11	16.8
23	5.59	14	15.1
29	5.07	17	13.6
38	4.37	22	11.5
47	3.85	28	9.70
62	3.21	38	7.60
81	2.71	50	6.15
93	2.36	75	4.40
		99	3.50



Reference: H. A. Hyman, Phys. Rev. A 20, 855 (1979).

Tabular and Graphical Data C-4.13b. Calculated cross sections for electron-impact ionization for excited states of Hg.

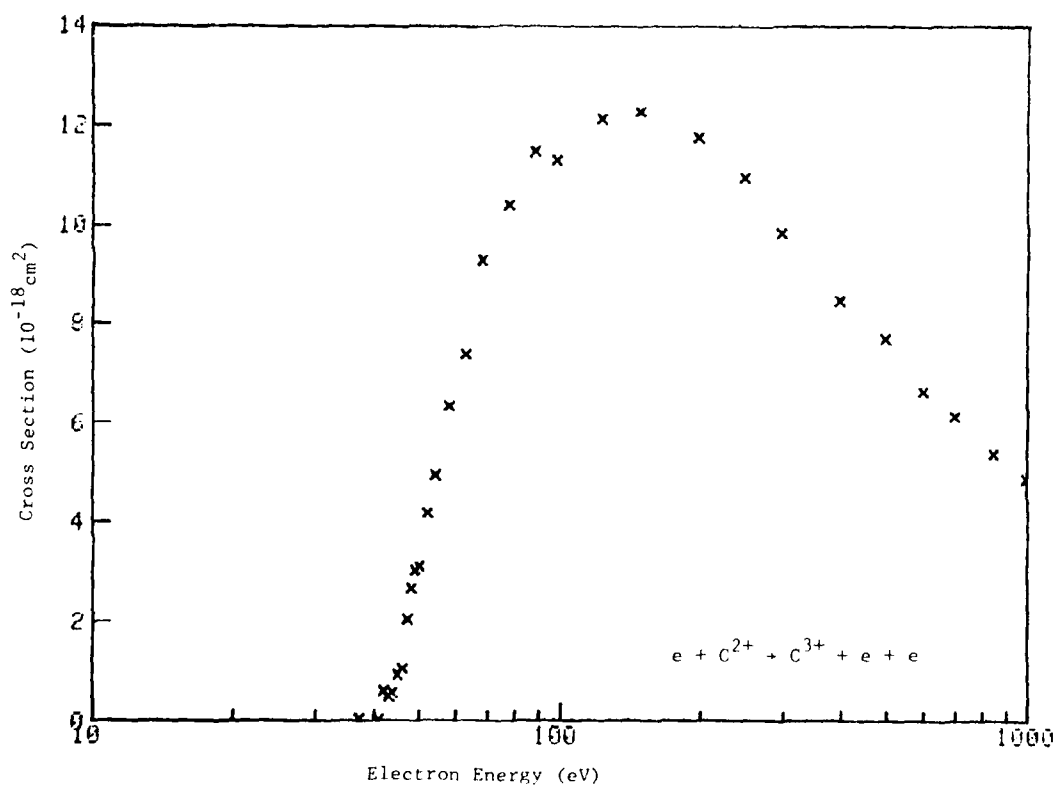
Hg(7p ³ p)		Hg(6d ³ D)	
Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10 ⁻¹⁶ cm ²	eV	10 ⁻¹⁶ cm ²
1.9	13.3	1.8	17.1
2.0	18.4	1.8	22.8
2.2	25.2	2.0	31.5
2.4	32.2	2.2	43.9
2.8	42.1	2.7	59.1
3.3	48.6	3.2	65.4
3.6	50.7	3.7	68.0
4.3	52.0	4.3	67.1
5.0	51.2	5.3	64.3
6.3	48.1	6.6	59.6
8.3	42.6	8.2	54.6
11	37.2	11	46.0
15	30.3	16	34.7
21	23.8	23	25.8
27	19.7	34	18.8
35	16.0	47	14.0
47	12.7	63	10.9
61	9.95	81	8.67
80	7.90	98	7.22
100	6.36		



Reference: H. A. Hyman, Phys. Rev. A 20, 855 (1979).

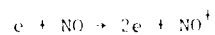
Tabular and Graphical Data C-4.14. Cross sections for electron-impact
ionization of C^{2+} .
 $e + C^{2+} \rightarrow C^{3+} + e + e$

Electron Energy eV	Cross Section 10^{-18} cm^2	Electron Energy eV	Cross Section 10^{-18} cm^2
37.1	.02	68	9.32
41	.00	78	10.42
42	.58	88	11.52
43	.47	98	11.33
44	.57	123	12.16
45	.92	148	12.31
46	1.04	198	11.78
47	2.05	248	10.98
48	2.66	298	9.85
49	3.03	398	8.50
50	3.09	498	7.70
52	4.18	598	6.64
54	4.95	698	6.13
58	6.35	848	5.38
63	7.40	998	4.87



Reference: P. R. Woodruff, M. C. Hublet, M. F. A Harrison, and E. Brook,
J. Phys. B 11, L679 (1978).

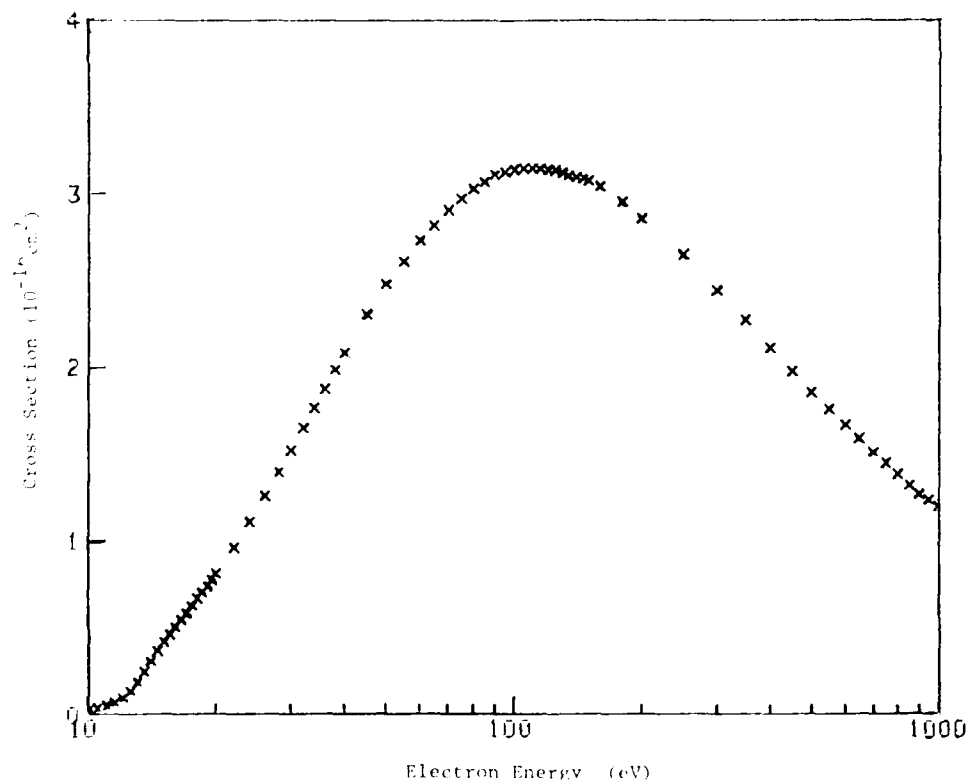
Tabular and Graphical Data C-4.15. Cross sections for electron-impact ionization of NO.



Electron Energy	Cross Section	Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2	eV	10^{-16} cm^2
145	3.116	50.0	1.57	145	3.09
150	3.117	50.0	1.65	150	3.06
160	3.053	54.0	1.77	160	3.04
170	3.040	56.0	1.87	170	2.96
180	3.040	58.0	1.99	200	2.86
190	3.054	60.0	2.09	250	2.65
200	3.131	65.0	2.31	300	2.45
250	3.159	70.0	2.48	350	2.27
300	3.188	75.0	2.61	400	2.11
350	3.305	80.0	2.74	450	1.96
400	3.369	85.0	2.82	500	1.86
450	3.435	90.0	2.91	550	1.76
500	3.461	95.0	2.97	600	1.67
550	3.507	100.0	3.04	650	1.59
600	3.545	105.0	3.07	700	1.51
650	3.585	110.0	3.12	750	1.45
700	3.605	115.0	3.15	800	1.39
750	3.606	120.0	3.14	850	1.33
800	3.591	125.0	3.14	900	1.28
850	3.571	130.0	3.12	950	1.24
900	3.554	135.0	3.11	1000	1.21
950	3.540	140.0	3.10		

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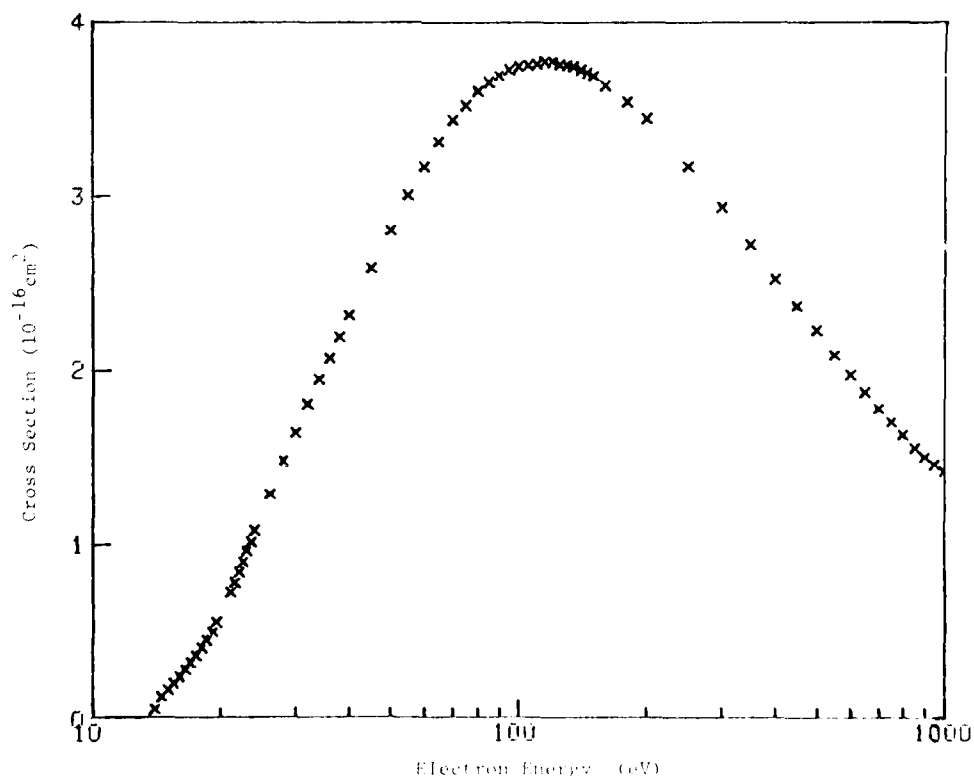
Reference: D. Rapp and P. Englander-Golden, J. Chem. Phys. 43, 1464 (1965).

Tabular and Graphical Data C-4.16. Cross sections for electron-impact ionization of N_2O .
 $e + N_2O \rightarrow 2e + N_2O^+$

Electron Energy	Cross Section	Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2	eV	10^{-16} cm^2
14.0	0.0537	40.0	0.31	200	1.15
14.5	0.111	45.0	0.59	300	1.44
15.0	0.153	50.0	0.81	350	1.71
15.5	0.199	55.0	1.01	400	2.03
16.0	0.233	60.0	1.16	450	2.37
16.5	0.260	65.0	1.30	500	2.69
17.0	0.311	70.0	1.44	550	2.99
17.5	0.302	75.0	1.53	600	3.27
18.0	0.404	80.0	1.61	650	3.53
18.5	0.451	85.0	1.66	700	3.73
19.0	0.496	90.0	1.70	750	3.76
19.5	0.551	95.0	1.73	800	3.75
20.0	0.743	100	1.75	850	3.75
21.0	0.781	105	1.76	900	3.71
22.0	0.840	110	1.77	950	3.70
23.0	0.890	115	1.75	1000	3.64
23.5	0.959	120	1.76		
24.0	1.01	125	1.77		
24.5	1.07	130	1.76		
25.0	1.23	135	1.75		
26.0	1.47	140	1.73		
27.0	1.64	145	1.71		
28.0	1.80	150	1.70		
29.0	1.94	160	1.64		
30.0	2.07	180	1.55		
35.0	2.19	200	1.45		

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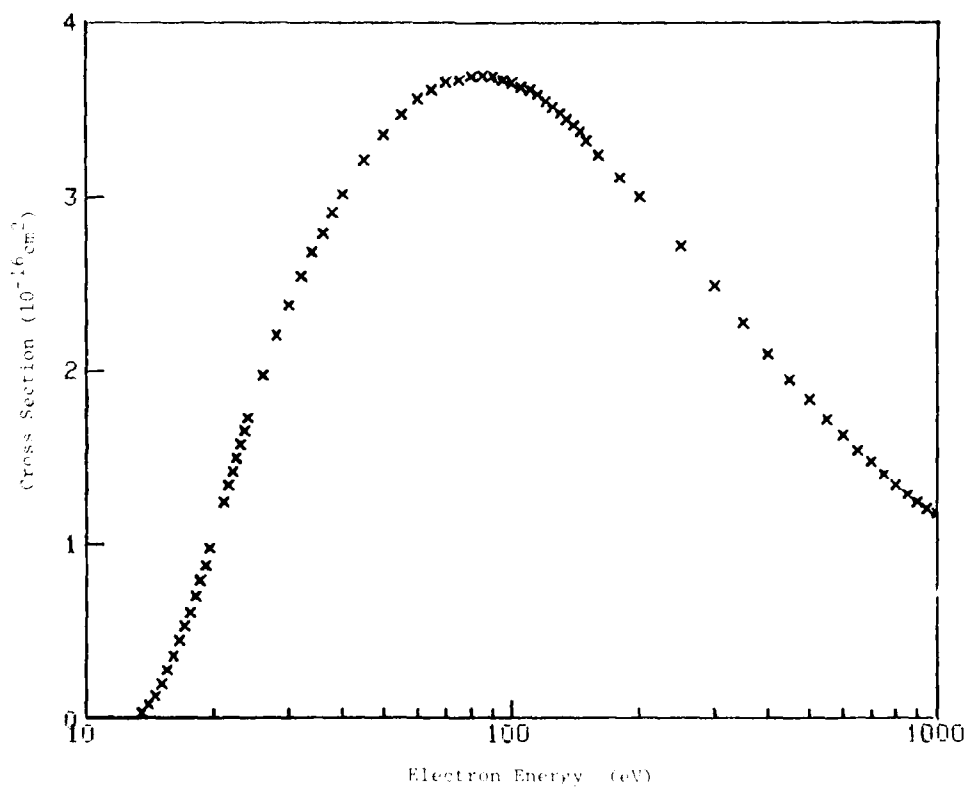
Reference: D. Rapp and P. Englander-Golden, J. Chem. Phys. 43, 1464 (1965)

Tabular and Graphical Data C-4.17. Cross sections for electron-impact ionization of CH_4 .

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
13.5	0.0343	33.0	2.91	200	3.01
14.0	0.0739	40.0	3.02	250	2.72
14.5	0.130	45.0	3.21	300	2.49
15.0	0.190	50.0	3.36	350	2.27
15.5	0.275	55.0	3.48	400	2.09
16.0	0.361	60.0	3.56	450	1.94
16.5	0.445	65.0	3.62	500	1.83
17.0	0.531	70.0	3.65	550	1.72
17.5	0.610	75.0	3.68	600	1.63
18.0	0.700	80.0	3.70	650	1.54
18.5	0.793	85.0	3.70	700	1.47
19.0	0.880	90.0	3.70	750	1.40
19.5	0.977	95.0	3.68	800	1.34
21.0	1.24	100	3.66	850	1.28
21.5	1.34	105	3.63	900	1.24
22.0	1.42	110	3.62	950	1.21
22.5	1.50	115	3.59	1000	1.18
23.0	1.58	120	3.56		
23.5	1.65	125	3.52		
24.0	1.72	130	3.48		
25.0	1.77	135	3.45		
26.0	1.80	140	3.41		
30.0	2.38	145	3.38		
32.0	2.54	150	3.33		
34.0	2.68	160	3.25		
36.0	2.79	180	3.12		

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Reference: U. Rapp and P. Englander-Golden, J. Chem. Phys. 43, 1464 (1965).

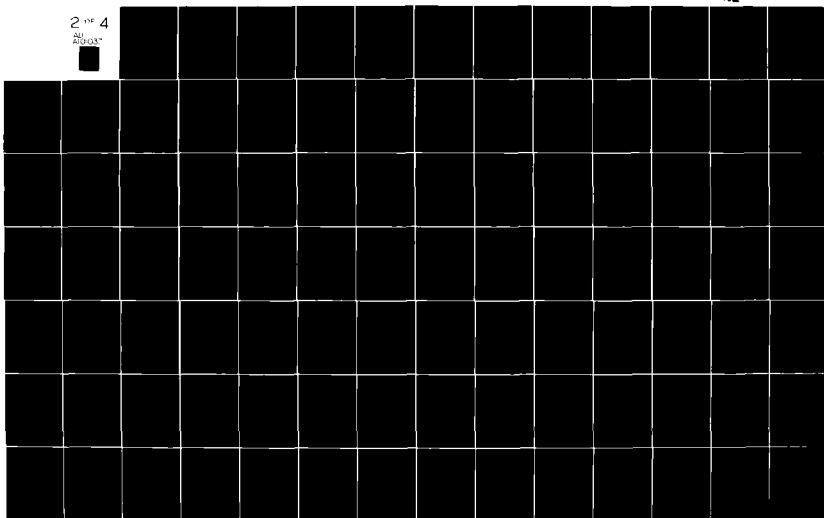
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DRSMI-RH-81-4-VOL-8

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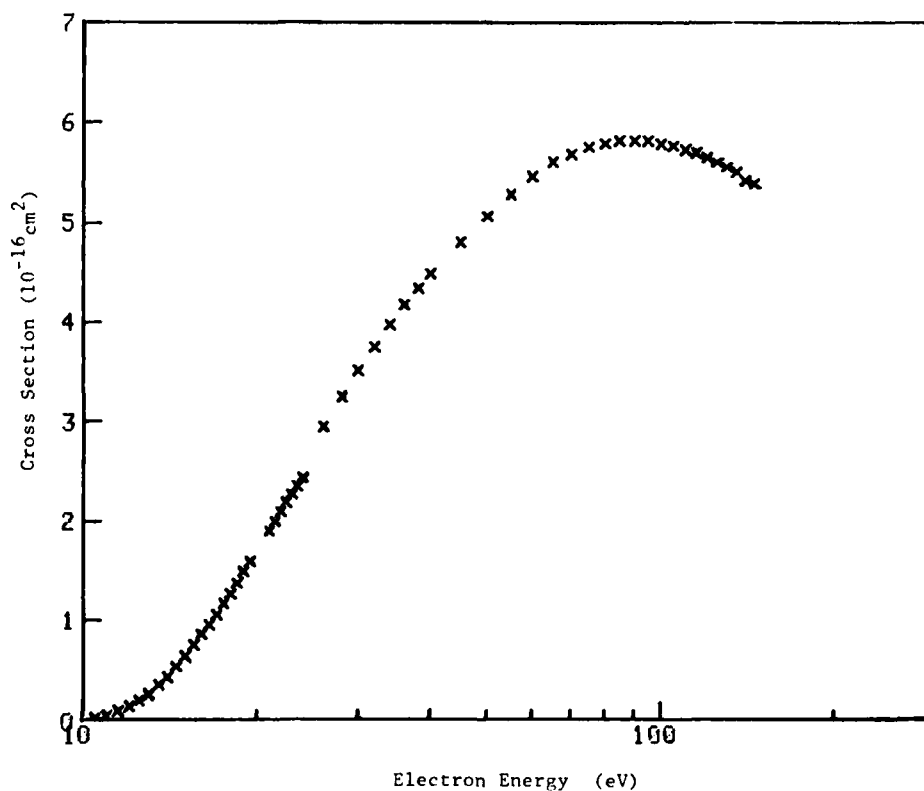


Tabular and Graphical Data C-4.18. Cross sections for electron-impact ionization of C_2H_4 .

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
10.5	0.0114	21.5	2.00	75.0	5.76
11.0	0.0449	22.0	2.09	80.0	5.80
11.5	0.0871	22.5	2.18	85.0	5.83
12.0	0.134	23.0	2.27	90.0	5.83
12.5	0.193	23.5	2.35	95.0	5.83
13.0	0.263	24.0	2.44	100	5.79
13.5	0.345	26.0	2.95	105	5.77
14.0	0.431	28.0	3.25	110	5.74
14.5	0.533	30.0	3.52	115	5.70
15.0	0.642	32.0	3.76	120	5.66
15.5	0.752	34.0	3.98	125	5.61
16.0	0.861	36.0	4.18	130	5.56
16.5	0.959	38.0	4.35	135	5.51
17.0	1.06	40.0	4.50	140	5.42
17.5	1.16	45.0	4.80	145	5.39
18.0	1.27	50.0	5.07		
18.5	1.37	55.0	5.29		
19.0	1.49	60.0	5.47		
19.5	1.59	65.0	5.61		
21.0	1.90	70.0	5.69		

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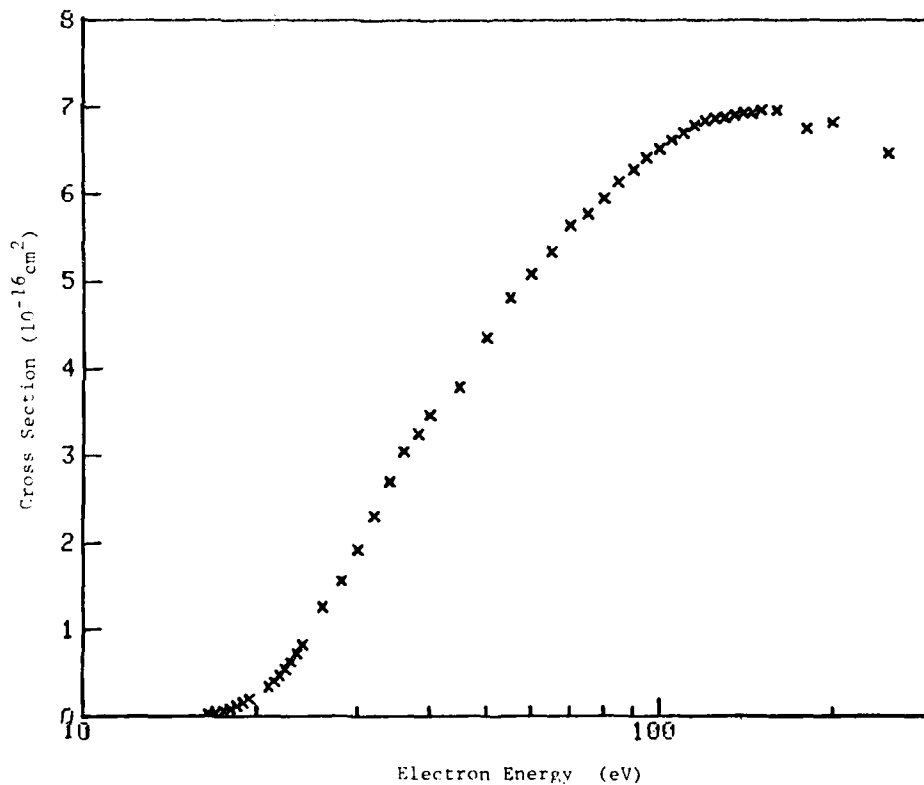
Reference: D. Rapp and P. Englander-Golden, J. Chem. Phys. 43, 1464 (1965).

Tabular and Graphical Data C-4.19. Cross sections for electron-impact ionization of SF_6 .

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
16.5	0.0202	34.0	2.70	115	6.79
17.0	0.0352	36.0	3.04	120	6.86
17.5	0.0554	38.0	3.26	125	6.87
18.0	0.0836	40.0	3.47	130	6.88
18.5	0.110	45.0	3.79	135	6.92
19.0	0.155	50.0	4.35	140	6.93
19.5	0.194	55.0	4.81	145	6.94
21.0	0.331	60.0	5.09	150	6.97
21.5	0.392	65.0	5.34	160	6.97
22.0	0.456	70.0	5.65	180	6.76
22.5	0.537	75.0	5.77	200	6.63
23.0	0.621	80.0	5.95	250	6.48
23.5	0.715	85.0	6.14	300	6.09
24.0	0.820	90.0	6.28		
26.0	1.26	95.0	6.42		
28.0	1.56	100	6.53		
30.0	1.93	105	6.63		
32.0	2.31	110	6.71		

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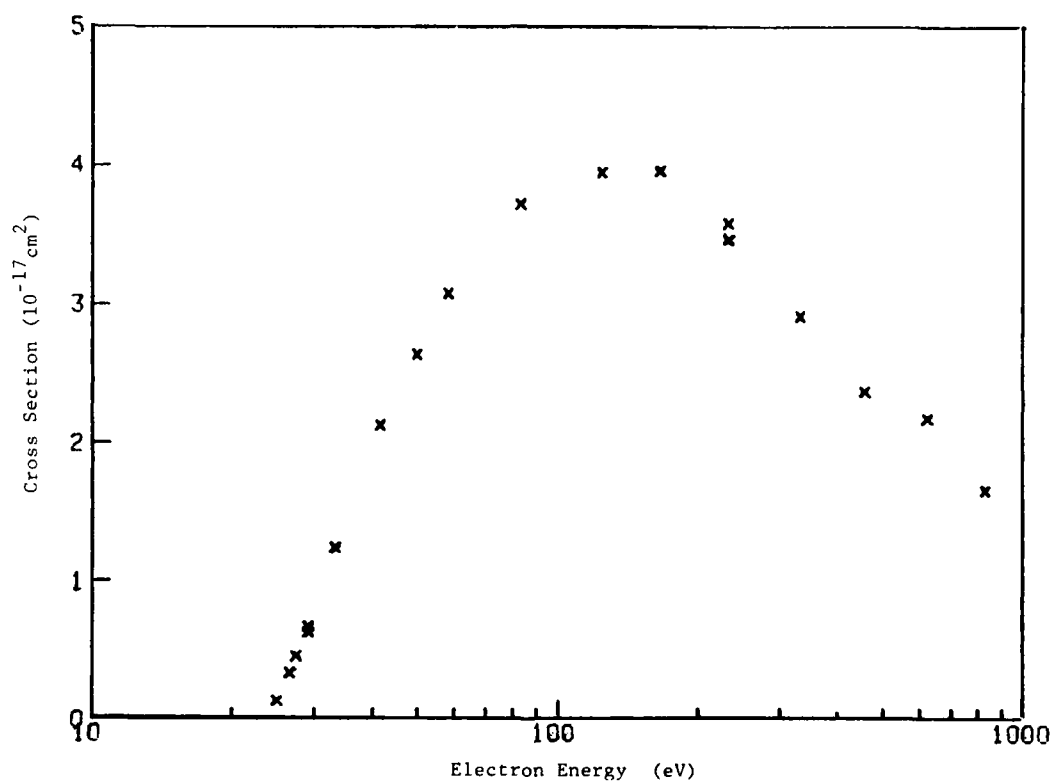


Reference: D. Rapp and P. Englander-Golden, J. Chem. Phys. 43, 1464 (1965).

Tabular and Graphical Data C-4.20. Cross sections for electron-impact
ionization of CO_2^+ ions.

Electron Energy eV	Cross Section 10^{-17} cm^2	Electron Energy eV	Cross Section 10^{-17} cm^2
24.90	0.126	124.5	3.95
26.50	0.326	166.0	3.96
27.40	0.450	232.4	3.47
29.00	0.626	232.4	3.58
29.00	0.662	332.0	2.91
33.20	1.24	456.5	2.37
41.50	2.13	622.5	2.18
49.80	2.64	830.0	1.66
58.10	3.08		
83.00	3.72		

Cont. Next Column



Reference: A. Muller, E. Salzborn, R. Frodl, R. Becker, and H. Klein, J. Phys. B 13,
L221 (1980).

Tabular Data C-4.21. Appearance potentials for electron-impact
ionization of rare-gas dimers.

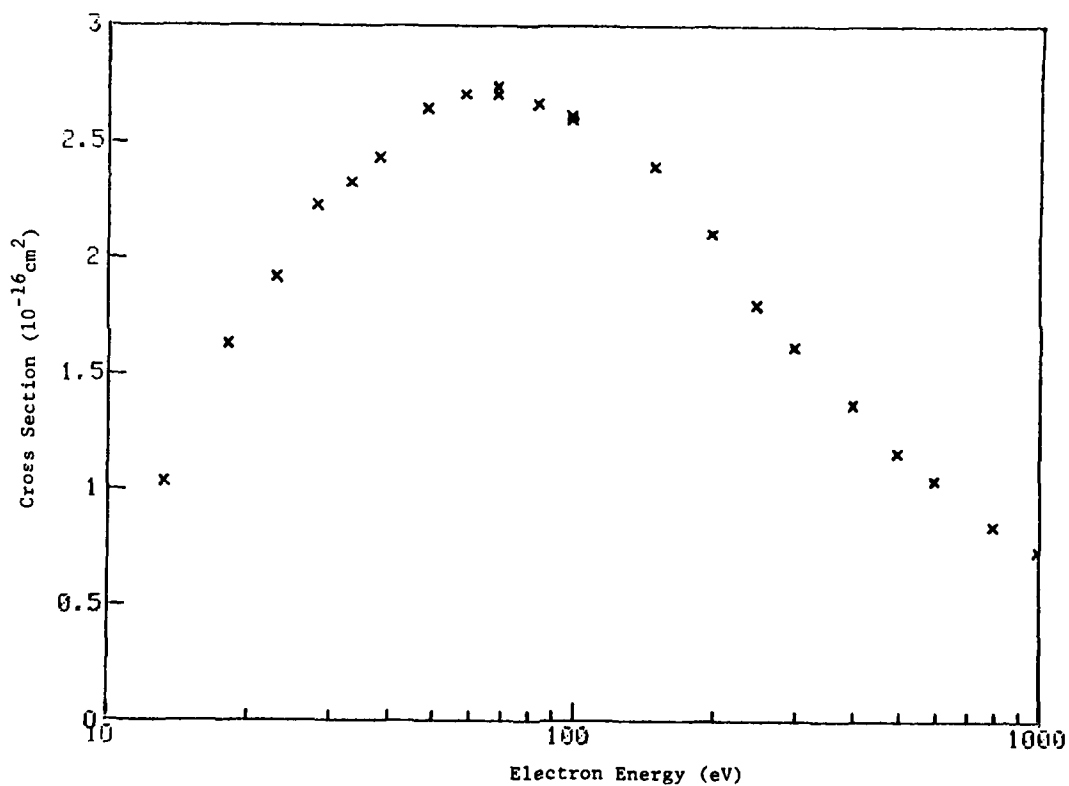
Dimer	Appearance Potential (eV)
Ar ₂	15.2 ± 0.2
ArKr	14.0 ± 0.2
Kr ₂	13.45 ± 0.3
KrXe	12.2 ± 0.2
Xe ₂	11.75 ± 0.3

Reference: H. Helm, K. Stephan and T. D. Mark, Phys. Rev. A 19,
2154 (1979).

Tabular and Graphical Data C-4.22. Cross sections for electron-impact detachment from F^- .

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
13.2	1.03	98.0	2.60
18.0	1.63	148	2.40
23.0	1.92	197	2.11
28.0	2.23	247	1.80
33.0	2.33	296	1.62
38.0	2.44	396	1.37
48.0	2.65	495	1.16
58.0	2.71	595	1.04
68.0	2.74	794	0.840
83.0	2.71	990	0.730
98.0	2.67		
98.0	2.62		

Cont. Next Column



Reference: B. Peart, R. Forrest, and K. T. Dolder, J. Phys. B 12, L115 (1979).

Tabular and Graphical Data C-4.23. Cross sections for electron-impact detachment from O^- .

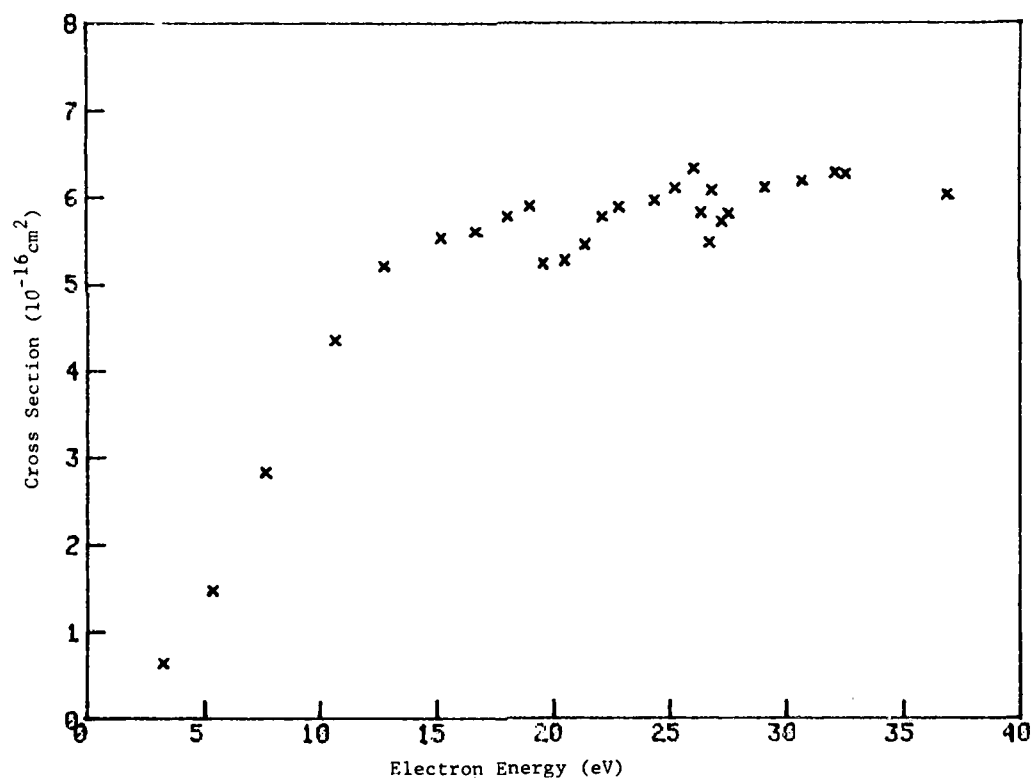
Inclined beams

90% confidence limits of random error

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
3.28	0.640	20.5	5.27	27.3	5.72
5.37	1.48	21.4	5.46	27.5	5.81
7.62	2.83	22.1	5.78	29.1	6.12
10.6	4.35	22.8	5.89	30.7	6.19
12.8	5.21	24.4	5.97	32.1	6.29
15.2	5.54	25.2	6.10	32.5	6.28
16.7	5.61	26.0	6.34	36.9	6.03
18.1	5.78	26.3	5.82		
19.0	5.90	26.8	6.08		
19.6	5.24	26.7	5.48		

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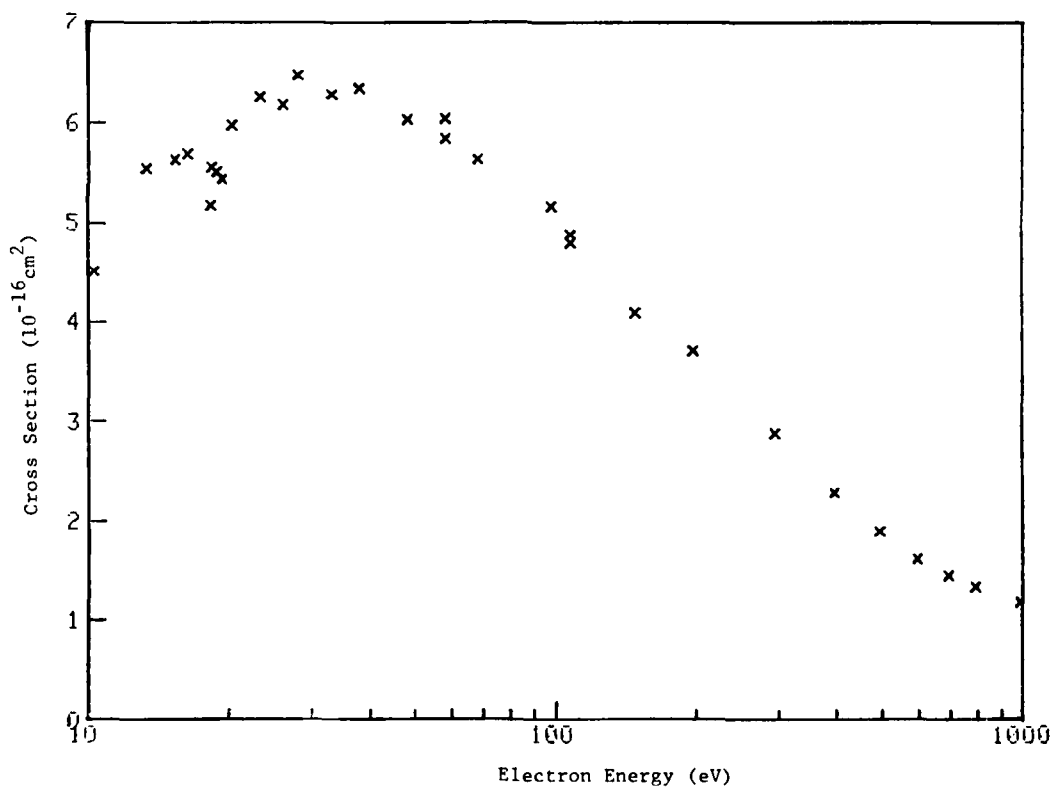


Reference: B. Peart, R. A. Forrest, and K. Dolder, J. Phys. B 12, 2735 (1979).

Tabular and Graphical Data C-4.24 Cross sections for electron-
impact detachment from O^- .

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16}cm^2	eV	10^{-16}cm^2
10.3	4.51	58.0	5.85
13.3	5.54	68.0	5.64
15.3	5.63	98.0	5.16
16.3	5.69	108	4.88
18.3	5.18	108	4.80
18.3	5.56	148	4.09
18.8	5.51	197	3.71
19.3	5.44	296	2.88
20.3	5.98	396	2.28
23.3	6.26	495	1.89
26.1	6.18	595	1.62
28.1	6.48	694	1.45
33.3	6.29	794	1.33
38.0	6.35	990	1.18
48.0	6.04		
58.0	6.05		

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Reference: B. Peart, R. Forrest, and K. T. Dolder, J. Phys. B 12, 847 (1979).

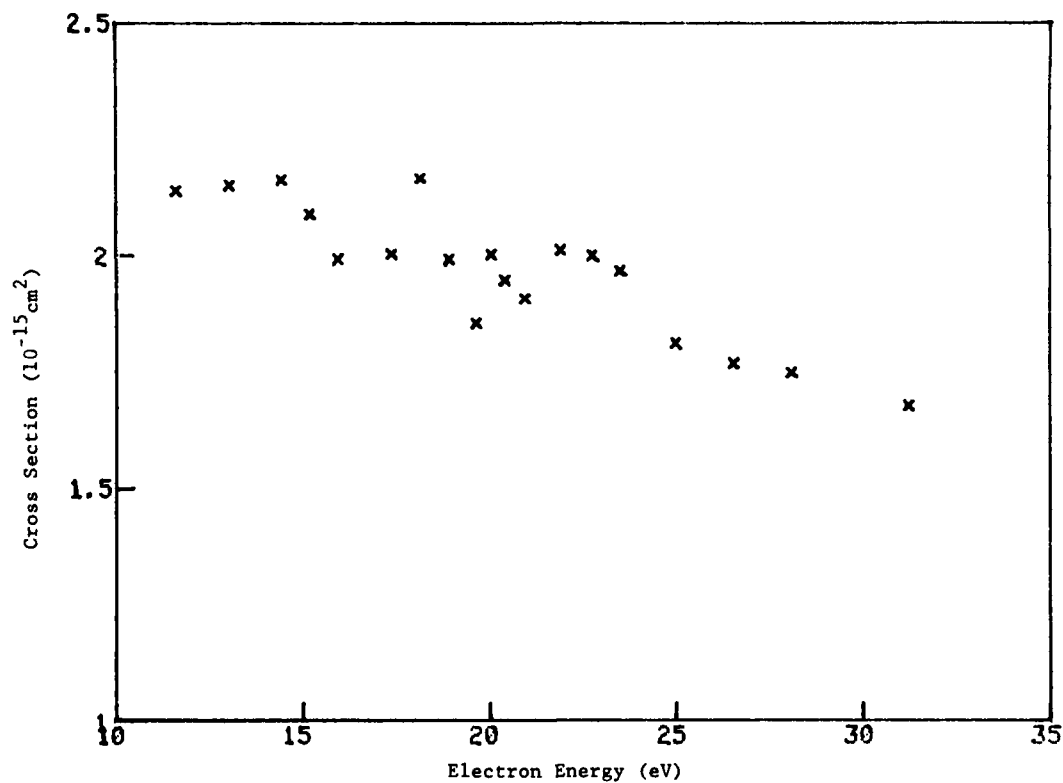
Tabular and Graphical Data C-4.25. Cross sections for electron-
impact detachment from C^- .

Inclined beams

90% confidence limits of random error

Electron Energy eV	Cross Section 10^{-15}cm^2	Electron Energy eV	Cross Section 10^{-15}cm^2
11.6	2.14	20.4	1.95
13.0	2.15	20.9	1.91
14.4	2.16	21.9	2.01
15.2	2.09	22.7	2.00
15.9	1.99	23.5	1.97
17.4	2.01	25.0	1.81
18.1	2.17	26.5	1.77
18.9	1.99	28.1	1.75
19.6	1.86	31.2	1.68
20.0	2.00		

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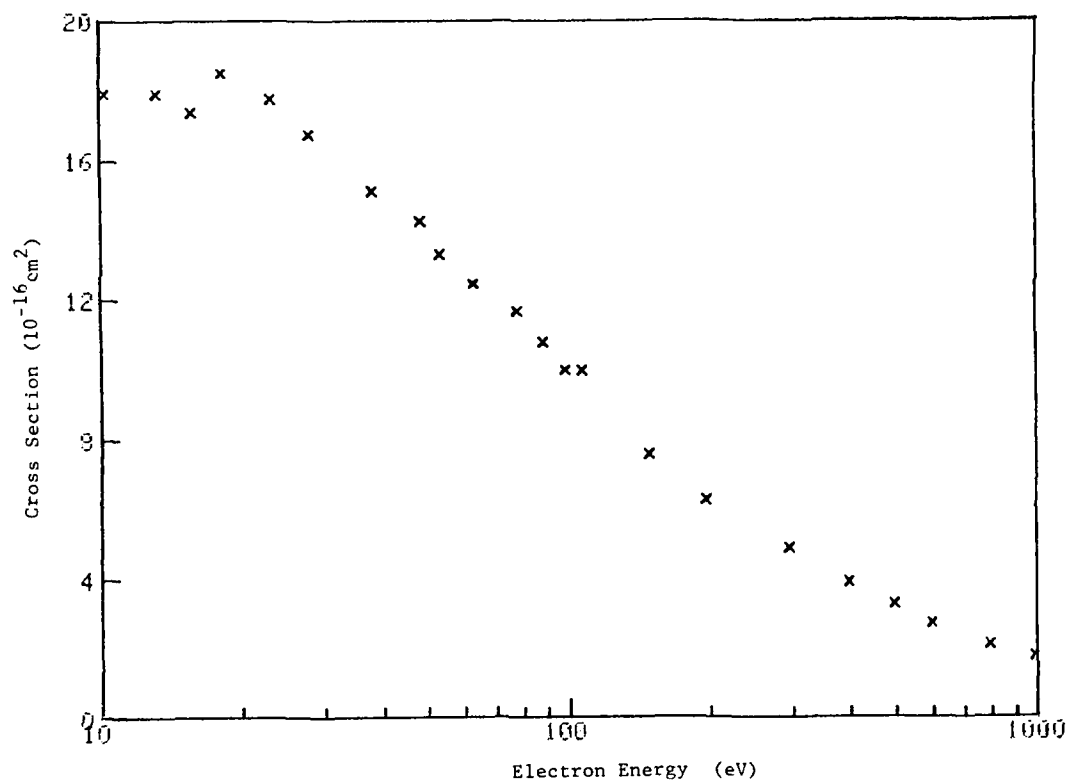


Reference: B. Peart, R. A. Forrest, and K. Dolder, J. Phys. B 12, 2735 (1979).

Tabular and Graphical Data C-4.26. Cross sections for electron-
impact detachment from C^- :

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
10.2	17.9	98.0	9.97
13.2	17.9	107	9.96
15.7	17.4	148	7.55
18.2	16.5	197	6.23
23.2	17.8	296	4.87
28.0	16.7	396	3.87
38.0	15.1	495	3.22
48.0	14.2	595	2.70
53.0	13.3	794	2.07
63.0	12.4	990	1.73
78.0	11.6		
88.0	10.8		

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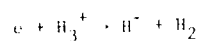
Reference: B. Peart, R. Forrest, and K. T. Dolder, J. Phys. B 12, 847 (1979).

C-5. ELECTRON-ION RECOMBINATION

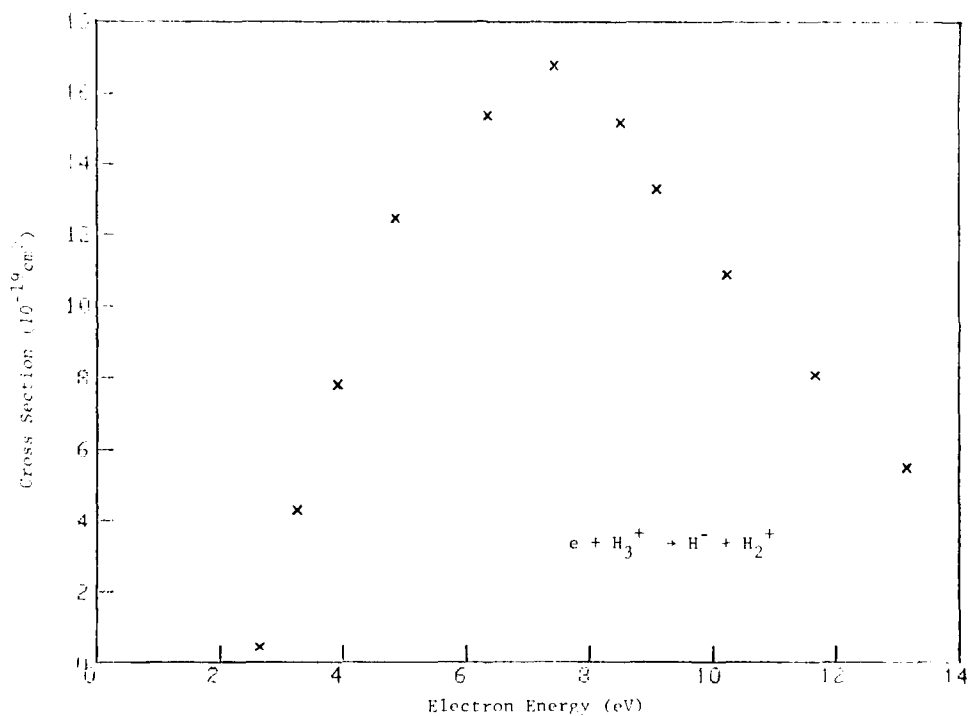
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Tabular and Graphical Data C-5.1. Cross sections for
dissociative recombination of electrons with H_3^+ to form H^- .



Electron Energy	Cross Section
eV	10^{-19} cm^2
2.64	0.421
3.75	4.30
3.91	7.80
4.85	12.5
6.34	15.4
7.40	16.8
8.51	15.2
9.09	13.3
10.2	10.9
11.7	8.08
13.2	5.51



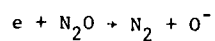
Reference: B. Peart, R. A. Forrest, and K. Dolder, J. Phys. B 12, 3441 (1979)

C-6. NEGATIVE ION FORMATION BY ELECTRON IMPACT

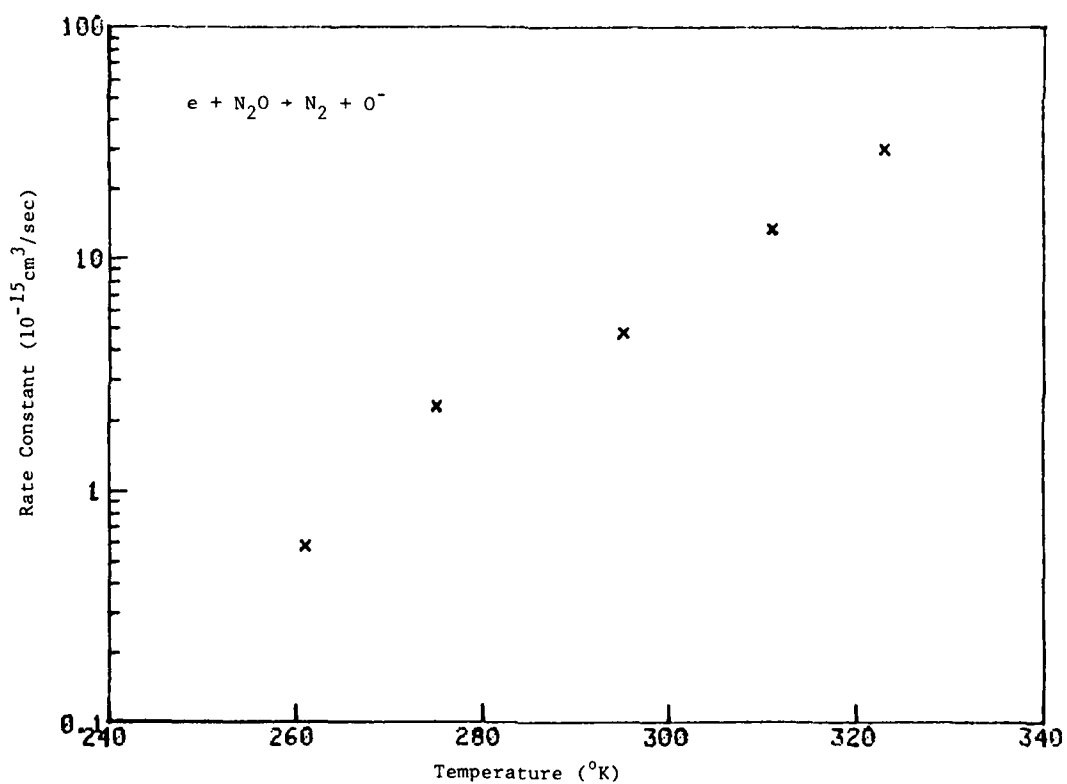
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C-6.1.	Cross sections for dissociative attachment of electrons to N_2O	2999
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C-6.4.	Total electron attachment cross sections for $CClF_3$	3002
C-6.5.	The effect of vibrational and rotational excitation on threshold dissociative attachment cross sections in H_2 and D_2	3003
C-6.6.	Increase in dissociative attachment cross sections in SF_6 due to vibrational excitation of the gas prior to the collision.....	3005

Tabular and Graphical Data C-6.1. Cross sections for dissociative
attachment of electrons to N_2O .



Temperature	Rate Constant
K	$10^{-15} \text{ cm}^3/\text{mole-sec}$
323	30.0
311	13.5
295	4.80
275	2.30
261	0.580

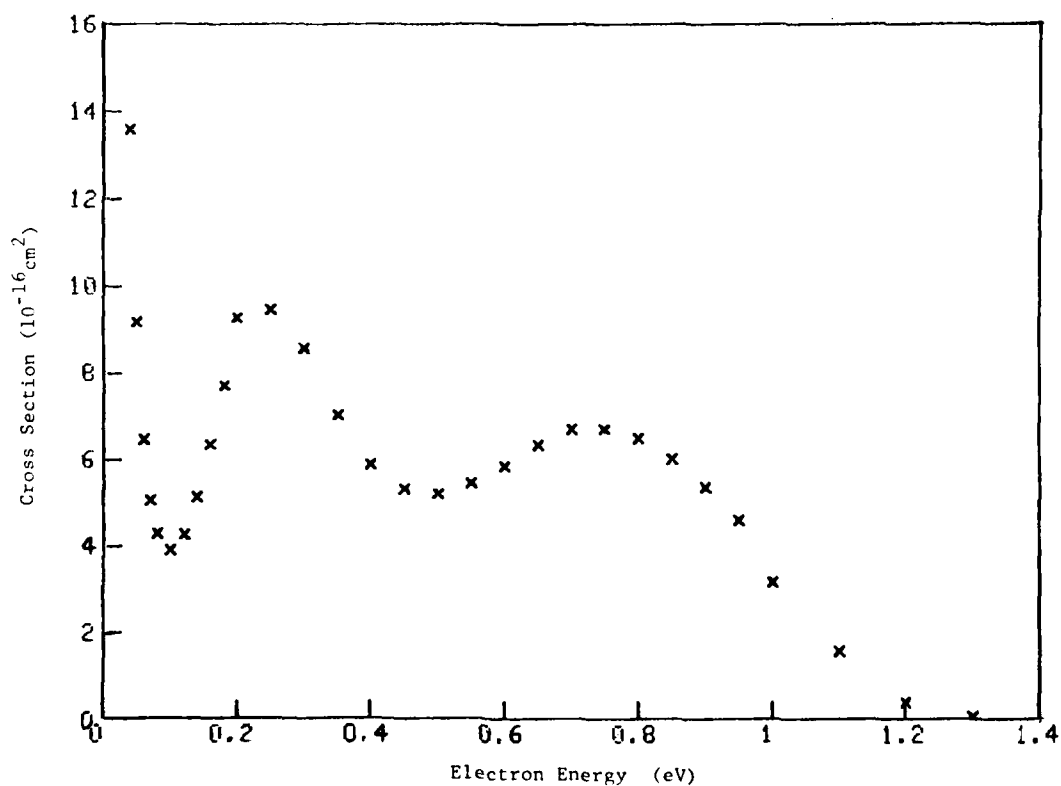


Reference: H. Shimamori and R. W. Fessenden, J. Chem. Phys. 70, 1137 (1979).

Tabular and Graphical Data C-6.2. Total electron attachment cross
sections for CCl_3F .

Electron Energy	Cross Section	Electron Energy	Cross Section
eV	10^{-16} cm^2	eV	10^{-16} cm^2
0.040	13.6	0.50	5.19
0.050	9.10	0.55	5.44
0.060	6.46	0.60	5.83
0.070	5.04	0.65	6.32
0.080	4.27	0.70	6.67
0.10	3.59	0.75	6.70
0.12	4.24	0.80	6.47
0.14	5.12	0.85	5.99
0.16	6.34	0.90	5.36
0.18	7.69	0.95	4.59
0.20	9.26	1.00	3.15
0.25	9.46	1.1	1.56
0.30	8.55	1.2	0.390
0.35	7.02	1.3	0.0500
0.40	5.67		
0.45	5.36		

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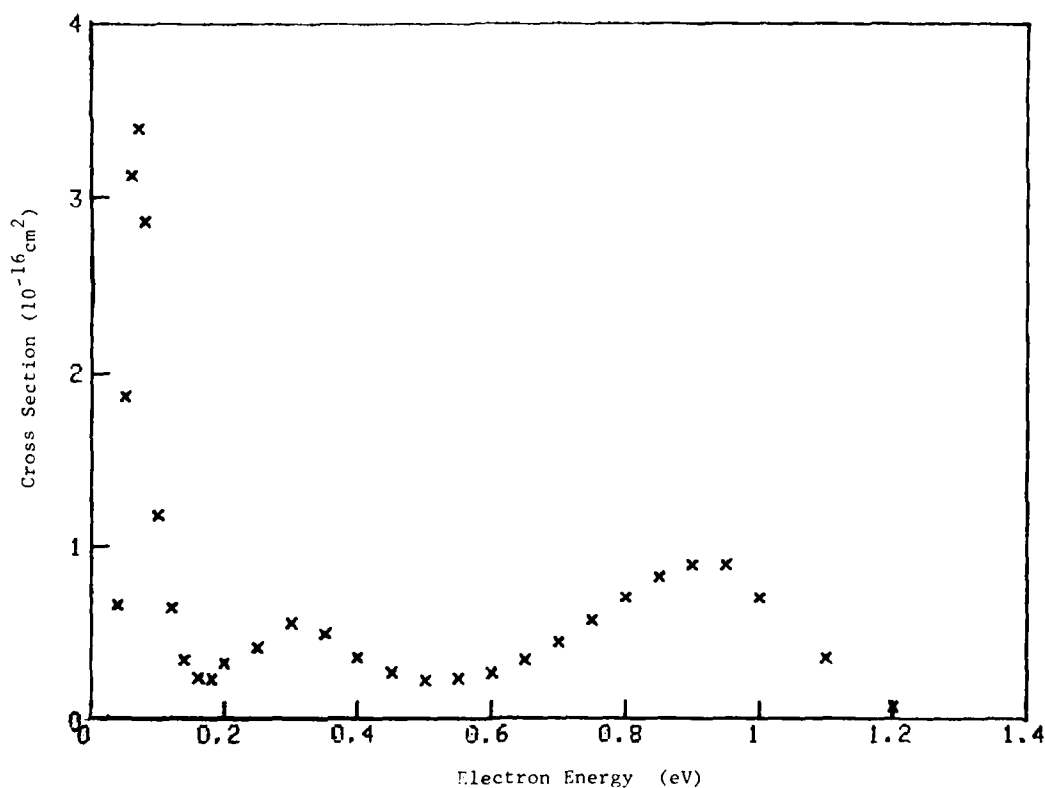


Reference: D. L. McCorkle, A. A. Christodoulides, L. G. Christophorou, and I. Szamrej,
J. Chem. Phys. 72, 4049 (1980).

Tabular and Graphical Data C-6.3. Total electron attachment cross
sections for CCl_2F_2 .

Electron Energy eV	Cross Section 10^{-16} cm^2	Electron Energy eV	Cross Section 10^{-16} cm^2
0.040	0.600	0.50	0.220
0.050	1.80	0.55	0.230
0.060	3.13	0.60	0.260
0.070	3.40	0.65	0.340
0.080	2.67	0.70	0.440
0.10	1.18	0.75	0.570
0.12	0.640	0.80	0.700
0.14	0.340	0.85	0.820
0.16	0.240	0.90	0.890
0.18	0.230	0.95	0.890
0.20	0.320	1.00	0.700
0.25	0.410	1.1	0.350
0.30	0.550	1.2	0.0700
0.35	0.490		
0.40	0.350		
0.45	0.260		

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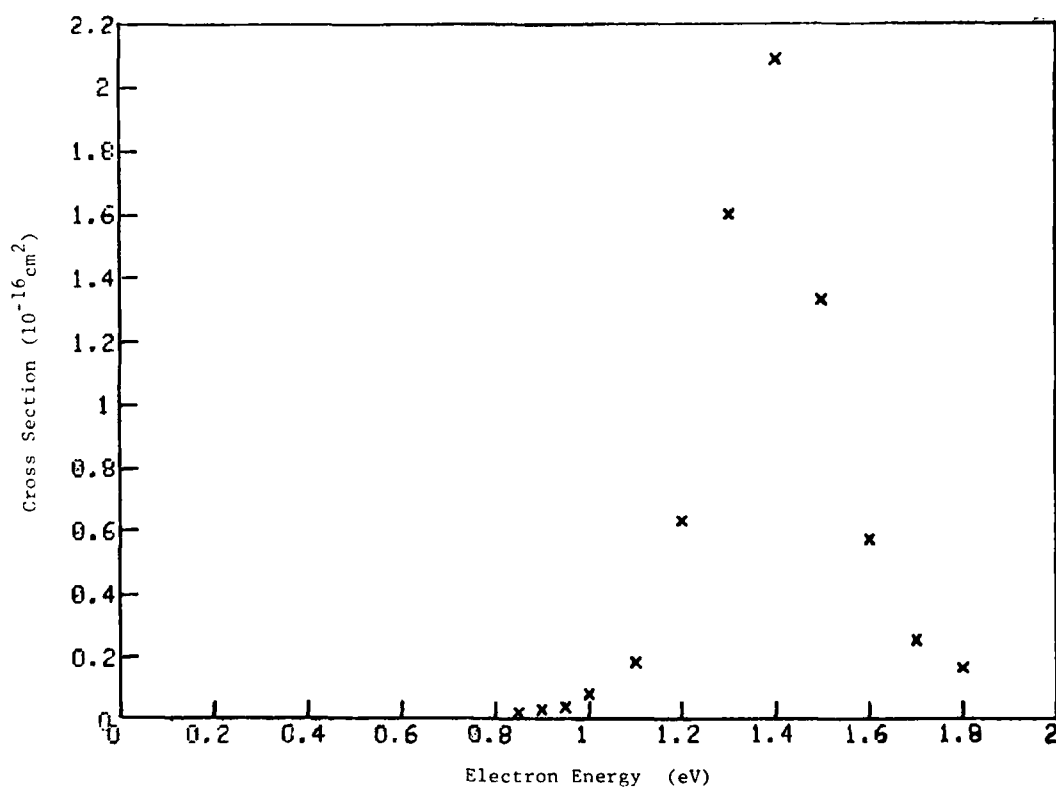


Reference: D. L. McCorkle, A. A. Christodoulides, L. G. Christophorou, and I. Szamrej,
J. Chem. Phys. 72, 4049 (1980).

Tabular and Graphical Data C-6.4. Total electron attachment cross
sections for CClF_3 .

Electron Energy eV	Cross Section 10^{-16}cm^2	Electron Energy eV	Cross Section 10^{-16}cm^2
0.85	0.0200	1.4	2.09
0.90	0.0300	1.5	1.33
0.95	0.0400	1.6	0.570
1.00	0.0800	1.7	0.250
1.1	0.180	1.8	0.160
1.2	0.630		
1.3	1.60		

Cont. Next Column



Reference: D. L. McCorkle, A. A. Christodoulides, L. G. Christophorou, and I. Szamrej,
J. Chem. Phys. 72, 4049 (1980).

Tabular Data C-6.5. The effect of vibrational and rotational excitation on threshold dissociative attachment cross sections in H_2 and D_2 .

Tabular Data C-6.5a. Experimental ratio of threshold dissociative attachment cross sections for vibrationally excited molecules to that for molecules in the vibrational ground state.

v	H_2		D_2	
	Internal Energy eV	$\sigma_{DA}^v / \sigma_{DA}^{v=0}$	Internal Energy eV	$\sigma_{DA}^v / \sigma_{DA}^{v=0}$
1	.49	32.	.36	42.
2	.98	560	.70	900.
3	1.44	5.6×10^3	1.05	1.2×10^4
4	1.86	3.8×10^4	1.36	1.0×10^5
5			1.67	7.2×10^5

Tabular Data C-6.5b. Experimental ratio of threshold dissociative attachment cross sections for rotationally excited molecules to that for molecules in the rotational ground state.

j	H_2	
	Internal Energy eV	$\sigma_{DA}^j / \sigma_{DA}^{j=0}$
5	.23	1.6
7	.41	4.2

Ground-state dissociative attachment cross sections at 300°K, taken from G. Schulz and R. K. Asundi, Phys. Rev. 158, 25 (1967)

$$\sigma_{DA}(H_2) = 1.6 \times 10^{-21} \text{ cm}^2$$

$$\sigma_{DA}(D_2) = 8 \times 10^{-24} \text{ cm}^2$$

Reference: M. Allan and S. F. Wong, Phys. Rev. Lett 41, 1791 (1978).

Table 1. Calculated ratio of threshold dissociative attachment cross sections for vibrationally excited molecules to that for molecules in the vibrational ground state.

H_2	Internal Energy (eV)	$\sigma_{vib}/\sigma_{g.s.}$	H_2	Internal Energy (eV)	$\sigma_{vib}/\sigma_{g.s.}$
	0.91	0.8		0.91	0.8
	1.1	0.85		1.1	0.85
	1.46	1.8×10^{-1}		1.46	1.8×10^{-1}
	1.89	2.1×10^{-1}		1.89	2.1×10^{-1}
				1.68	1.6×10^{-1}

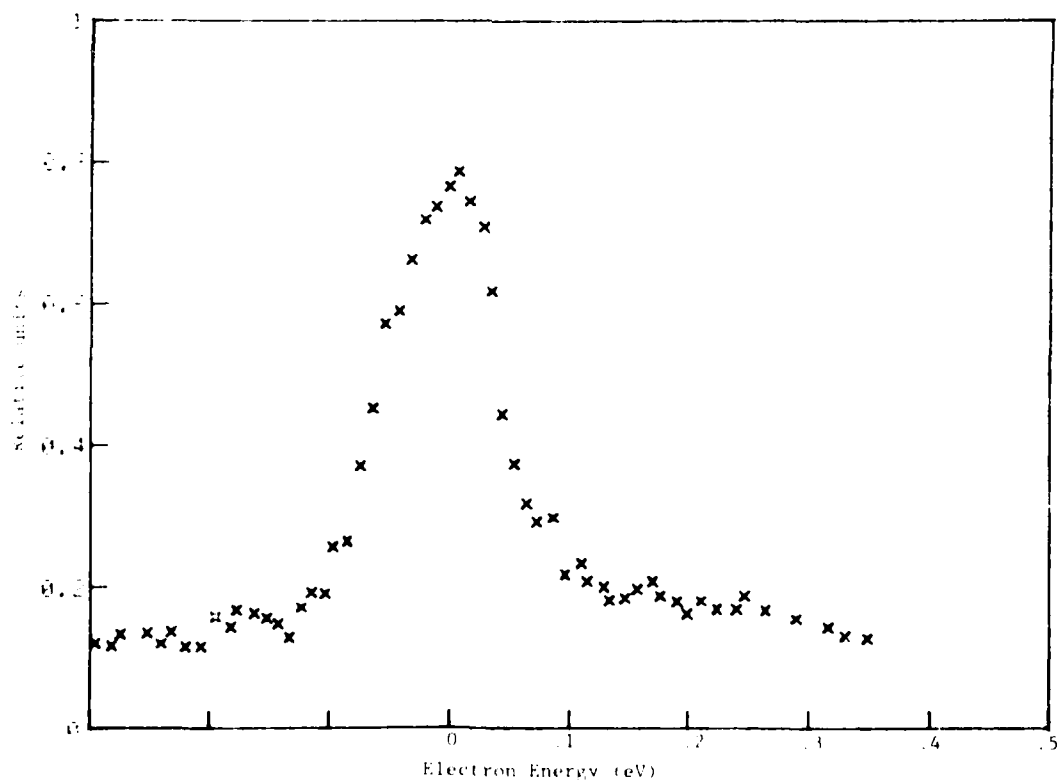
Reference: 1. N. Bardsley and J. M. Wadehra, Phys. Rev. A 20, 1199 (1979).

Tabular and Graphical Data C-6.6. Increase in dissociative attachment cross sections in SF_6 due to vibrational excitation of the gas prior to the collision.

- a. Temperature dependence of the cross section for dissociative attachment to SF_6 to form SF_5^- . Taken from Fig. 1 of the reference.

Temperature (°K)	Electron Energy (eV)	0	0.1	0.2	0.3
300		6.5	3.6	5.4	6.4
330		8.3	5.0	6.4	7.5
355		14.	7.0	8.6	9.3
420		44.	16.	15.	-
500		130.	29.	19.	-
607		290.	43.	23.	-
740		490.	76.	30.	-
880		780.	-	-	-

- b. Laser enhancement of dissociative attachment cross section for electrons on $^{32}\text{SF}_6$ to produce $^{32}\text{SF}_5^-$. The radiation selectively excites the ν_3 vibrational mode of $^{32}\text{SF}_6$ to produce the enhanced signal.



Reference: G. L. Chen and P. J. Chantry, J. Chem. Phys. 71, 3897 (1979).

D. PHOTON COLLISION PROCESSES IN GASES

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The data presented in this chapter either extend or supersede the data given previously in Chapter D of Volume II, pages 639-713, and Volume IV, pages 1917-2078.

ACKNOWLEDGEMENTS

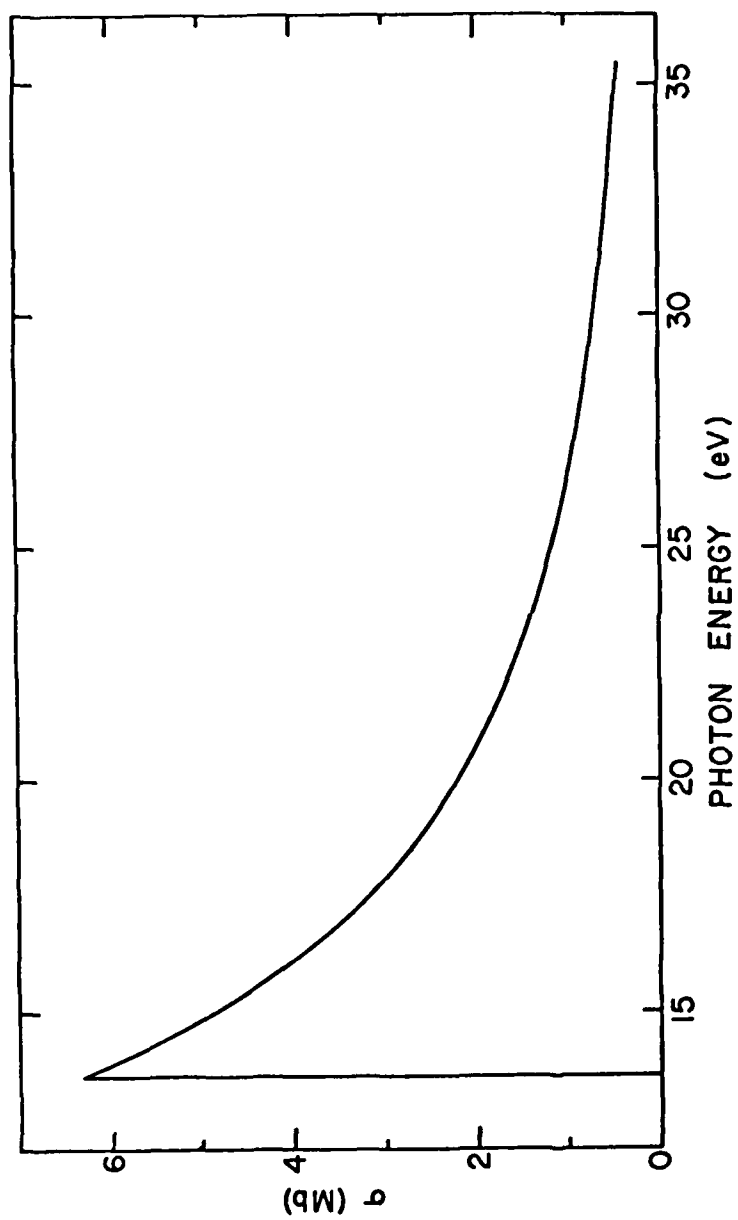
This chapter on photon collision processes in gases was put together with the aid of a number of scientists. Particularly significant were the contributions of Dr. Joseph Berkowitz, of Argonne National Laboratory, whose book Photoabsorption, Photoionization, and Photoelectron Spectroscopy (Academic Press, New York, 1979) provided us with a wealth of references and critically evaluated data on atoms and molecules. We gratefully acknowledge being allowed access to the manuscript prior to publication as well as Dr. Berkowitz providing us with a number of large-size versions of figures from his book.

In addition we acknowledge the contributions of Professor C.E. Brion, of the University of British Columbia, for providing us with a complete set of reprints, spanning a decade, of his very extensive work on partial and total cross sections of atoms and molecules.

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Graphical Data D-1.1
Theoretical Photoionization Cross Section of Atomic H.

Tabular Data D-1.2

Branching Ratios and Partial Cross Section for
the Photoionization of Ar 3s

Energy loss (eV)	Measured intensity ratio ^b 3s/3p	Branching ratio ^c		Total photoionization cross-section (Mb) ^d	Argon 3s photoionization ^e cross-section σ (Mb)
		3s	3p		
32.8	0.031	0.030 (6)	0.970 (6)	16.7	0.500 (100)
33.8	0.023	0.022 (9)	0.978 (10)	14.2	0.310 (127)
35.8	0.023	0.022 (8)	0.978 (10)	9.4	0.210 (75)
37.8	0.012	0.016 (10)	0.983 (13)	5.9	0.094 (59)
39.8	0.000	0.000 (4)	1.000 (4)	3.7	0.000 (14)
41.8	0.012	0.011 (11)	0.989 (11)	2.2	0.024 (24)
44.8	0.061	0.057 (43)	0.943 (44)	1.25	0.070 (53)
46.8	0.103	0.093 (35)	0.907 (46)	0.97	0.091 (34)
49.8	0.124	0.102 (40)	0.826 (59)	0.92	0.094 (37)
51.8	0.129	0.106 (68)	0.782 (65)	1.01	0.110 (70)
54.8	0.164	0.123 (28)	0.754 (40)	1.19	0.146 (33)
59.8	0.226	0.155 (38)	0.685 (38)	1.37	0.213 (52)
64.8	0.267	0.177 (35)	0.658 (42)	1.45	0.257 (51)
69.8	0.266	0.177 (37)	0.664 (48)	1.48	0.262 (55)
74.8	0.223	0.153 (68)	0.687 (68)	1.48	0.227 (101)

Values in parentheses represent the uncertainties in the derived quantities.

^b This represents the 3s/3p intensity ratio corrected for analyzer transmission efficiency.

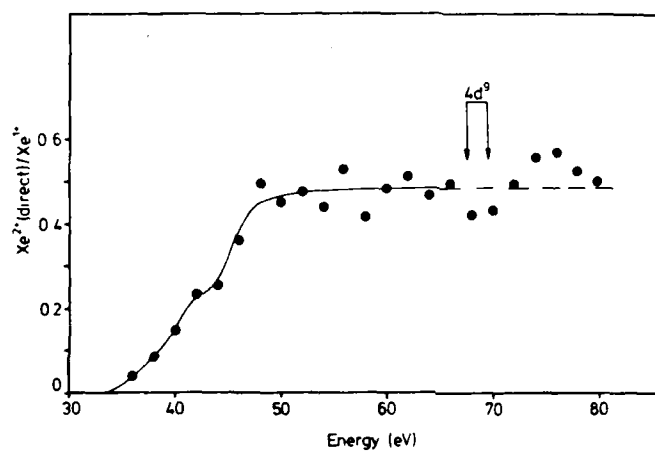
^c Branching ratio (3s) = 3s/(3s + 3p + 2+) etc. The data for multiple ionization are taken from ref. 30. This correction has been made above 46.8 eV.

^d From data of West and Marr.

^e $\sigma(\text{Mb}) = 1.0975 \times 10^2(dI/dE) (\text{eV})^{-1}$.

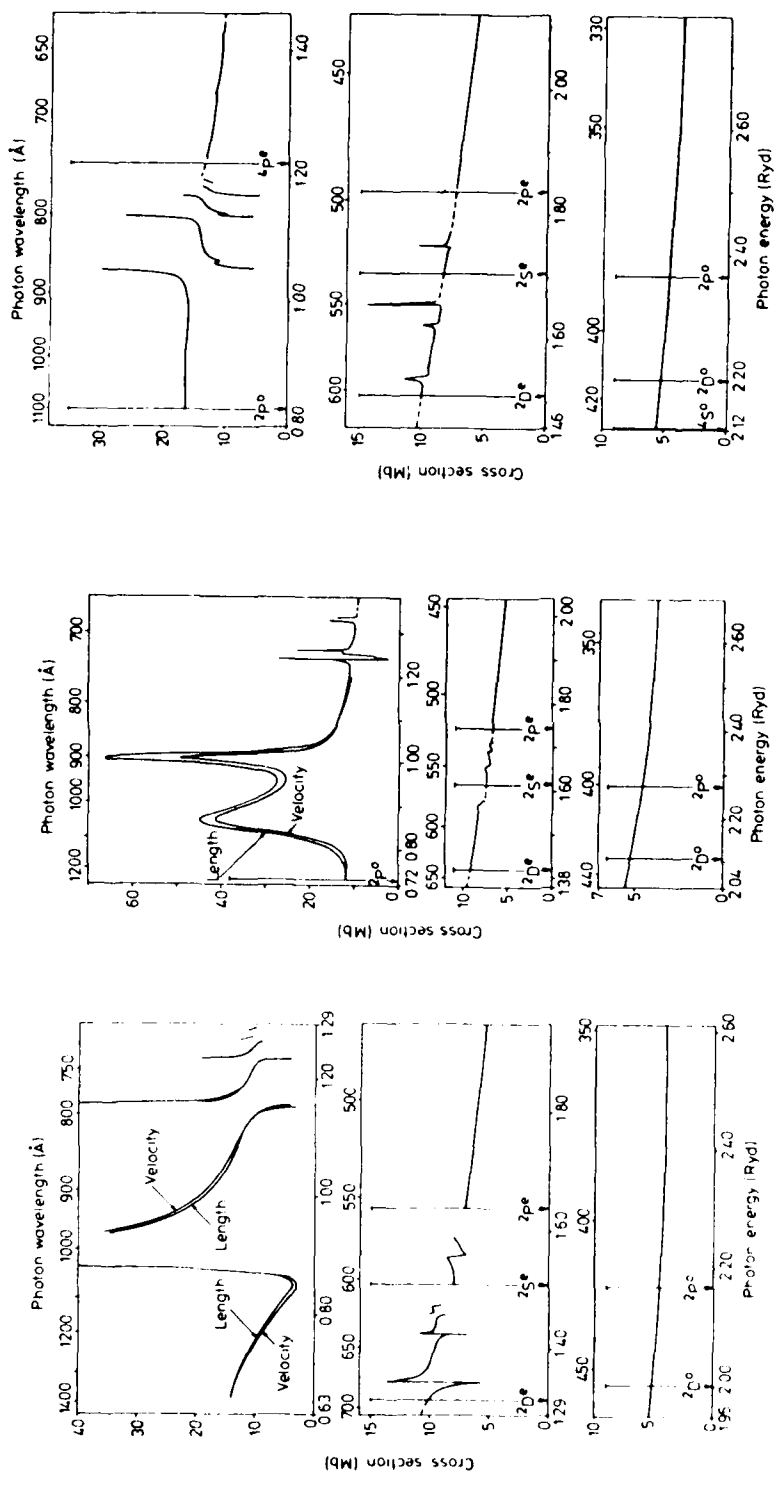
Note: The value in parentheses represent the uncertainties.

Reference: These data were taken from K. H. Tan and C. E. Brion, J. Electron Spectrosc. 13, 77 (1978) except for the total cross section which was taken from J. B. West and G. V. Marr, Proc. Roy. Soc. A 349, 397 (1976).



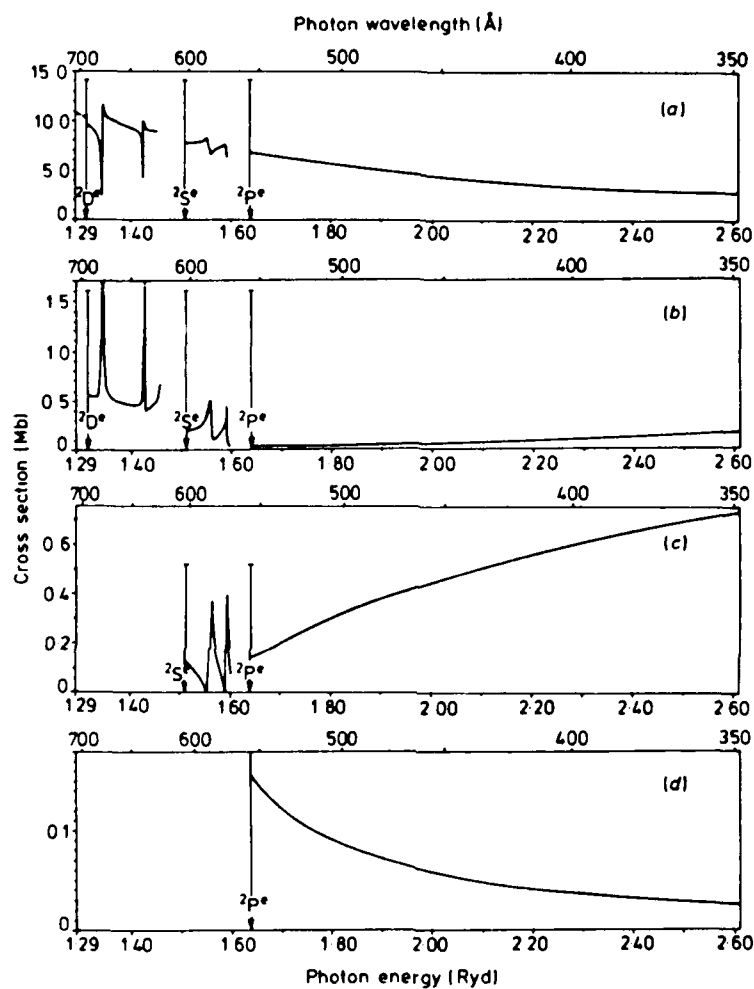
Graphical Data D-1.3

Ratio of direct double photoionization to single ionization for Xe



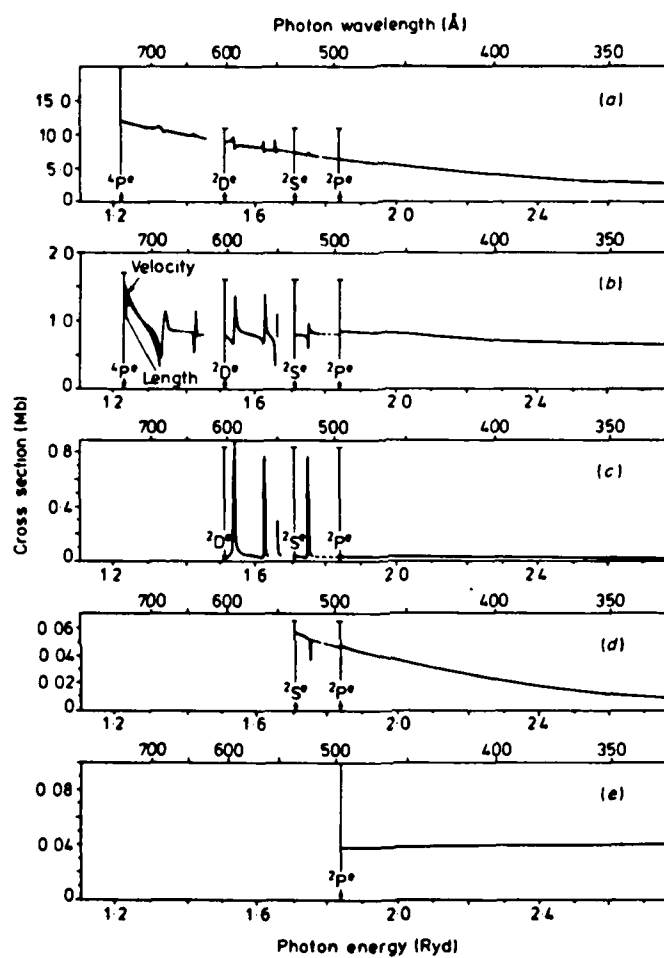
Graphical Data D-1.4

Total photoionization cross section for $C(2P)$, $C(1D)$, and $C(1S)$.



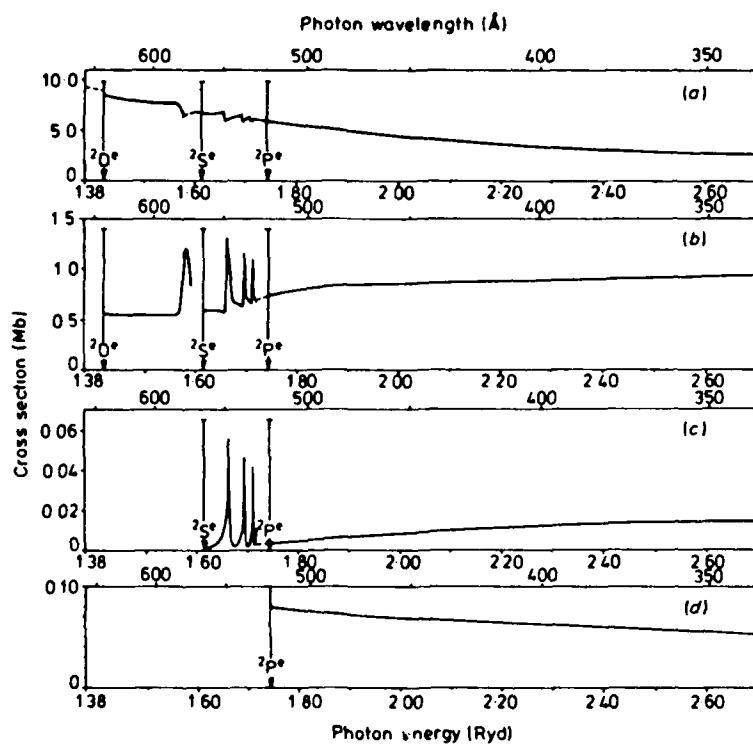
Graphical Data D-1.5

Partial photoionization cross section for $C(1S)$ leaving the ion in the a) $2s^2 2p \ ^2P$, b) $2s 2p^2 \ ^2D$, c) $2s 2p^2 \ ^2S$, and d) $2s 2p^2 \ ^2P$ states.



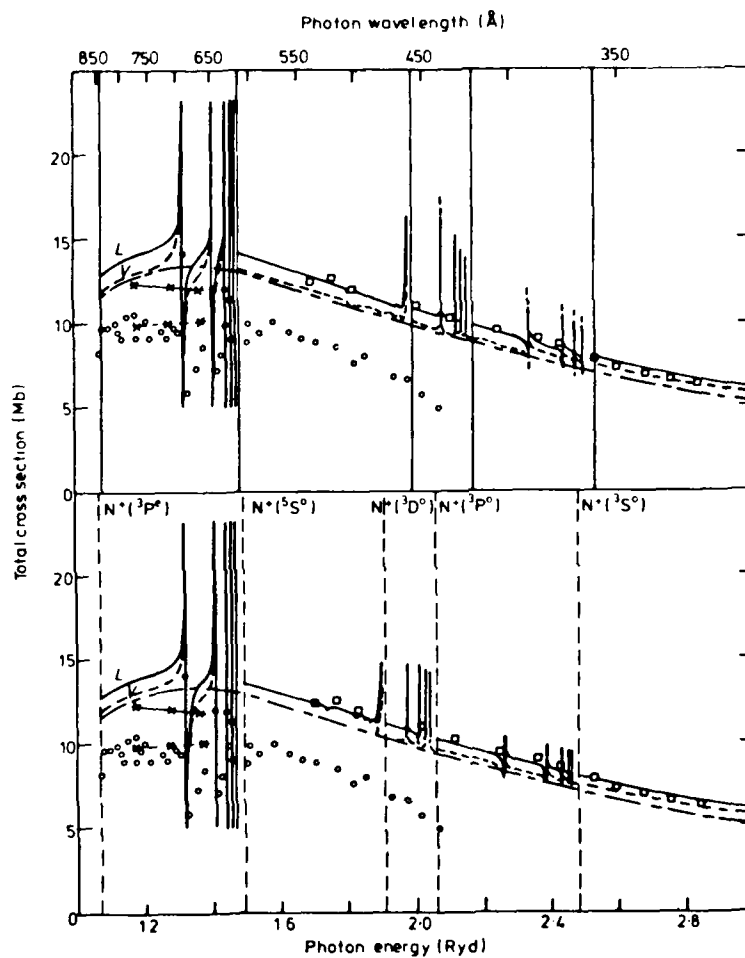
Graphical Data D-1.6

Partial photoionization cross section for $C(^3P)$ leaving the ion in the a) $2s^2 2p \ ^2P$, b) $2s \ 2p^2 \ ^4P$, c) $2s \ 2p^2 \ ^2D$, and d) $2s \ 2p^2 \ ^2S$, and e) $2s \ 2p^2 \ ^2P$ states.



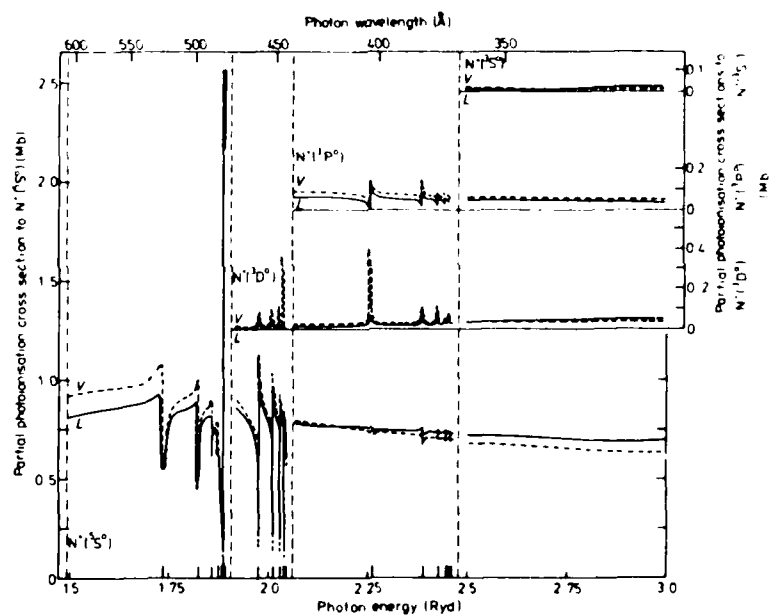
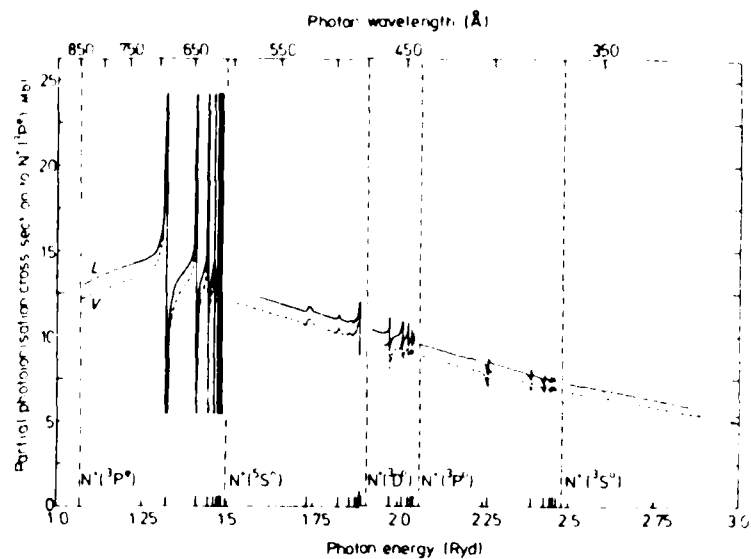
Graphical Data D-1.7

Partial photoionization cross section for $C(^1D)$ leaving the ion in the a) $2s^2 2p \ ^2P$, b) $2s 2p^2 \ ^2D$, c) $2s 2p^2 \ ^2S$, and d) $2s 2p^2 \ ^2P$ states.



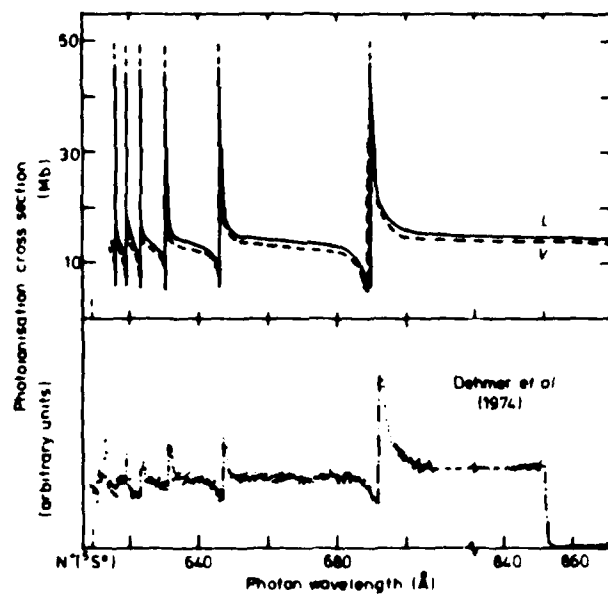
Graphical Data D-1.8

Theoretical total photoionization cross section for $N(^1S)$ in length (L) and velocity (V) formulations in two theoretical approximations compared with various other experimental and theoretical results.



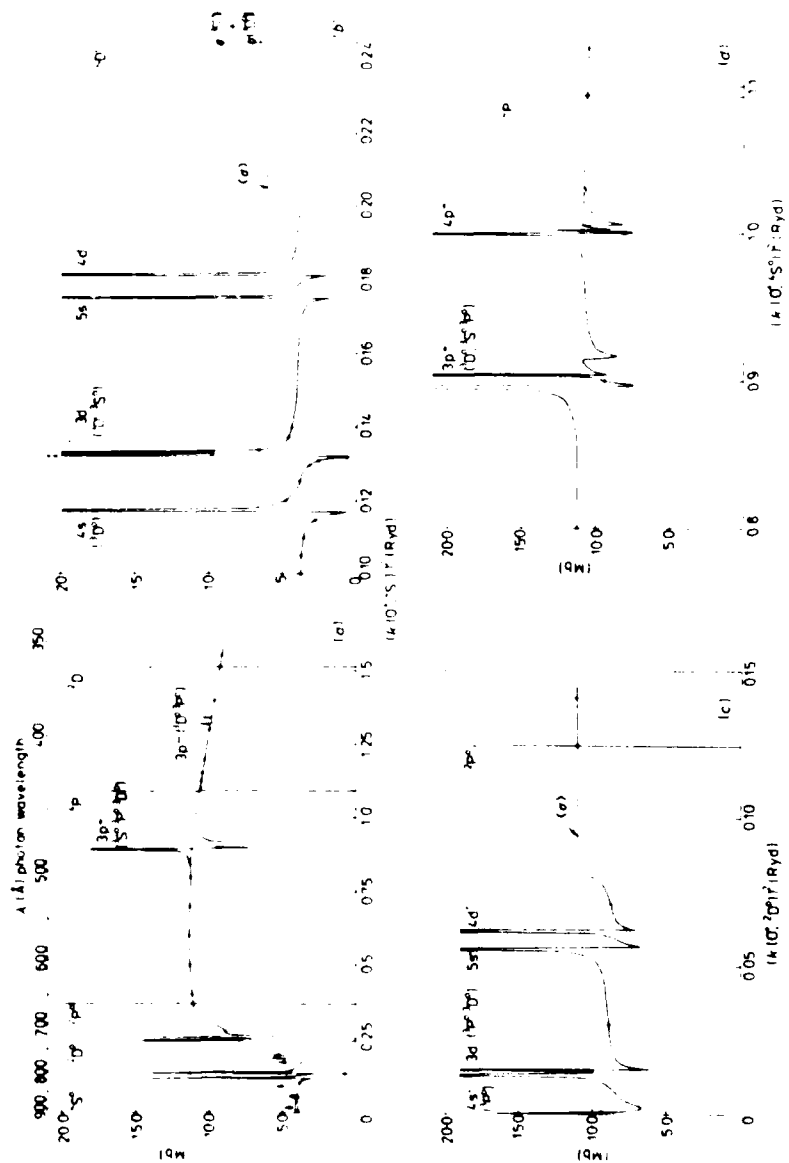
Graphical Data D-1.9

Photoionization cross section of $N(1S^\circ)$ to $N^+(1P)$ corresponding to the ejection of a 2p electron and the various N^+ states corresponding to ejection of a 2s electron in length (L) and velocity (V).



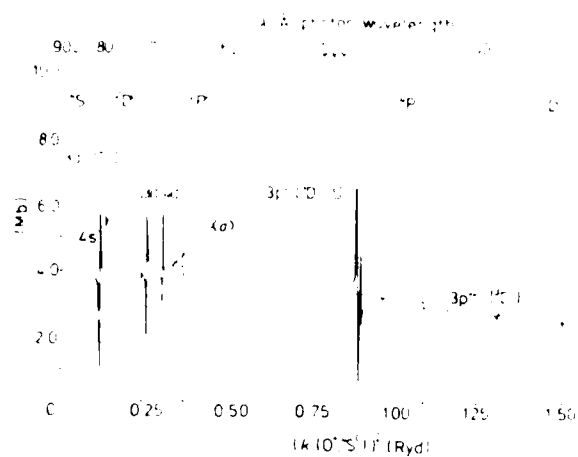
Graphical Data D-1.10

Photoionization cross section of ground state $N(1S)$ in the neighborhood of the $(2s2p^2\ ^3S)np\ ^1P$ autoionizing series.



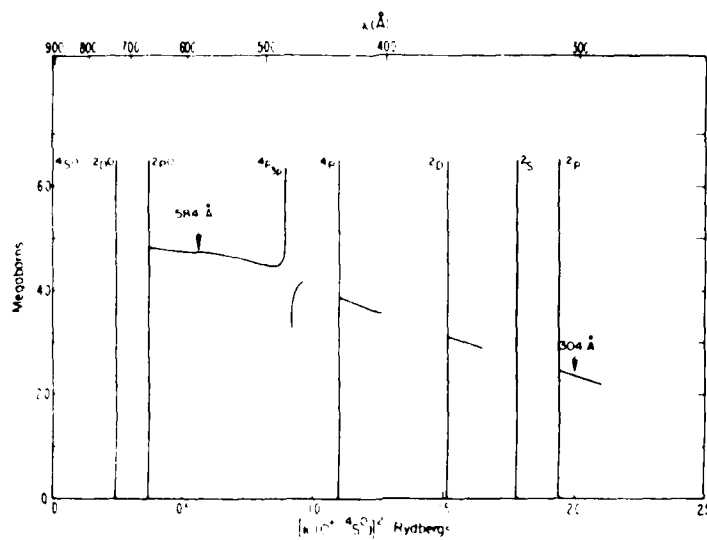
Graphical Data D-1.11

Photoionization cross section for atomic O. (a) Total cross section, dashed line is averaged over resonances, points are experimental, (b), (c), (d) are details of the resonance regions.



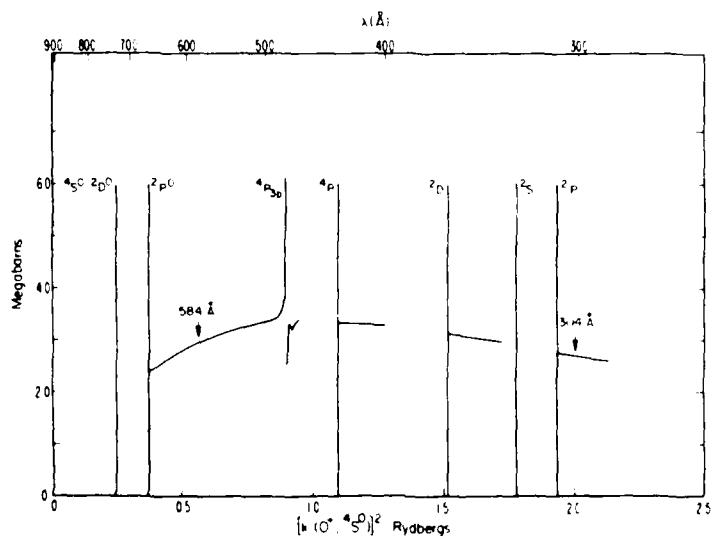
Graphical Data U-1.12

Photoionization cross section for the ground ($1P$) state of O into the ground ($1S$) of O^+ .



Graphical Data D-1.13

Photoionization cross section for the ground $3P$ state of atomic O to the $3D$ state of O^+ .



Graphical Data D-1.14

Photoionization cross section for the ground ^3P state of atomic O to the $^3P^\circ$ state of O^+ .

Tabular Data D-1.15

Photoionization Cross Sections and Branching Ratios
in Various Theoretical Approximations for C^+

Sum of the $3p \rightarrow kd, ks$ cross sections at various energies (in 10^{-18} cm^2).

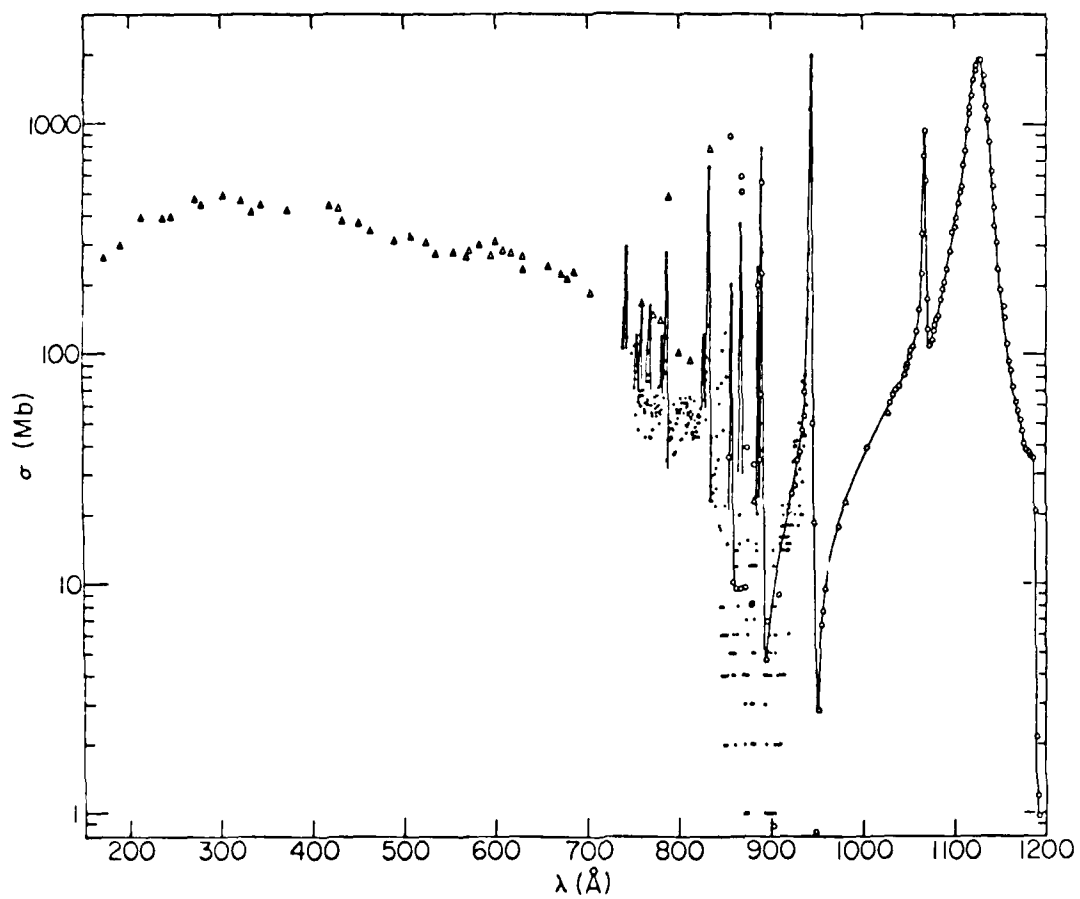
Photon energy (eV)	With relaxation effects				Without relaxation effects			
	Lowest-order HF Length	Velocity	Correlated ^a Length	Velocity	Lowest-order HF Length	Velocity	Correlated ^a Length	Velocity
16.62	56.14	37.05	35.91	34.00	60.40	43.06	39.40	35.99
18	51.66	33.30	36.13	34.71	51.10	35.60	40.10	36.36
20	42.60	26.41	35.38	31.62	38.79	25.92	39.38	35.49
21.2	37.22	22.46	34.32	30.10	32.48	21.11	37.81	34.02
23	29.98	17.32	32.06	29.57	24.33	15.17	33.94	30.55
25	23.17	12.71	28.45	23.34	16.97	10.13	27.52	24.87
27	17.51	9.11	24.17	21.46	11.31	6.49	20.33	18.38
30	11.03	5.30	17.40	12.89	5.87	3.18	11.52	10.37
32	7.75	3.54	13.01	10.94	3.64	1.95	7.17	6.44
35	4.45	1.89	7.87	5.16	1.69	0.92	3.19	2.85
40	1.61	0.70	2.83	2.19	0.68	0.50	0.93	0.85
45	0.57	0.36	0.88	0.68	0.42	0.45	0.38	0.38

^aIncludes correlations from coupled-equations method.

Ratios of the ^1D and ^1S cross sections to ^3P at various energies

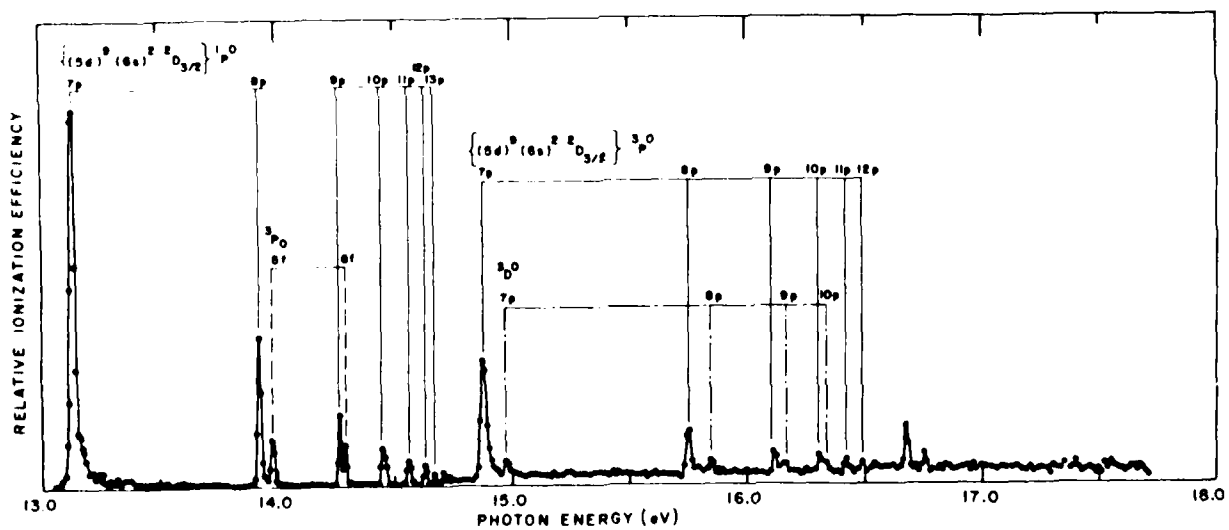
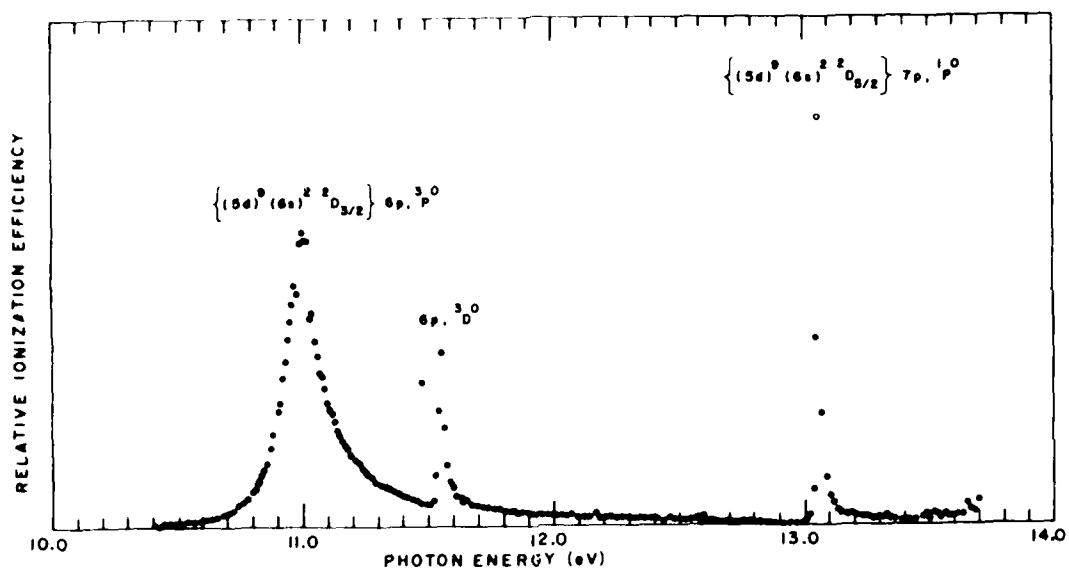
Photon energy (eV)	^{1D} ratio				^{1S} ratio			
	With relaxation effects Lowest-order HF Length	Velocity	Correlated Length	Velocity	Without relaxation effects Lowest-order HF Length	Velocity	Correlated ^a Length	Velocity
18	0.82	0.97	0.66	0.69	1.1	1.2	0.79	0.85
19	0.67	0.82	0.66	0.68	0.95	1.1	0.77	0.82
20	0.55	0.70	0.67	0.70	0.83	0.96	0.75	0.80
21.2	0.45	0.58	0.68	0.72	0.70	0.83	0.74	0.79
22	0.40	0.53	0.70	0.74	0.63	0.76	0.74	0.79
23	0.35	0.47	0.72	0.76	0.55	0.68	0.74	0.78
18	0.24	0.25	0.11	0.14	0.32	0.31	0.15	0.18
19	0.18	0.20	0.11	0.13	0.27	0.28	0.14	0.16
20	0.14	0.16	0.11	0.13	0.22	0.24	0.14	0.16
21.2	0.11	0.13	0.11	0.12	0.18	0.20	0.13	0.14
22	0.09	0.12	0.11	0.12	0.15	0.17	0.12	0.14
23	0.08	0.10	0.11	0.12	0.13	0.15	0.12	0.13

Reference: These data are from E.R. Brown, S.L. Carter, and H.P. Kelly,
Phys. Rev. A **22**, 1987 (1980).



Graphical Data D-1.16

Photoabsorption cross sections for atomic Hg.



Graphical Data D-1.17

Relative photoionization cross section of atomic Hg

Tabular Data D-1.18

Photoionization Cross Sections for Singly Charged Positive Ions

HV	Li+	Be+	B+	HV	C+	N+	O+	F+
6	0.000E 00	0.000E 00	0.000E 00	15	0.000E 00	0.000E 00	0.000E 00	0.000E 00
7	0.000E 00	0.000E 00	0.000E 00	20	0.000E 00	0.000E 00	0.000E 00	0.000E 00
8	0.000E 00	0.000E 00	0.000E 00	25	3.405E 00	0.000E 00	0.000E 00	0.000E 00
9	0.000E 00	0.000E 00	0.000E 00	30	2.388E 00	6.071E 00	0.000E 00	0.000E 00
10	0.000E 00	0.000E 00	0.000E 00	35	3.377E 00	4.808E 00	0.000E 00	0.000E 00
15	0.000E 00	0.000E 00	0.000E 00	40	2.751E 00	5.190E 00	6.571E 00	8.177E 00
20	0.000E 00	1.398E 00	0.000E 00	45	2.277E 00	4.350E 00	5.624E 00	7.541E 00
25	0.000E 00	1.054E 00	2.259E 00	50	1.910E 00	3.670E 00	5.862E 00	6.827E 00
30	0.000E 00	8.055E-01	1.872E 00	60	1.395E 00	2.674E 00	4.473E 00	6.307E 00
35	0.000E 00	5.278E-01	1.550E 00	70	1.059E 00	2.009E 00	3.442E 00	5.115E 00
40	0.000E 00	4.985E-01	1.294E 00	80	8.284E-01	1.550E 00	2.688E 00	4.137E 00
45	0.000E 00	4.027E-01	1.092E 00	90	6.636E-01	1.225E 00	2.132E 00	3.360E 00
50	0.000E 00	3.304E-01	9.306E-01	100	5.413E-01	9.890E-01	1.719E 00	2.749E 00
60	0.000E 00	2.310E-01	6.957E-01					
70	0.000E 00	1.685E-01	5.368E-01					
80	2.554E 00	1.269E-01	4.242E-01					
90	2.035E 00	9.812E-02	3.420E-01					
100	1.613E 00	7.764E-02	2.800E-01					

HV	Ne+	Na+	Mg+	Al+	Si+	P+	S+
15	0.000E 00	0.000E 00	0.000E 00	0.000E 00	1.206E 00	0.000E 00	0.000E 00
20	0.000E 00	0.000E 00	2.541E-01	4.400E-01	2.455E-01	1.272E 00	0.000E 00
25	0.000E 00	0.000E 00	2.279E-01	4.709E-01	6.216E-01	4.092E-01	1.192E 00
30	0.000E 00	0.000E 00	1.954E-01	4.438E-01	6.751E-01	8.118E-01	5.453E-01
35	0.000E 00	0.000E 00	1.664E-01	4.020E-01	6.854E-01	8.878E-01	9.730E-01
40	0.000E 00	0.000E 00	1.422E-01	3.600E-01	6.679E-01	9.291E-01	1.068E 00
45	8.100E 00	0.000E 00	1.226E-01	3.221E-01	6.362E-01	9.359E-01	1.135E 00
50	7.846E 00	7.014E 00	1.067E-01	2.890E-01	5.980E-01	9.189E-01	1.166E 00
60	6.925E 00	8.166E 00	8.287E-02	2.359E-01	5.177E-01	8.441E-01	1.146E 00
70	6.576E 00	7.811E 00	6.289E 00	1.960E-01	4.425E-01	7.488E-01	1.065E 00
80	5.594E 00	7.401E 00	6.854E 00	1.654E-01	3.767E-01	6.527E-01	9.592E-01
90	4.720E 00	6.333E 00	6.557E 00	5.409E 00	3.211E-01	5.649E-01	8.505E-01
100	3.972E 00	5.363E 00	6.559E 00	5.769E 00	2.750E-01	4.886E-01	7.490E-01

Accuracy: These theoretical data should be good to $\pm 20\%$ except near thresholds where they were somewhat worse.

Reference: The above data were taken from R. F. Reilman and S. T. Manson, Astrophysical Jour. Supp. 40, 815 (1979).

Tabular Data D-1.19

Photoionization Cross Sections for Singly
Charged Positive Ions

HV	CL+	A+	K+	CA+	SC+	TI+	V+
15	0.000E 00	0.000E 00	0.000E 00	2.036E-01	3.066E 00	1.650E-01	1.416E-01
20	0.000E 00	0.000E 00	0.000E 00	1.729E-01	3.311E 00	5.100E 00	5.480E 00
25	3.782E 00	0.000E 00	0.000E 00	1.400E-01	3.144E 00	5.356E 00	6.573E 00
30	1.145E 00	3.034E 00	0.000E 00	1.142E-01	2.786E 00	5.181E 00	6.822E 00
35	6.618E-01	1.133E 00	2.605E 00	9.480E-02	2.399E 00	4.713E 00	6.587E 00
40	1.098E 00	7.584E-01	1.240E 00	8.015E-02	2.052E 00	4.188E 00	6.108E 00
45	1.207E 00	1.189E 00	9.036E-01	1.228E 00	1.757E 00	3.694E 00	5.555E 00
50	1.296E 00	1.308E 00	1.341E 00	1.019E 00	2.558E 00	3.252E 00	5.016E 00
60	1.370E 00	1.486E 00	1.588E 00	1.604E 00	2.036E 00	3.354E 00	4.563E 00
70	1.342E 00	1.543E 00	1.709E 00	1.780E 00	2.380E 00	2.872E 00	4.054E 00
80	1.257E 00	1.511E 00	1.708E 00	1.862E 00	2.316E 00	3.025E 00	3.936E 00
90	1.148E 00	1.428E 00	1.641E 00	1.847E 00	2.234E 00	2.803E 00	3.560E 00
100	1.033E 00	1.319E 00	1.546E 00	1.774E 00	2.117E 00	2.604E 00	3.239E 00

HV	CR+	MN+	FE+	CC+
15	8.594E 00	9.228E-02	0.000E 00	0.000E 00
20	9.152E 00	1.303E-01	1.160E-01	1.009E-01
25	9.410E 00	6.603E 00	6.095E 00	5.514E 00
30	9.397E 00	7.728E 00	7.408E 00	6.870E 00
35	9.116E 00	8.336E 00	8.335E 00	7.987E 00
40	8.617E 00	8.462E 00	8.812E 00	8.743E 00
45	7.986E 00	8.247E 00	8.896E 00	9.120E 00
50	7.304E 00	7.847E 00	8.708E 00	9.180E 00
60	7.194E 00	6.881E 00	7.951E 00	8.731E 00
70	5.559E 00	6.742E 00	7.072E 00	7.988E 00
80	4.753E 00	5.740E 00	6.862E 00	7.199E 00
90	4.567E 00	5.058E 00	6.048E 00	6.993E 00
100	4.101E 00	4.882E 00	5.412E 00	6.292E 00

HV	NI+	CU+	ZN+
20	8.522E-02	7.056E 00	5.952E-02
25	4.970E 00	7.969E 00	6.250E-02
30	6.263E 00	8.766E 00	5.151E 00
35	7.451E 00	9.429E 00	6.248E 00
40	8.388E 00	9.916E 00	7.282E 00
45	9.005E 00	1.022E 01	8.148E 00
50	9.306E 00	1.034E 01	8.792E 00
60	9.219E 00	1.020E 01	9.410E 00
70	8.670E 00	9.664E 00	9.371E 00
80	7.987E 00	8.919E 00	8.699E 00
90	7.876E 00	8.832E 00	8.499E 00
100	7.109E 00	7.854E 00	7.650E 00

Accuracy: These theoretical data should be good to $\pm 20\%$ except near thresholds where they were somewhat worse.

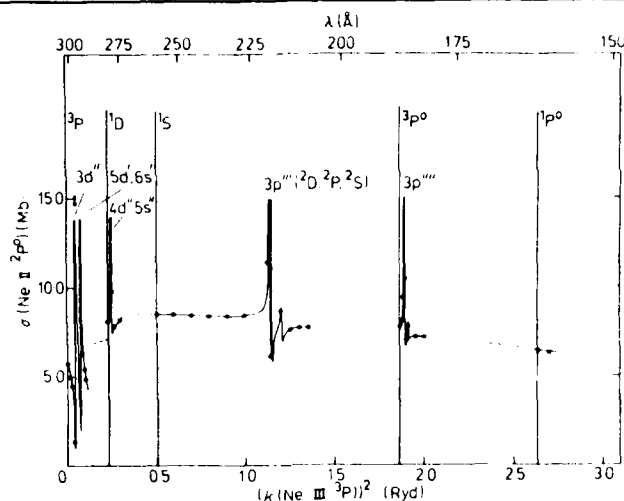
Reference: The above data were taken from R. F. Reilman and S. T. Manson, Astrophysical Jour. Supp. 40, 815 (1979).

Tabular Data D-1.20

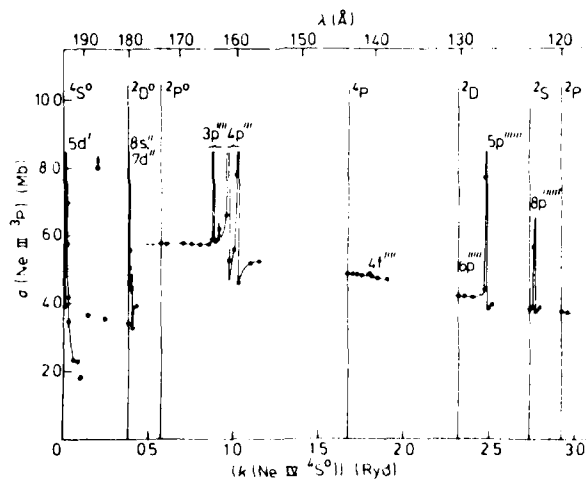
Photoionization Cross Section for Na^+

Threshold values of σ ($\text{Ne II}, 2P^0 \rightarrow \text{Ne III}, S, L, \dots$) in megabarns.

$S, L, \dots$	$E(S, L, \dots)$	$2P^0 \rightarrow 3P$	$2P^0 \rightarrow 1D$	$2P^0 \rightarrow 1S$	$2P^0 \rightarrow 3P^0$	$2P^0 \rightarrow 1P^0$	Total
$3P$	0.0	5.80	—	—	—	—	5.80
$1D$	0.2326	4.52	3.58	—	—	—	8.10
$1S$	0.5051	4.91	3.17	0.48	—	—	8.56
$3P^0$	1.8614	4.15	2.71	0.47	0.48	—	7.80
$1P^0$	2.6350	3.40	2.14	0.38	0.45	0.07	6.43



Graphical Data D-1.21 Photoionisation cross section of the ground state $2P^0$ of Ne II. The averaged value obtained by numerical integration over the resonances $3d$, $5d$ and $6s$ is 6.37 Mb.



Graphical Data D-1.22 Photoionisation cross section of the ground state $1P$ of Ne III. The averaged value over the resonance $5d$ is 3.95 Mb and the averaged value over the resonances $8s$ and $7d$ is 4.75 Mb.

Reference: The above theoretical data were taken from A. K. Pradhan, J. Phys. B 12, 3317 (1979).

D-2. PHOTOABSORPTION, PHOTOIONIZATION, AND PHOTODISSOCIATION CROSS SECTION
OF MOLECULES AND POSITIVE MOLECULAR IONS (MONOMERS)

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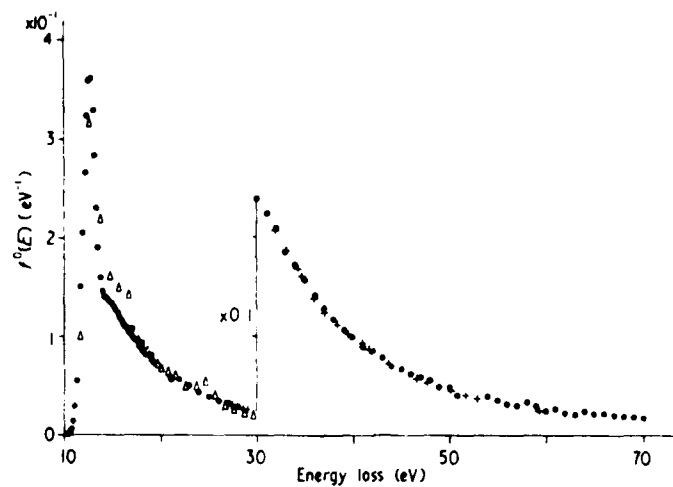
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Tabular Data D-2.1

Photoabsorption and Photoionization Oscillator
Strengths (Cross Sections) for H_2

$f^{(0)}(10^{-2} \text{ eV}^{-1})$			$f^{(0)}(10^{-2} \text{ eV}^{-1})$			$f^{(0)}(10^{-2} \text{ eV}^{-1})$		
E(eV)	Absorption† H_2^+		E(eV)	Absorption† H_2^+		E(eV)	Absorption† H_2^+	
10.0	0.08		20	6.95	6.82	40	0.99	0.89
10.5	0.26		21	6.25	6.13	42	0.84	0.75
11.0	1.70		22	5.51	5.40	44	0.71	0.63
11.5	7.66		23	4.92	4.82	46	0.65	0.58
12.0	20.7		24	4.38	4.30	48	0.61	0.54
12.5	34.1		25	3.92	3.83	50	0.48	0.42
13.0	32.5		26	3.49	3.42	52	0.41	0.36
13.5	21.2		27	3.24	3.15	54	0.39	0.34
14.0	14.9		28	2.89	2.81	56	0.32	0.28
14.5	13.8		29	2.68	2.58	58	0.33	0.28
15.0	13.3	0.53	30	2.46	2.35	60	0.24	0.20
15.5	12.6	4.73	31	2.27	2.15	62	0.23	0.20
16.0	11.7	8.68	32	2.06	1.93	64	0.23	0.20
16.5	10.8	9.74	33	1.90	1.76	66	0.21	0.18
17.0	10.1	9.63	34	1.72	1.58	68	0.19	0.17
17.5	9.58	9.60	35	1.55	1.36	70	0.16	0.14
18.0	8.83	8.91	36	1.39	1.18			
18.5	8.25	8.50	37	1.31	1.14			
19.0	7.85	7.97	38	1.29	1.04			
19.5	7.25	7.41	39	1.10	0.94			

Reference: These data were taken from C. Backx, G.R. Wight, and
M.J. Van der Wiel, J. Phys. B 9, 315 (1976).



Graphical Data D-2.2

Photoabsorption oscillator strength (cross section) for H_2 .

Tabular Data D-2.3

Photofragmentation Ratios for H⁺ and D⁺

HeV	H ⁺ H ₂ ⁺	D ⁺ D ₂ ⁺	HeV	H ⁺ H ₂ ⁺	D ⁺ D ₂ ⁺	HeV	H ⁺ H ₂ ⁺	D ⁺ D ₂ ⁺
($\times 10^{-2}$)	($\times 10^{-2}$)	($\times 10^{-2}$)	($\times 10^{-2}$)	($\times 10^{-2}$)	($\times 10^{-2}$)	($\times 10^{-2}$)	($\times 10^{-2}$)	($\times 10^{-2}$)
18.0	0.33	0.20	22.0	2.04	0.66	30	4.44	2.15
18.2	0.72	0.29	22.2	1.96	0.65	31	5.36	3.21
18.4	1.17	0.31	22.4	1.87	0.76	32	6.96	4.38
18.6	1.08	0.53	22.6	1.97	0.55	33	7.95	6.92
18.8	1.43	0.53	22.8	1.92	0.58	34	8.65	8.40
19.0	1.52	0.52	23.0	1.97	0.78	35	9.49	8.65
19.2	1.57	0.65	23.2	1.85	0.52	36	9.29	8.25
19.4	1.59	0.48	23.4	2.08	0.84	37	7.96	7.32
19.6	1.70	0.60	23.6	2.21	0.79	38	7.78	7.07
19.8	1.86	0.48	23.8	2.03	0.76	39	8.74	8.04
20.0	1.72	0.70	24.0	2.00	0.67	40	9.94	9.21
20.2	1.90	0.64	24.2	1.98	0.88	45	12.1	11.4
20.4	1.94	0.74	24.4	2.04	0.60	50	13.6	11.7
20.6	1.79	0.60	24.6	2.12	0.89	55	15.4	12.8
20.8	1.88	0.75	24.8	2.23	0.56	60	16.8	13.4
21.0	1.86	0.62	25.0	2.04	0.85	65	14.4	13.8
21.2	1.88	0.64	26.0	2.21	0.94	70	17.6	14.8
21.4	1.94	0.63	27.0	2.46	1.12			
21.6	1.92	0.76	28.0	3.22	1.56			
21.8	1.92	0.77	29.0	3.92	1.85			

Reference: These data were taken from C. Backx, G.R. Wight, and M.J. Van der Wiel, J. Phys. B 9, 315 (1976).

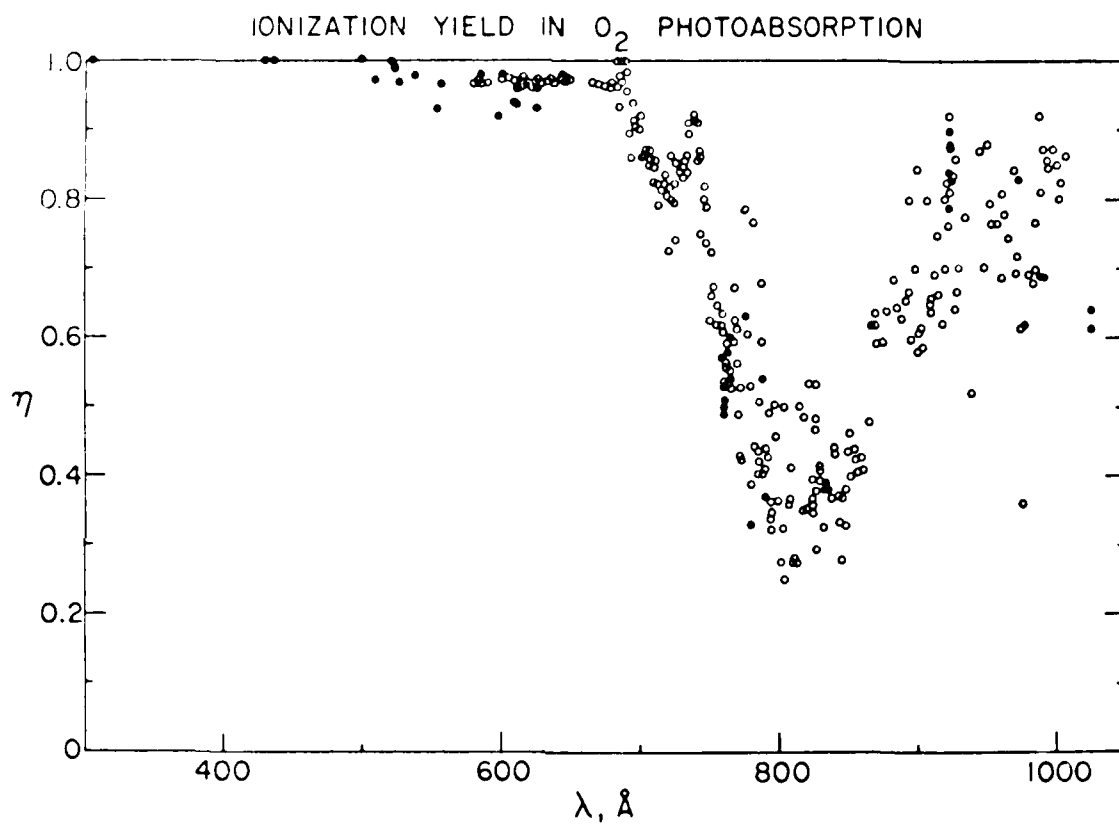
Tabular Data D-2.4

Photoabsorption Cross Sections for N₂, CO, NO, O₂, and N₂O

λ , Å	Cross Section (Mb) $\times 10^{-17}$ cm ²				
	N ₂	CO	NO	O ₂	N ₂ O
20	0.18		0.29	0.39	0.38
25	0.25		0.37	0.47	0.45
30	0.32		0.46	0.57	0.56
35	0.39		0.54	0.67	0.65
40	0.46		0.63	0.77	0.75
45	0.55		0.73	0.87	0.85
50	0.64		0.82	0.97	0.95
55	0.75		0.92	1.07	1.05
60	0.85		1.02	1.17	1.15
65	0.95		1.12	1.27	1.25
70	1.05		1.22	1.37	1.35
75	1.15		1.32	1.47	1.45
80	1.25		1.42	1.57	1.55
85	1.35		1.52	1.67	1.65
90	1.45		1.62	1.77	1.75
95	1.55		1.72	1.87	1.85
100	1.65		1.82	1.97	1.95
105	1.75	3.00	1.92	2.07	2.05
110	1.85	3.15	2.02	2.17	2.15
115	1.95	3.30	2.12	2.27	2.25
120	2.05	3.45	2.22	2.37	2.35
125	2.15	3.60	2.32	2.47	2.45
130	2.25	3.75	2.42	2.57	2.55
135	2.35	3.90	2.52	2.67	2.65
140	2.45	4.05	2.62	2.77	2.75
145	2.55	4.20	2.72	2.87	2.85
150	2.65	4.35	2.82	2.97	2.95
155	2.75	4.50	2.92	3.07	3.05
160	2.85	4.65	3.02	3.17	3.15
165	2.95	4.80	3.12	3.27	3.25
170	3.05	4.95	3.22	3.37	3.35
175	3.15	5.10	3.32	3.47	3.45
180	3.25	5.25	3.42	3.57	3.55
185	3.35	5.40	3.52	3.67	3.65
190	3.45	5.55	3.62	3.77	3.75
195	3.55	5.70	3.72	3.87	3.85
200	3.65	5.85	3.82	3.97	3.95
205	3.75	6.00	3.92	4.07	4.05
210	3.85	6.15	4.02	4.17	4.15
215	3.95	6.30	4.12	4.27	4.25
220	4.05	6.45	4.22	4.37	4.35
225	4.15	6.60	4.32	4.47	4.45
230	4.25	6.75	4.42	4.57	4.55
235	4.35	6.90	4.52	4.67	4.65
240	4.45	7.05	4.62	4.77	4.75
245	4.55	7.20	4.72	4.87	4.85
250	4.65	7.35	4.82	4.97	4.95
255	4.75	7.50	4.92	5.07	5.05
260	4.85	7.65	5.02	5.17	5.15
265	4.95	7.80	5.12	5.27	5.25
270	5.05	7.95	5.22	5.37	5.35
275	5.15	8.10	5.32	5.47	5.45
280	5.25	8.25	5.42	5.57	5.55
285	5.35	8.40	5.52	5.67	5.65
290	5.45	8.55	5.62	5.77	5.75
295	5.55	8.70	5.72	5.87	5.85
300	5.65	8.85	5.82	5.97	5.95
305	5.75	9.00	5.92	6.07	6.05
310	5.85	9.15	6.02	6.17	6.15
315	5.95	9.30	6.12	6.27	6.25
320	6.05	9.45	6.22	6.37	6.35
325	6.15	9.60	6.32	6.47	6.45
330	6.25	9.75	6.42	6.57	6.55
335	6.35	9.90	6.52	6.67	6.65
340	6.45	10.05	6.62	6.77	6.75
345	6.55	10.20	6.72	6.87	6.85
350	6.65	10.35	6.82	6.97	6.95
355	6.75	10.50	6.92	7.07	7.05
360	6.85	10.65	7.02	7.17	7.15
365	6.95	10.80	7.12	7.27	7.25
370	7.05	10.95	7.22	7.37	7.35
375	7.15	11.10	7.32	7.47	7.45
380	7.25	11.25	7.42	7.57	7.55
385	7.35	11.40	7.52	7.67	7.65
390	7.45	11.55	7.62	7.77	7.75
395	7.55	11.70	7.72	7.87	7.85
400	7.65	11.85	7.82	7.97	7.95
405	7.75	12.00	7.92	8.07	8.05
410	7.85	12.15	8.02	8.17	8.15
415	7.95	12.30	8.12	8.27	8.25
420	8.05	12.45	8.22	8.37	8.35
425	8.15	12.60	8.32	8.47	8.45
430	8.25	12.75	8.42	8.57	8.55
435	8.35	12.90	8.52	8.67	8.65
440	8.45	13.05	8.62	8.77	8.75
445	8.55	13.20	8.72	8.87	8.85
450	8.65	13.35	8.82	8.97	8.95
455	8.75	13.50	8.92	9.07	9.05
460	8.85	13.65	9.02	9.17	9.15
465	8.95	13.80	9.12	9.27	9.25
470	9.05	13.95	9.22	9.37	9.35
475	9.15	14.10	9.32	9.47	9.45
480	9.25	14.25	9.42	9.57	9.55
485	9.35	14.40	9.52	9.67	9.65
490	9.45	14.55	9.62	9.77	9.75
495	9.55	14.70	9.72	9.87	9.85
500	9.65	14.85	9.82	9.97	9.95
505	9.75	15.00	9.92	10.07	10.05
510	9.85	15.15	10.02	10.17	10.15
515	9.95	15.30	10.12	10.27	10.25
520	10.05	15.45	10.22	10.37	10.35
525	10.15	15.60	10.32	10.47	10.45
530	10.25	15.75	10.42	10.57	10.55
535	10.35	15.90	10.52	10.67	10.65
540	10.45	16.05	10.62	10.77	10.75
545	10.55	16.20	10.72	10.87	10.85
550	10.65	16.35	10.82	10.97	10.95
555	10.75	16.50	10.92	11.07	11.05
560	10.85	16.65	11.02	11.17	11.15
565	10.95	16.80	11.12	11.27	11.25
570	11.05	16.95	11.22	11.37	11.35
575	11.15	17.10	11.32	11.47	11.45
580	11.25	17.25	11.42	11.57	11.55
585	11.35	17.40	11.52	11.67	11.65
590	11.45	17.55	11.62	11.77	11.75
595	11.55	17.70	11.72	11.87	11.85
600	11.65	17.85	11.82	11.97	11.95
605	11.75	18.00	11.92	12.07	12.05
610	11.85	18.15	12.02	12.17	12.15
615	11.95	18.30	12.12	12.27	12.25
620	12.05	18.45	12.22	12.37	12.35
625	12.15	18.60	12.32	12.47	12.45
630	12.25	18.75	12.42	12.57	12.55
635	12.35	18.90	12.52	12.67	12.65
640	12.45	19.05	12.62	12.77	12.75
645	12.55	19.20	12.72	12.87	12.85
650	12.65	19.35	12.82	12.97	12.95
655	12.75	19.50	12.92	13.07	13.05
660	12.85	19.65	13.02	13.17	13.15
665	12.95	19.80	13.12	13.27	13.25
670	13.05	19.95	13.22	13.37	13.35
675	13.15	20.10	13.32	13.47	13.45
680	13.25	20.25	13.42	13.57	13.55
685	13.35	20.40	13.52	13.67	13.65
690	13.45	20.55	13.62	13.77	13.75
695	13.55	20.70	13.72	13.87	13.85
700	13.65	20.85	13.82	13.97	13.95
705	13.75	21.00	13.92	14.07	14.05
710	13.85	21.15	14.02	14.17	14.15
715	13.95	21.30	14.12	14.27	14.25
720	14.05	21.45	14.22	14.37	14.35
725	14.15	21.60	14.32	14.47	14.45
730	14.25	21.75	14.42	14.57	14.55
735	14.35	21.90	14.52	14.67	14.65
740	14.45	22.05	14.62	14.77	14.75
745	14.55	22.20	14.72	14.87	14.85
750	14.65	22.35	14.82	14.97	14.95
755	14.75	22.50	14.92	15.07	15.05
760	14.85	22.65	15.02	15.17	15.15
765	14.95	22.80	15.12	15.27	15.25
770	15.05	22.95	15.22	15.37	15.35
775	15.15	23.10	15.32	15.47	15.45
780	15.25	23.25	15.42	15.57	15.55
785	15.35	23.40	15.52	15.67	15.65
790	15.45	23.55	15.62	15.77	15.75
795	15.55	23.70	15.72	15.87	15.85
800	15.65	23.85	15.82	15.97	15.95
805	15.75	24.00	15.92	16.07	16.05
810	15.85	24.15	16.02	16.17	16.15
815	15.95	24.30	16.12	16.27	16.25
820	16.05	24.45	16.22	16.37	16.35
825	16.15	24.60	16.32	16.47	16.45
830	16.25	24.75	16.42	16.57	16.55
835	16.35	24.90	16.52	16.67	16.65
840	16.45	25.05	16.62	16.77	16.75
845	16.55	25.20	16.72	16.87	16.85
850	16.65	25.35	16.82	16.97	16.95
855	16.75	25.50	16.92	17.07	17.05
860	16.85	25.65	17.02	17.17	17.15
865	16.95	25.80	17.12	17.27	17.25
870	17.05	25.95	17.22	17.37	17.35
875	17.15	26.10	17.32	17.47	17.45
880	17.25	26.25	17.42	17.57	17.55
885	17.35	26.40	17.52	17.67	17.65
890	17.45	26.55	17.62	17.77	17.75
895	17.55	26.70	17.72	17.87	17.85
900	17.65	26.85	17.82	17.97	17.95
905	17.75	27.00	17.92	18.07	18.05
910	17.85	27.15	18.02	18.17	18.15
915	17.95	27.30	18.12	18.27	18.25
920	18.05	27.45	18.22	18.37	18.35
925	18.15	27.60	18.32	18.47	18.45
930	18.25	27.75	18.42	18.57	18.55
935	18.35	27.90	18.52	18.67	18.65
940	18.45	28.05	18.62	18.77	18.75
945	18.55	28.20	18.72	18.87	18.85
950	18.65	28.35	18.82	18.97	18.95
955	18.75	28.50	18.92	19.07	19.05
960	18.85	28.65	19.02	19.17	19.15
965	18.95	28.80	19.12	19.27	19.25
970	19.05	28.95	19.22	19.37	19.35
975	19.15	29.10	19.32	19.47	19.45
980	19.25	29.25	19.42	19.57	19.55
985	19.35	29.40	19.52	19.67	19.65
990	19.45	29.55	19.62	19.77	19.75
995	19.55	29.70	19.72	19.87	19.85
1000	19.65	29.85	19.82	19.97	19.95

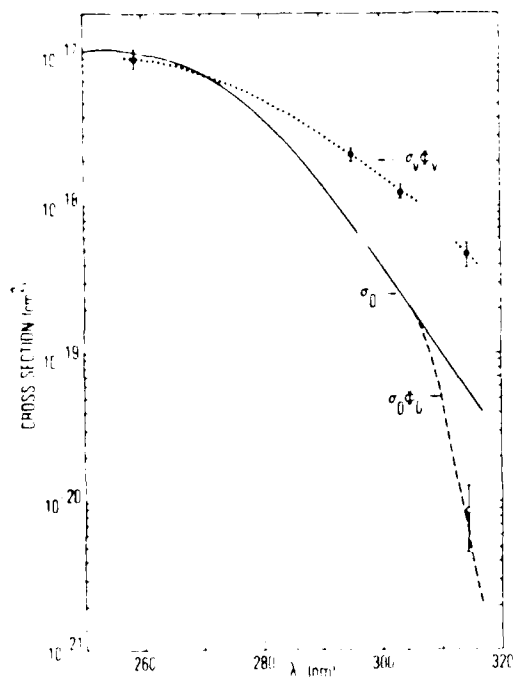
Accuracy: The quoted accuracy is 10% absolute and 2% relative.

Reference: The data are from B.H. Cole and P.G. Dexter, J. Phys. B 11, 1011 (1978).



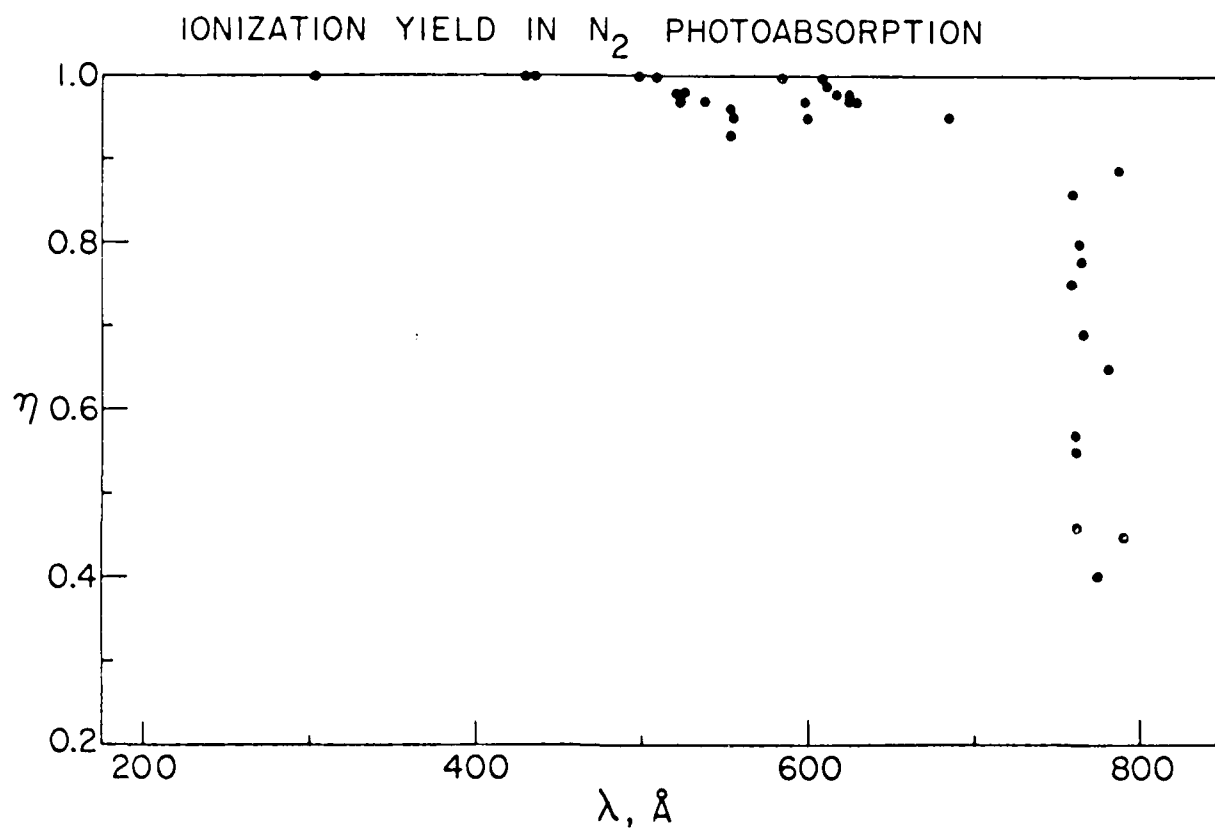
Graphical Data D-7.5

Quantum yield of ionization, η , for O_2 .



Graphical Data D-2.6

Cross section for photodissociative production of O(1D) from vibrationally excited and room temperature O₂.



Graphical Data D-1.7

Quantum yield of ionization, η , for N_2

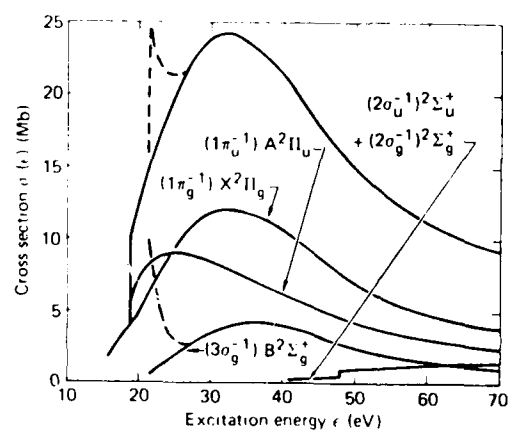
Tabular Data D-2.3

Branching Ratios in the Photoionization of N

Energy loss E (eV)	Electronic state of N ₂ ⁺				Oscillator strength f ₀ for total photoabsorption (eV) ⁻¹
	X ² Σ _g ⁺	4 ¹ Π _g	B ² Σ _g ⁺	Z	
18	34.8	65.2			0.227
19	32.7	67.3			0.217
20	37.8	56.5	5.7		0.206
21	36.0	56.3	7.7		0.200
22	34.0	57.0	9.0		0.207
23	32.3	58.6	9.1		0.225
24	35.0	56.0	9.0		0.219
25	39.5	51.5	9.0		0.212
27	41.6	51.6	6.8		0.207
30	43.1	47.9	9.0		0.189
33	36.1	49.0	11.9	3.0	0.156
35	31.9	51.0	12.8	4.4	0.141
37	31.3	51.9	11.6	5.3	0.129
40	28.0	50.4	12.3	9.3	0.112
42	26.7	44.1	11.5	9.3	8.2 0.105
45	26.1	45.5	9.3	5.7	12.7 0.097
47	28.1	42.0	10.1	4.5	15.4 0.097
50	26.5	36.3	11.0	6.4	19.9 0.090

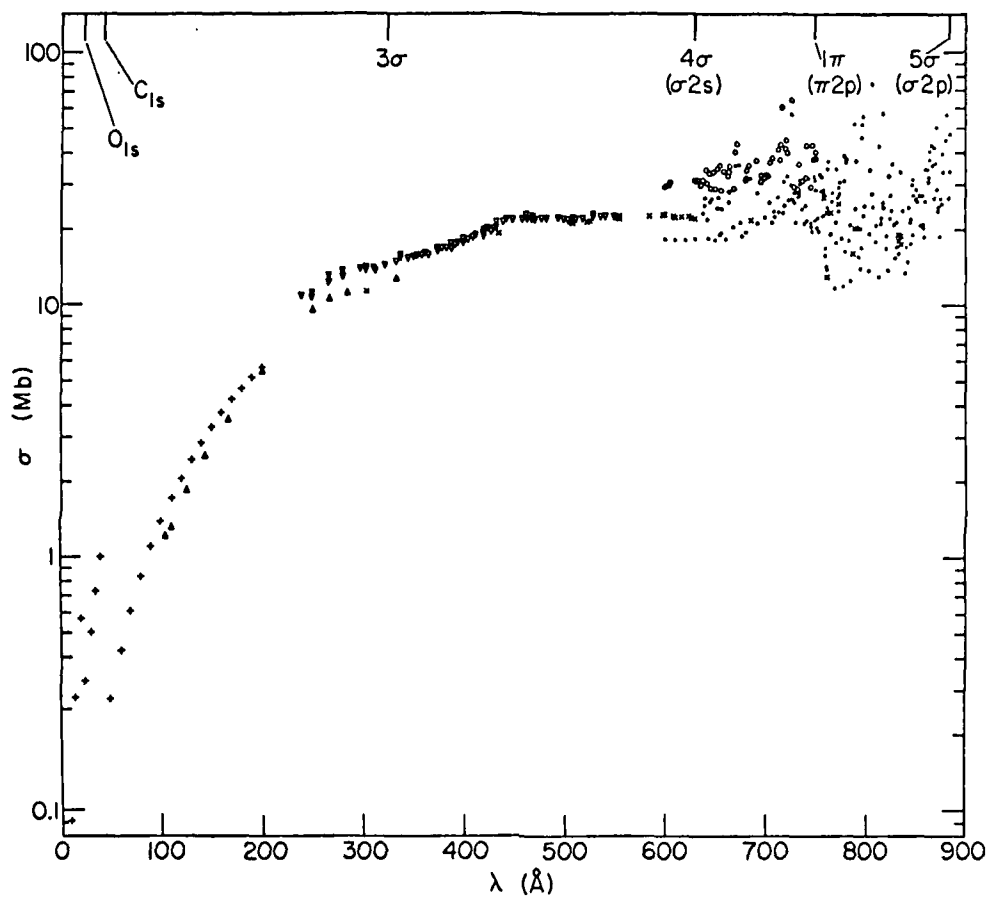
Note: Partial Oscillator Strengths for individual ionic states can be obtained from the product of the branching ratio (given in %) and the f₀ value given in the last column of the table. Oscillator strengths in (eV)⁻¹ can be converted into cross sections in Mb (10⁻¹⁸ cm²) by multiplying by 109.75.

Reference: The above data were taken from A. Hammett, W. Stoll, and C.F. Brion, J. Electron Spectrosc. 8, 367 (1976).



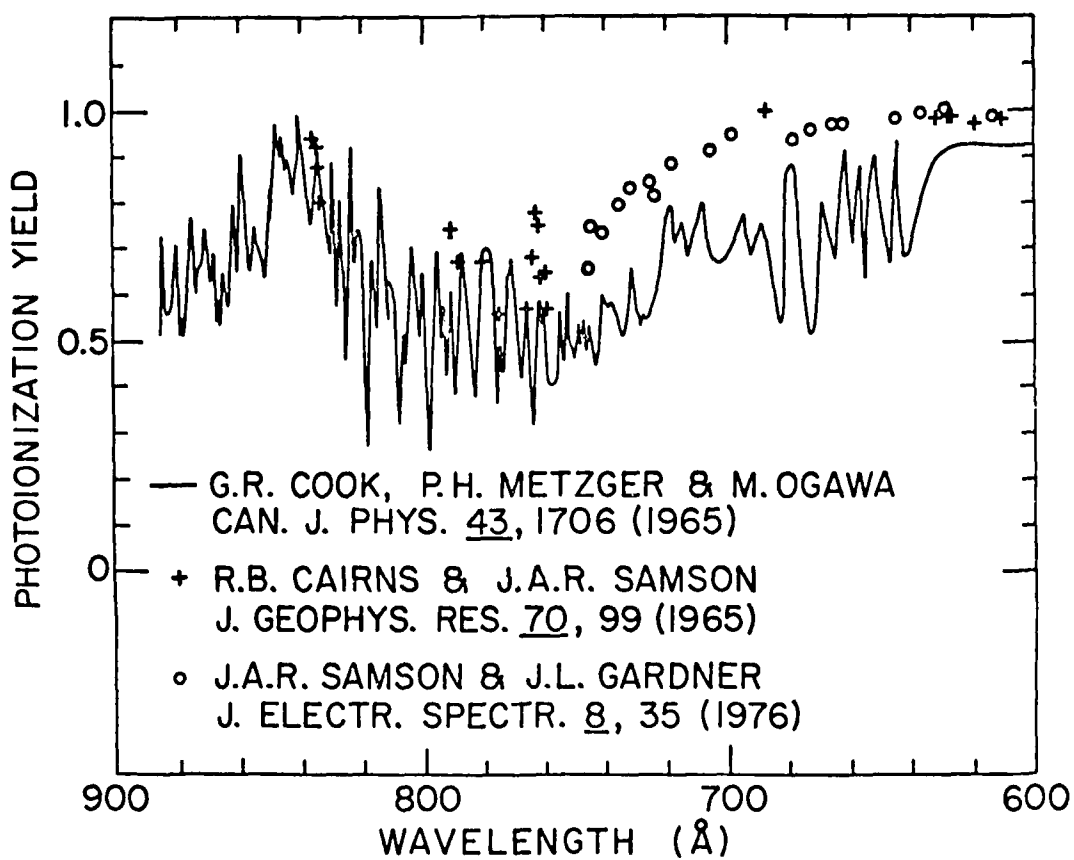
Graphical Data D-2.9

Theoretical total and partial photoionization cross sections of F_2^+ for production of various states of F_2^+ .



Graphical Data D-2.10

Photoabsorption cross section of CO.



Graphical Data D-2.11

Quantum yield of ionization, η , for CO.

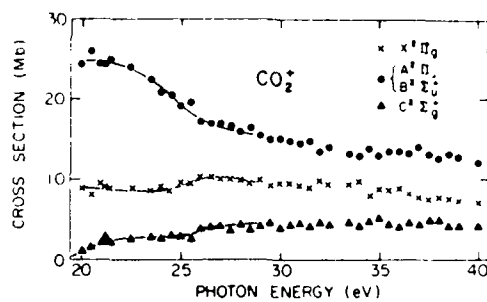
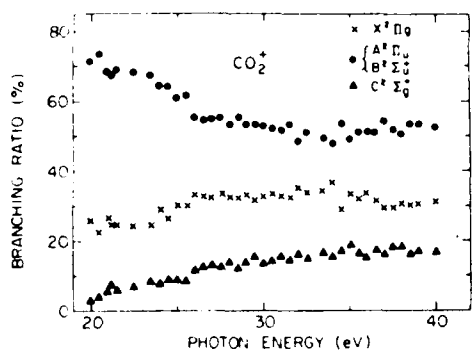
Tabular Data D-2.12

Branching Ratios in the Photoionization of CO

Energy loss E (eV)	Electronic state of CO ⁺						Oscillator strength f ₀ for total photoabsorption ^b (eV) ⁻¹
	X ² Σ ⁺	A ² Π	B ² Σ ⁺	C ² Σ ⁺	W	² Σ _g ⁺ O ₂ N ⁺	
18	50	50					0.231
19	44	56					0.215
20	41.5	58.5					0.213
21	34.0	57.4	8.6				0.218
22	29.8	60.1	10.1				0.219
25	33.0	51.0	16.0				0.210
27	27.7	48.0	19.2	5.1			0.203
30	23.5	46.0	26.5	4.0			0.171
33	18.9	47.3	29.8	4.0			0.154
35	17.2	43.8	28.8	3.4	6.8		0.142
38	15.7	45.9	27.4	3.1	7.9		0.129
40	15.8	46.5	26.5	2.7	8.5		0.122
42	14.5	44.2	23.8	2.4	6.7	8.4	0.116
45	14.7	41.0	23.0	2.9	7.0	11.4	0.110
47	15.2	39.8	22.8	2.7	7.0	12.5	0.103
50	13.0	39.6	24.2	2.8	7.0	13.4	0.097

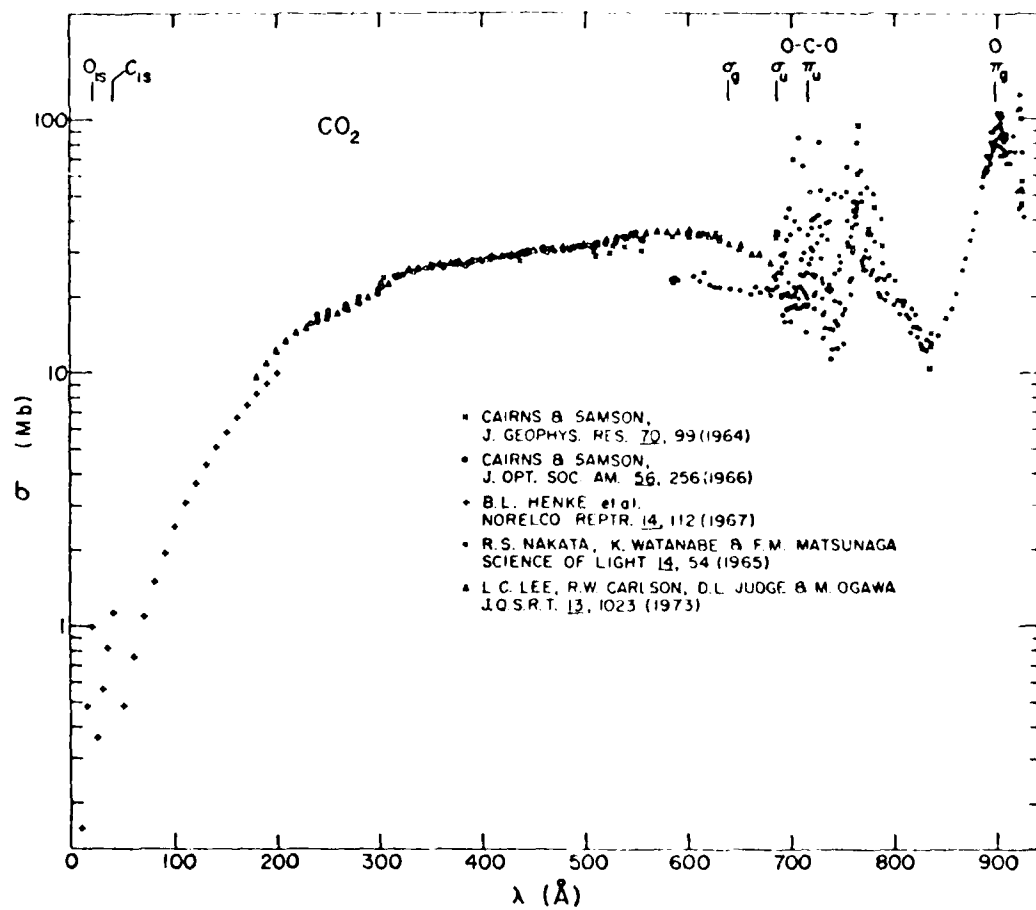
Note: Partial oscillator strengths for individual ionic states can be obtained from the product of the branching ratio (given in %) and the f₀ value given in the last column of the table. Oscillator strengths in (eV)⁻¹ can be converted into cross sections in Mb (10⁻¹⁸ cm²) by multiplying by 109.75.

Reference: The above data were taken from A. Hamnett, W. Stoll, and C.E. Brion, J. Electron Spectrosc. 8, 367 (1976).



Graphical Data D-2.13

Partial photoionization cross sections and branching ratios for production of various states of CO_2^+ in the photoionization of CO_2 .



Graphical Data D-2.14

Photoabsorption cross section of CO₂.

Tabular Data D-1.15

Photoabsorption and Partial Photionization Oscillator
Strengths (Cross Sections) for CO

Energy (eV)	Partial oscillator strength (10^{-2} eV^{-1})			Total MET	Total photoabsorption (10^{-2} eV^{-1})
	$A^2\Pi_p$	$A^2\Pi_u + B^2\Sigma_u$	$C^2\Sigma_g$		
21.2	9.37	22.07	2.00		33.44
22	8.41	21.03	2.59		32.35
23	9.07	17.85	3.33		30.25
24	10.20	16.33	2.62		29.16
25	9.94	14.91	2.76		27.61
26	9.65	14.08	2.34	1.08	27.15
27	9.44	14.02	2.70	1.08	26.97
28	8.20	13.17	3.69	1.31	26.33
29	8.05	12.98	3.38	1.56	25.97
30	7.40	12.50	3.32	2.30	25.51
31	7.60	12.27	3.68	1.71	24.51
32	7.41	11.61	3.71	1.98	24.69
34	7.73	10.62	3.86	1.93	24.15
36	7.24	10.38	3.14	3.14	24.15
38	6.25	10.05	3.12	3.12	22.32
40	5.81	8.82	2.61	2.81	20.06
41	5.90	8.19	1.90	3.04	19.04
42	5.41	7.75	1.80	2.89	18.04
44	4.89	6.85	1.63	2.93	16.31
46	4.34	6.42	1.19	2.99	14.94
48	4.14	5.76	1.03	3.69	14.76
50	4.05	5.50	1.30	3.63	14.49
52	3.89	4.89	1.15	4.46	14.40
54	3.34	4.74	1.26	4.60	13.94
56	3.56	4.36	1.32	3.84	13.21
58	3.37	3.84	1.20	3.73	12.03
60	2.87	3.31	1.10	3.74	11.02

$\sigma(\text{Mb}) = 109.75 (d\sigma/dE \text{ eV}^{-1})$

Note: MET refers to multiple electron transitions.

Reference: These data were taken from C.E. Brion and K.H. Tan, Chem. Phys. 34, 141 (1978).

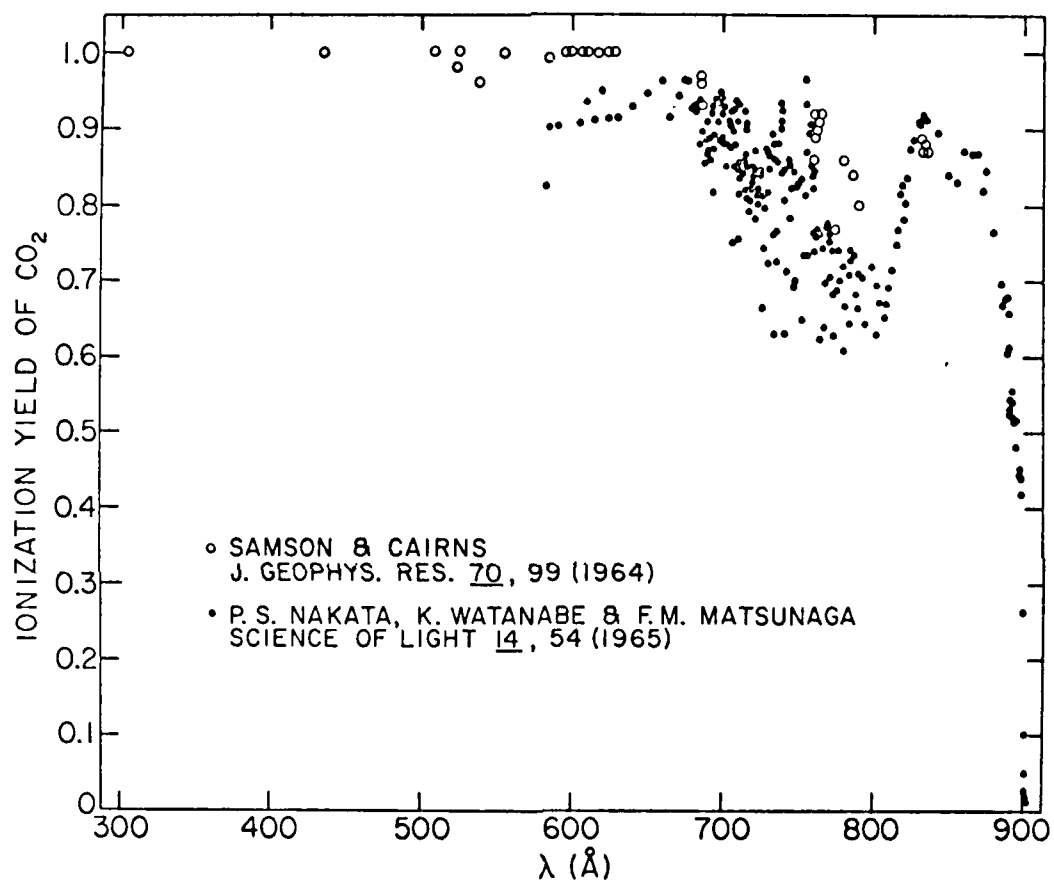
Tabular Data D-2.16

Photoabsorption and Photofragmentation Oscillator
Strengths (Cross Sections) for CO

Energy (eV)	Ionisation efficiency	Oscillator strength ($\times 10^{-2}$ eV $^{-1}$)					
		absorption	CO $_2^+$	CO $^+$	O $^+$	C $^+$	CO $_2^+$
8		0.5					
9		0.9					
10		1.4					
11		17.0					
12	0	32.3	0				
13	0.02	48.3	0.97				
14	0.19	52.0	9.80				
15	0.63	24.3	15.2				
16	0.69	28.5	19.5	0			
17	0.81	33.0	26.4	0.05	0		
18	0.93	30.6	28.1	0.05	0.03		
19	1.05	32.0	33.0	0.11	0.15		
20	1.02	32.0	31.7	0.05	0.62		
21	1.02	32.5	31.5	0.27	1.10		
22	1.01	32.3	30.7	0.38	1.24		
23	1.03	30.4	29.2	0.41	1.38		
24	1.05	27.8	26.7	0.77	1.46		
25	1.01	27.8	25.0	0.11	1.60		
26	1.02	27.1	23.8	1.62	1.88		
27	0.95	26.8	21.1	2.22	1.94	0	
28	0.98	26.4	20.6	2.90	2.15	0.04	
29	0.96	25.9	19.3	3.11	2.27	0.02	
30	0.96	25.8	18.8	3.26	2.49	0.13	
31	0.95	25.2	18.5	3.11	2.57	0.24	
32	0.99	23.6	17.7	2.94	2.65	0.44	-0.02
33	0.97	23.3	16.4	2.69	2.62	0.72	0.02
34	1.03	22.9	16.6	2.78	2.93	1.13	0.01
35	1.03	22.0	15.7	2.59	2.97	1.39	0
36	0.99	22.3	15.0	2.34	3.07	1.56	0.01
37	1.02	21.6	14.7	2.35	3.15	1.70	-0.03
38	1.02	21.0	14.2	2.14	3.23	1.79	-0.02
39	1.01	20.5	13.7	2.01	3.19	1.75	0.02
40	0.96	20.2	12.9	1.83	2.97	1.61	0.02
41	1.00	19.4	13.0	1.76	2.95	1.53	0.06
42	1.01	18.6	12.8	1.66	2.78	1.41	0.07
43	0.94	17.9	11.6	1.50	2.37	1.18	0.07
44	1.00	17.0	11.9	1.48	2.31	1.16	0.10
45	0.98	16.4	11.3	1.41	2.16	1.09	0.07
46	1.00	15.8	11.1	1.42	2.07	1.09	0.10
47	0.94	15.2	10.0	1.29	1.90	0.96	0.07
48	0.98	15.3	10.4	1.33	2.07	1.07	0.10
49	0.97	14.6	9.63	1.34	2.03	1.05	0.09
50	0.98	14.8	9.48	1.46	2.24	1.20	0.09
51	0.99	14.8	9.36	1.56	2.38	1.25	0.09
52	1.01	14.3	8.90	1.55	2.48	1.33	0.12
53	1.07	13.2	7.85	1.61	2.45	1.35	0.11
56	0.94	12.9	7.35	1.48	2.53	1.41	0.13
58	1.03	12.1	6.75	1.53	2.58	1.45	0.16
60	0.96	11.6	5.99	1.34	2.30	1.30	0.14
62	[1.00]	10.0	5.44	1.16	2.09	1.16	0.15
64	[1.00]	9.07	4.93	1.10	1.89	1.03	0.11
66	[1.00]	8.78	4.62	0.99	1.84	0.99	0.12
68	[1.00]	8.41	4.63	0.96	1.77	0.94	0.12
70	[1.00]	7.82	4.30	0.91	1.67	0.88	0.08
72	[1.00]	7.82	4.09	0.86	1.62	0.87	0.09
74	[1.00]	7.02	3.79	0.80	1.53	0.81	0.10
76	[1.00]	6.63	3.56	0.73	1.47	0.77	0.11
78	[1.00]	6.30	3.30	0.73	1.47	0.74	0.08
80	[1.00]	6.02	3.16	0.68	1.40	0.71	0.07

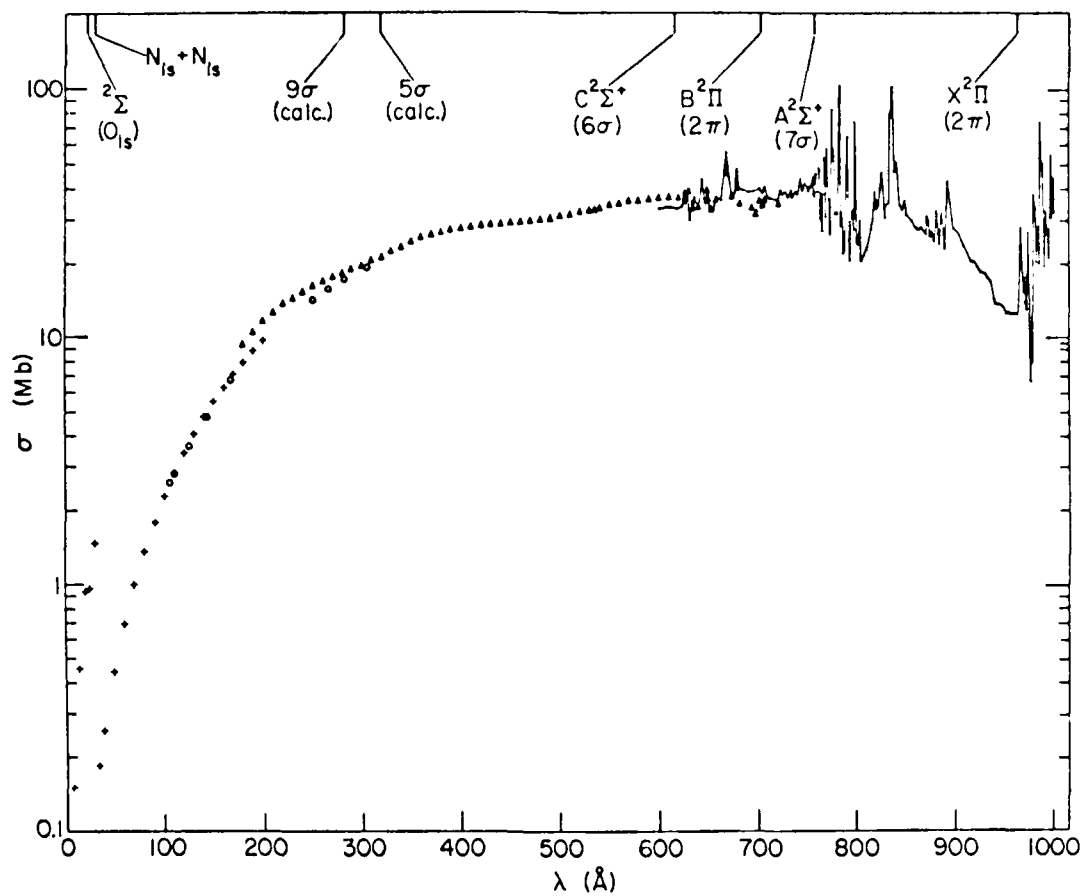
$\sigma(10^{-18} \text{ cm}^2) = 109.75 \text{ df/df (eV}^{-1}\text{)}$

Reference: These data were taken from A.P. Hitchcock, C.E. Brion,
and M.J. Van der Wiel, Chem. Phys. 45, 461 (1980).



Graphical Data D-2.17

Quantum yield of ionization for CO₂ .



Graphical Data D-2.18

Photoabsorption cross section for N O.

Tabular Data D-2.19

Photoabsorption and Partial Photoionization Oscillator
Strengths (Cross Sections) for N₂O

Energy (eV)	Partial oscillator strength (10^{-2} eV ⁻¹)					Total absorption (10^{-2} eV ⁻¹)
	X ² Π	A ² Σ ⁺	B ² Π	C ² Σ ⁺	24 eV	
17	23.22	18.24				41.46
19	15.58	15.21	6.30			37.08
19.6	14.80	12.63	9.02			36.08
21.2	11.66	8.34	8.67	5.00		33.35
22	10.55	6.08	8.32	7.03		31.98
23	9.86	5.68	8.07	6.28		29.89
24	9.73	5.44	7.73	6.00		28.61
25	10.20	5.10	7.65	4.54	0.85	28.34
26	9.57	4.92	7.11	4.37	1.64	27.33
27	9.78	4.88	7.06	4.35	1.36	27.15
28	9.55	4.51	6.90	3.97	1.86	26.51
29	8.89	4.97	6.80	3.66	1.82	26.15
30	8.80	4.65	6.72	3.63	1.81	25.88
32	8.61	4.56	6.33	3.54	1.52	25.33
34	8.43	3.72	6.20	3.72	1.73	24.78
36	7.34	3.02	5.40	3.02	1.08	21.59
38	6.89	2.95	4.33	2.75	1.18	19.68
40	6.35	2.81	4.11	2.61	0.75	18.68
41	6.25	2.32	3.75	2.50	1.43	17.86
42	6.16	2.64	3.52	2.63	1.06	17.59
44	5.26	2.20	3.39	2.54	0.85	16.95
46	5.08	2.22	3.02	2.38	0.95	15.85
48	5.14	2.02	2.65	2.33	0.62	15.58
50	4.60	1.65	2.67	2.08	0.59	14.85
52	4.44	1.86	2.43	2.15	0.57	14.31
54	4.26	1.51	2.20	2.07	0.97	13.76
56	3.63	1.55	2.20	1.69	0.52	12.94
58	3.22	1.15	1.95	1.60	0.46	11.48
60	2.95	0.98	1.75	1.42	0.66	10.93

(1 MB) 109.75 dB (eV⁻¹).

Note: MET refers to multiple electron transitions.

Reference: These data were taken from C.E. Brion and K.H. Tan, Chem.
Phys. 34, 141 (1978).

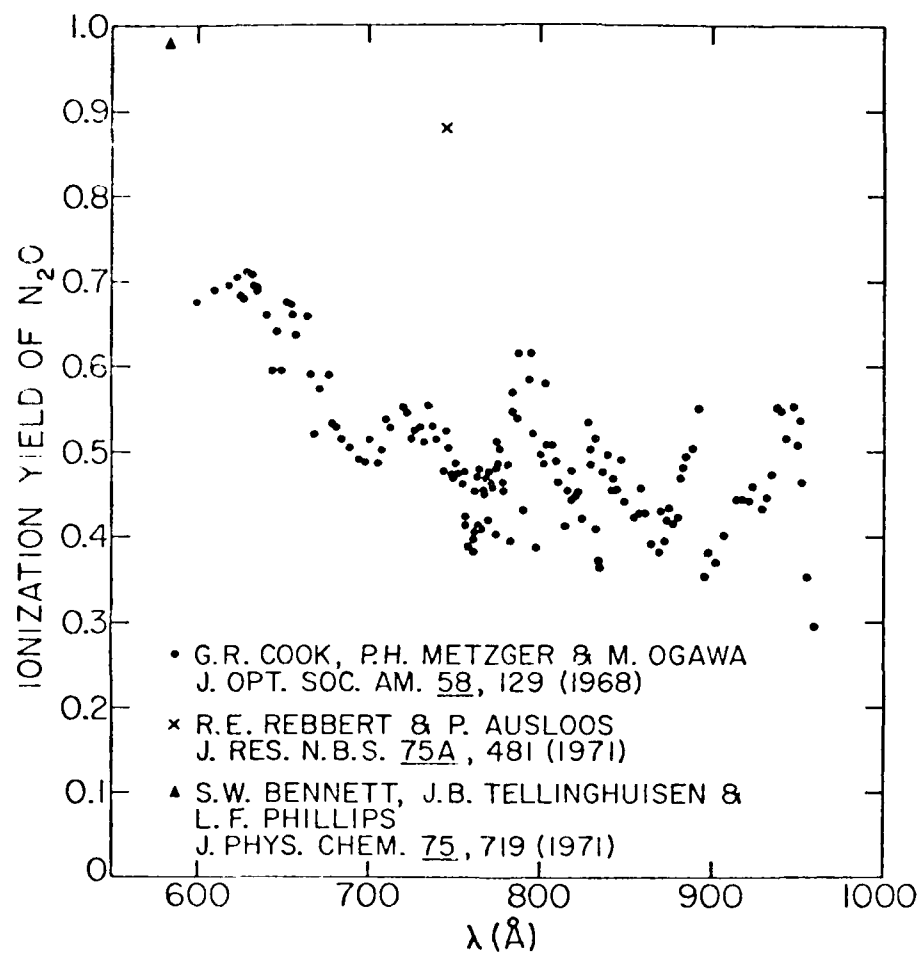
Tabular Data (continued)

Photoabsorption and vibrational orientation Oscillator strength, f_{osc} ; position, λ , in Å; 0

Wavelength, Å	Wavelength, Å	f_{osc}	f_{osc}	f_{osc}	f_{osc}	f_{osc}	f_{osc}
10	10.0	1.00	1.00	1.00	1.00	1.00	1.00
11	11.0	1.10	1.10	1.10	1.10	1.10	1.10
12	12.0	1.20	1.20	1.20	1.20	1.20	1.20
13	13.0	1.30	1.30	1.30	1.30	1.30	1.30
14	14.0	1.40	1.40	1.40	1.40	1.40	1.40
15	15.0	1.50	1.50	1.50	1.50	1.50	1.50
16	16.0	1.60	1.60	1.60	1.60	1.60	1.60
17	17.0	1.70	1.70	1.70	1.70	1.70	1.70
18	18.0	1.80	1.80	1.80	1.80	1.80	1.80
19	19.0	1.90	1.90	1.90	1.90	1.90	1.90
20	20.0	2.00	2.00	2.00	2.00	2.00	2.00
21	21.0	2.10	2.10	2.10	2.10	2.10	2.10
22	22.0	2.20	2.20	2.20	2.20	2.20	2.20
23	23.0	2.30	2.30	2.30	2.30	2.30	2.30
24	24.0	2.40	2.40	2.40	2.40	2.40	2.40
25	25.0	2.50	2.50	2.50	2.50	2.50	2.50
26	26.0	2.60	2.60	2.60	2.60	2.60	2.60
27	27.0	2.70	2.70	2.70	2.70	2.70	2.70
28	28.0	2.80	2.80	2.80	2.80	2.80	2.80
29	29.0	2.90	2.90	2.90	2.90	2.90	2.90
30	30.0	3.00	3.00	3.00	3.00	3.00	3.00
31	31.0	3.10	3.10	3.10	3.10	3.10	3.10
32	32.0	3.20	3.20	3.20	3.20	3.20	3.20
33	33.0	3.30	3.30	3.30	3.30	3.30	3.30
34	34.0	3.40	3.40	3.40	3.40	3.40	3.40
35	35.0	3.50	3.50	3.50	3.50	3.50	3.50
36	36.0	3.60	3.60	3.60	3.60	3.60	3.60
37	37.0	3.70	3.70	3.70	3.70	3.70	3.70
38	38.0	3.80	3.80	3.80	3.80	3.80	3.80
39	39.0	3.90	3.90	3.90	3.90	3.90	3.90
40	40.0	4.00	4.00	4.00	4.00	4.00	4.00
41	41.0	4.10	4.10	4.10	4.10	4.10	4.10
42	42.0	4.20	4.20	4.20	4.20	4.20	4.20
43	43.0	4.30	4.30	4.30	4.30	4.30	4.30
44	44.0	4.40	4.40	4.40	4.40	4.40	4.40
45	45.0	4.50	4.50	4.50	4.50	4.50	4.50
46	46.0	4.60	4.60	4.60	4.60	4.60	4.60
47	47.0	4.70	4.70	4.70	4.70	4.70	4.70
48	48.0	4.80	4.80	4.80	4.80	4.80	4.80
49	49.0	4.90	4.90	4.90	4.90	4.90	4.90
50	50.0	5.00	5.00	5.00	5.00	5.00	5.00
51	51.0	5.10	5.10	5.10	5.10	5.10	5.10
52	52.0	5.20	5.20	5.20	5.20	5.20	5.20
53	53.0	5.30	5.30	5.30	5.30	5.30	5.30
54	54.0	5.40	5.40	5.40	5.40	5.40	5.40
55	55.0	5.50	5.50	5.50	5.50	5.50	5.50
56	56.0	5.60	5.60	5.60	5.60	5.60	5.60
57	57.0	5.70	5.70	5.70	5.70	5.70	5.70
58	58.0	5.80	5.80	5.80	5.80	5.80	5.80
59	59.0	5.90	5.90	5.90	5.90	5.90	5.90
60	60.0	6.00	6.00	6.00	6.00	6.00	6.00
61	61.0	6.10	6.10	6.10	6.10	6.10	6.10
62	62.0	6.20	6.20	6.20	6.20	6.20	6.20
63	63.0	6.30	6.30	6.30	6.30	6.30	6.30
64	64.0	6.40	6.40	6.40	6.40	6.40	6.40
65	65.0	6.50	6.50	6.50	6.50	6.50	6.50
66	66.0	6.60	6.60	6.60	6.60	6.60	6.60
67	67.0	6.70	6.70	6.70	6.70	6.70	6.70
68	68.0	6.80	6.80	6.80	6.80	6.80	6.80
69	69.0	6.90	6.90	6.90	6.90	6.90	6.90
70	70.0	7.00	7.00	7.00	7.00	7.00	7.00
71	71.0	7.10	7.10	7.10	7.10	7.10	7.10
72	72.0	7.20	7.20	7.20	7.20	7.20	7.20
73	73.0	7.30	7.30	7.30	7.30	7.30	7.30
74	74.0	7.40	7.40	7.40	7.40	7.40	7.40

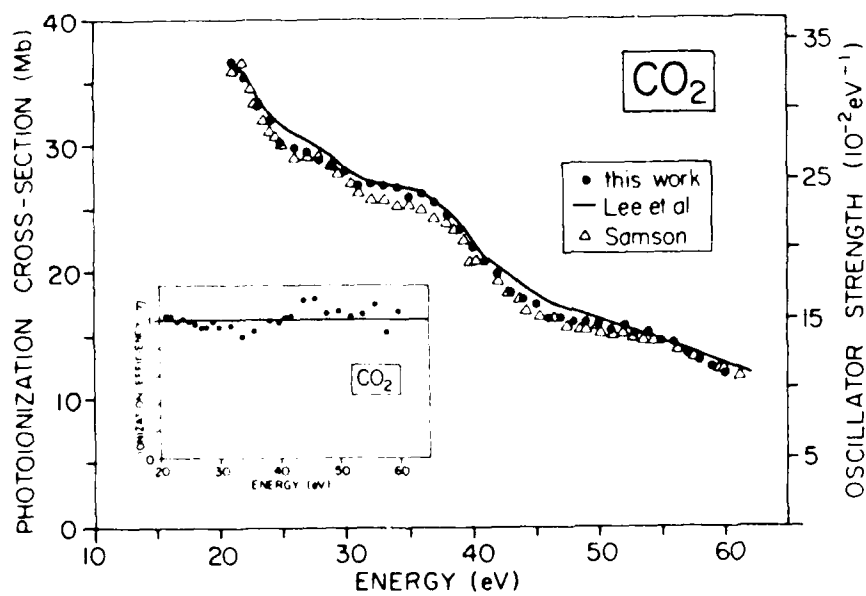
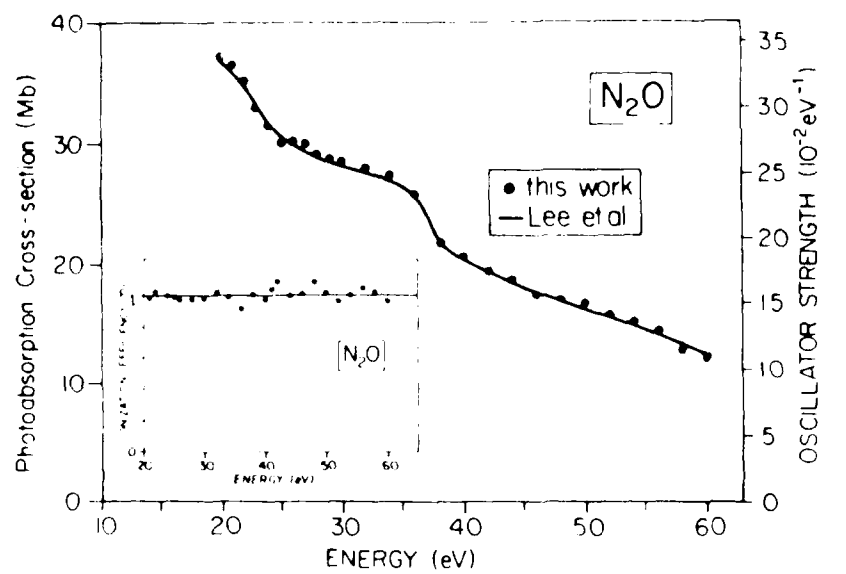
Continued on next page

Reference: These data were taken from A.P. Hitchcock, E.L. Prion, and M.L. Van der Wiel, Chem. Phys. 45, 461 (1960).



Graphical Data D-7.21

Quantum yield of ionization for N_2O .



Graphical Data D-2.22

Photoabsorption and Photoionization efficiency of N_2O and CO_2 .

Tabular Data D-2.23

Photoelectron Branching Ratios (%) for Various States in the
Photoionization of N_2O and CO_2

Photoelectron branching ratios (%) for electronic states of N_2O^+

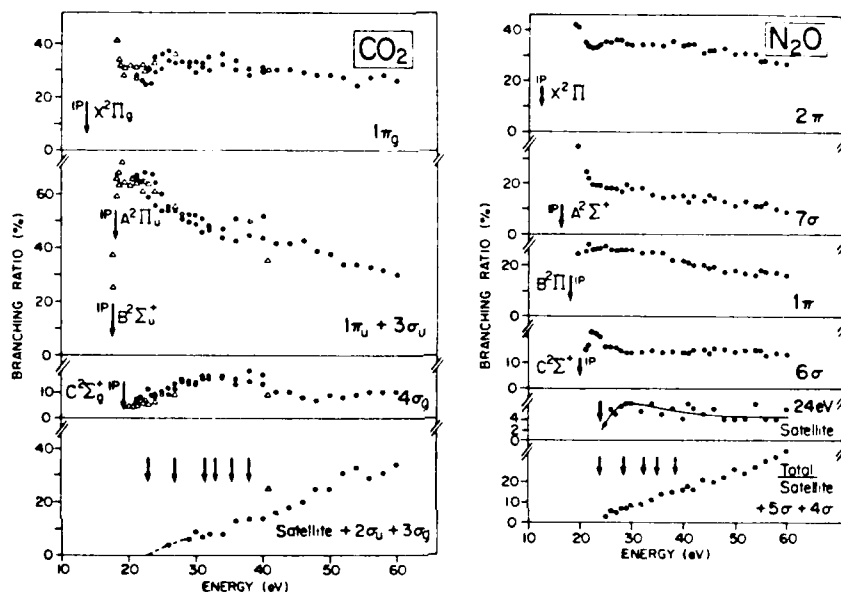
Energy (eV)	$X^2\Pi$	$A^2\Sigma^+$	$B^2\Pi$	$C^2\Sigma^+$	24 eV	Total MET
17.1	56	44	-	-	-	-
19.0	42	41	17	-	-	-
19.6	41	35	25	-	-	-
21.2	35	25	26	15	-	-
22	33	19	26	22	-	-
23	33	19	27	21	-	-
24	34	19	27	21	-	-
25	36	18	27	16	3	3
26	35	18	26	16	6	6
27	36	18	26	16	5	5
28	36	17	26	15	7	7
29	34	19	26	14	7	7
30	34	18	26	14	7	9
32	34	18	25	14	6	9
34	34	15	25	15	7	11
36	34	14	25	14	5	14
38	35	15	22	14	6	15
40	34	15	22	14	4	16
41	35	13	21	14	8	18
42	35	15	20	15	6	16
44	31	13	20	15	5	21
46	32	14	19	15	6	20
48	33	13	17	15	4	22
50	31	11	18	14	4	26
52	31	13	17	15	4	24
54	31	11	16	15	7	27
56	28	12	17	13	4	30
58	28	10	17	14	4	32
60	27	9	16	13	6	35

Photoelectron branching ratios (%) for electronic states of CO_2^+

Energy (eV)	$X^2\Pi_g$	$A^2\Pi_u + B^2\Sigma_u$	$C^2\Sigma_g$	Total MET
21.2	28	66	6	-
22	26	65	8	-
23	30	59	11	-
24	35	56	9	-
25	36	54	10	-
26	36	52	9	4
27	35	52	10	4
28	31	50	14	5
29	31	50	13	6
30	29	49	13	9
31	31	46	15	7
32	30	47	15	8
34	32	44	16	8
36	30	43	13	13
38	28	45	14	14
40	29	44	13	14
41	31	43	10	16
42	30	43	10	16
44	30	42	10	18
46	29	43	8	20
48	28	39	7	25
50	28	38	9	25
52	27	34	8	31
54	24	34	9	33
56	27	33	10	29
58	28	32	10	31
60	26	30	10	34

Note: MET refers to multiple electron transitions.

Reference: These data were taken from C.E. Brion and K.H. Tan, Chem. Phys. 34, 141 (1978).



Graphical Data D-2.24

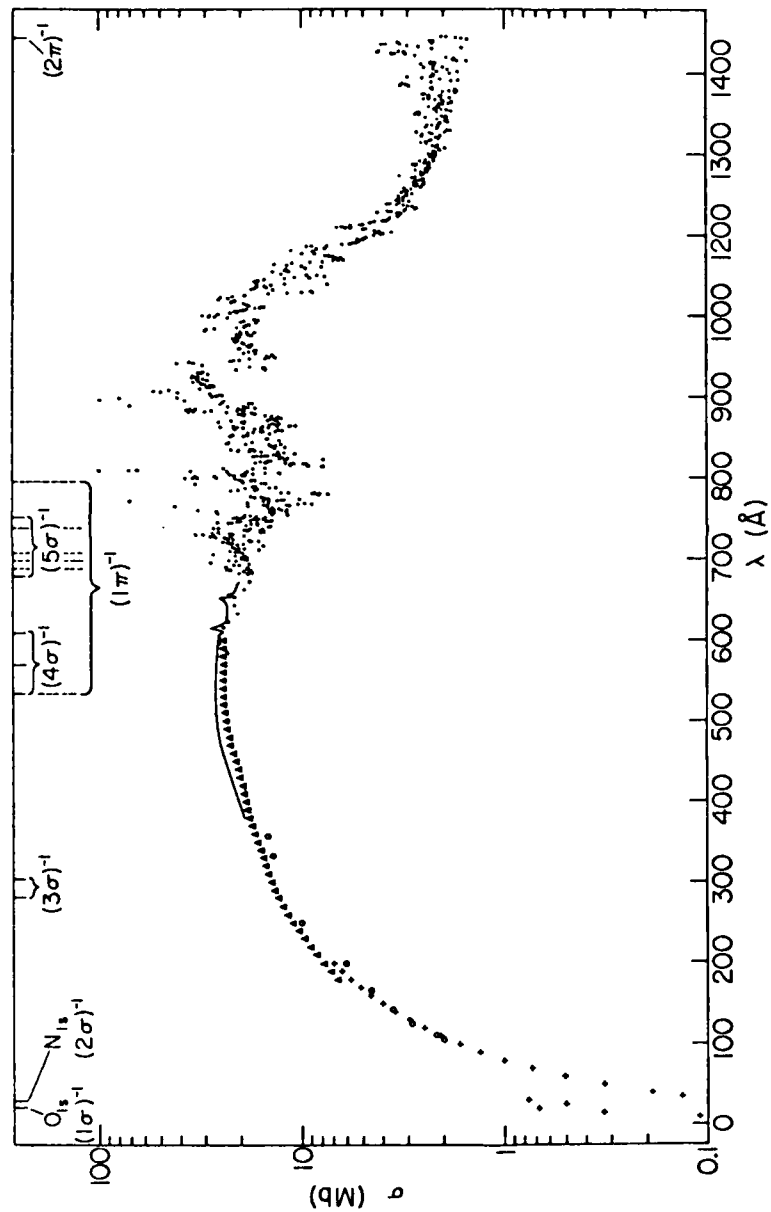
Photoionization branching ratios for CO_2 and N_2O .

Tabular Data D-2.25

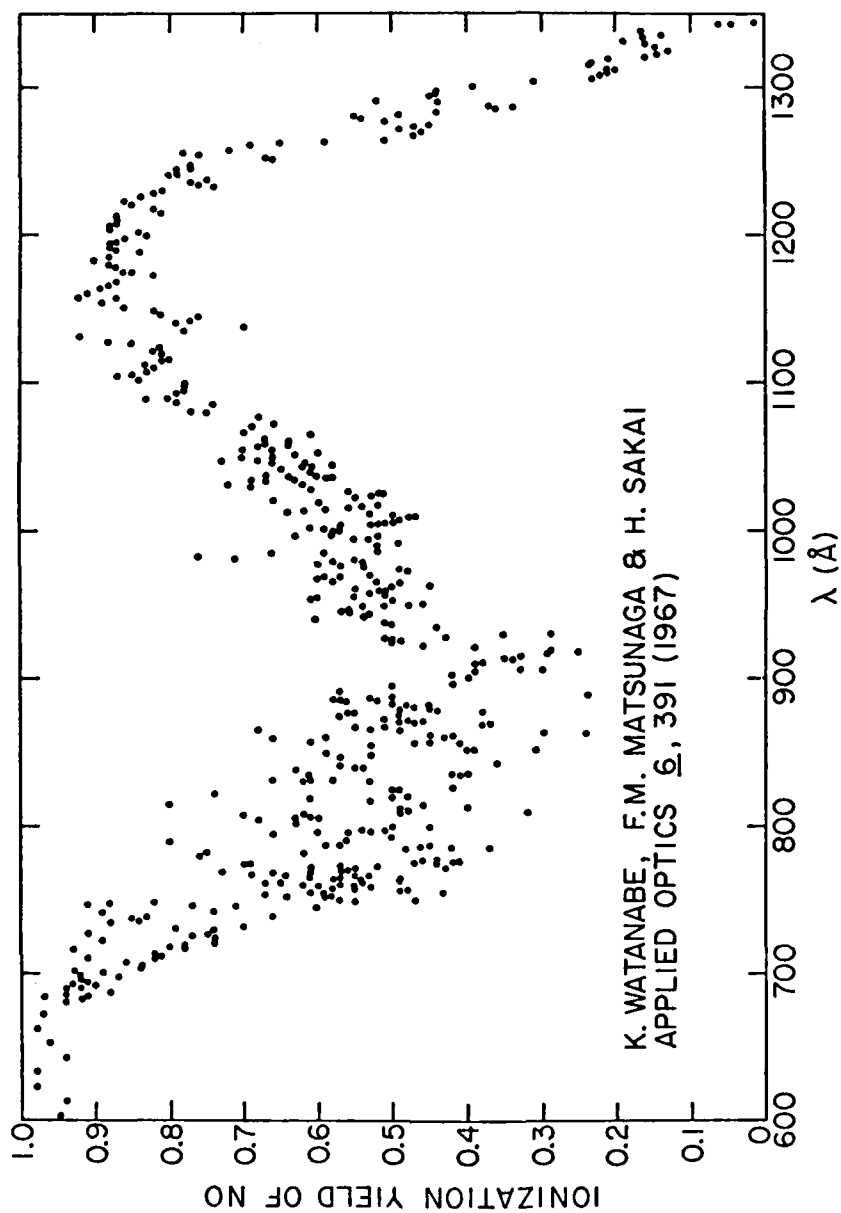
Photoabsorption Cross Section for NO
(Units of 10^{-18} cm^2)

$\lambda(\text{\AA})$	$\sigma(\text{Mb})$	$\lambda(\text{\AA})$	$\sigma(\text{Mb})$	$\lambda(\text{\AA})$	$\sigma(\text{Mb})$
672.5	18.9	600.0	24.2	110.0	2.08
668.0	19.2	575.0	24.2	100.0	1.68
660.0	19.8	550.0	24.4	90.0	1.33
657.5	20.1	525.0	24.2	80.0	1.01
656.5	21.2	500.0	23.7	70.0	0.74
654.0	20.3	475.0	23.0	60.0	0.51
653.5	21.0	450.0	21.2	50.0	0.33
651.0	23.4	425.0	20.0	40.0	0.19
650.0	22.8	400.0	18.8	35.0	0.136
646.5	21.9	375.0	17.5	30.0	0.77
640.0	21.4	350.0	15.8	25.0	0.50
630.0	21.3	325.0	14.5	20.0	0.68
625.0	21.5	300.0	13.2	15.0	0.33
622.0	21.9	275.0	12.0	9.89	0.109
619.0	22.8	250.0	10.5	8.265	0.067
617.0	23.7	225.0	8.8	6.199	0.030
616.0	24.7	200.0	6.98	4.153	0.0091
615.0	25.9	190.0	6.35	3.10	0.0038
614.0	25.6	180.0	5.75	2.48	0.0020
613.0	23.7	170.0	5.15	2.066	0.0011
611.0	22.8	160.0	4.57	1.550	0.0005
610.0	22.5	150.0	4.01	1.240	0.0002
607.5	22.8	140.0	3.49	0.827	0.0001
606.5	23.6	130.0	2.98	0.62	0.0000
603.5	23.2	120.0	2.51		

Reference: These data were taken from J. Berkowitz, Photoabsorption, Photoionization, and Photoelectron Spectroscopy (Academic Press, New York, 1979).



Graphical Data D-2.26
Photoabsorption cross section of NO.



Graphical Data D-2.27

Quantum yield of ionization for NO.

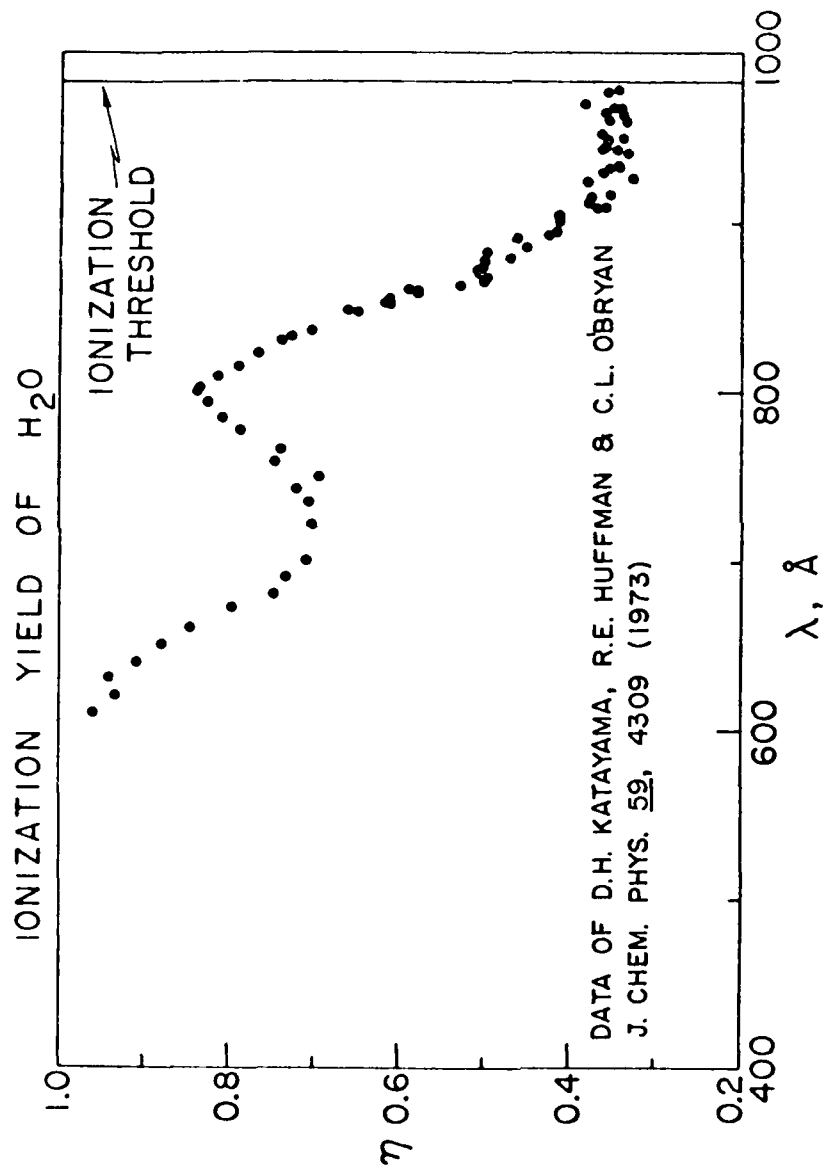
Table A Data D-2.123

Photoabsorption and Photofragmentation Oscillator Strengths (Cross Sections) for H₂O

Energy (eV)	Ionization threshold (eV)	Oscillator strength (10 ⁻¹⁸ cm ² /eV)			
		absorption	H ₂ O ⁺	OH ⁺	H ⁺
6		1.2			
7		2.77			
8		1.85			
9		5.32			
10		7.16	0		
11	0	8.40	0		
12	0.03	13.4	0.47		
13	0.23	17.1	4.00		
14	0.53	15.7	8.40		
15	0.75	15.1	11.30		
16	0.76	17.1	13.00	0	0
17	0.74	18.8	13.90	0	0
18	0.82	19.6	14.30	1.74	0.30
19	0.94	20.0	13.00	5.30	0.60
20	0.96	19.6	12.10	5.40	1.31
21	0.98	18.7	11.80	5.17	1.55
22	0.97	18.3	11.40	4.91	1.53
23	1.02	16.8	11.10	4.67	1.47
24	1.00	15.5	10.70	4.41	1.35
25	1.02	15.4	10.30	4.17	1.27
26	1.00	15.1	9.90	3.95	1.24
27	0.99	14.8	9.67	3.78	1.17
28	1.00	14.1	9.30	3.65	1.13
29	1.01	13.4	8.90	3.44	1.13
30	1.01	12.6	8.40	3.23	1.12
31	0.99	12.2	8.00	2.97	1.06
32	0.98	11.9	7.70	2.88	1.03
33	0.96	11.7	7.40	2.69	1.07
34	0.97	11.2	7.00	2.54	1.14
35	0.98	10.7	6.76	2.45	1.16
36	0.94	10.6	6.40	2.27	1.13
37	0.97	9.80	6.10	2.16	1.12
38	1.02	9.20	5.96	2.13	1.14
39	1.01	8.90	5.70	1.98	1.14
40	1.01	8.70	5.50	1.92	1.16
41	1.00	8.40	5.30	1.84	1.16
42	0.98	8.20	5.00	1.73	1.11
43	0.95	8.10	4.80	1.65	1.07
44	0.99	7.60	4.70	1.59	1.08
45	0.97	7.50	4.50	1.55	1.05
46	0.98	7.10	4.30	1.44	1.01
47	0.97	7.10	4.20	1.44	1.04
48	0.95	6.90	4.00	1.37	1.02
49	0.96	6.70	3.90	1.33	1.00
50	0.92	6.70	3.70	1.26	0.99
51	0.96	6.3	3.60	1.20	1.00
52	0.94	6.27	3.50	1.17	1.04
53	1.02	5.67	3.40	1.13	1.03
54	1.01	5.45	3.20	1.00	1.00
55	0.99	5.33	3.10	1.03	0.99
56	1.00	5.15	3.00	1.00	0.95
57	0.96	5.20	2.90	0.96	0.93
58	0.95	5.11	2.80	0.93	0.94
59	0.93	5.13	2.74	0.91	0.93
60	0.94	4.96	2.67	0.89	0.89

^a 10^{-18} cm² $\times 109.75$ eV / (eV²)

Reference: These data were taken from F. Zan, C.E. Brion, Ph. J. Van der Leeuw, and M.J. Van der Wiel, Chem. Phys., 29, 299 (1976).



Graphical Data D-2.29

Quantum yield of ionization for H_2O .

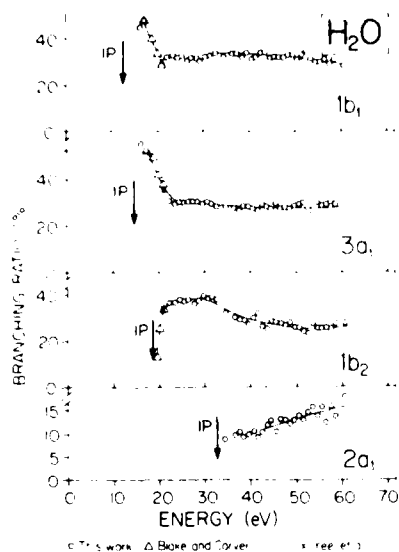
Tabular Data D-2.30

Partial Photoionization Oscillator Strengths
(Cross Sections) for the Valence Orbitals of
 H_2O

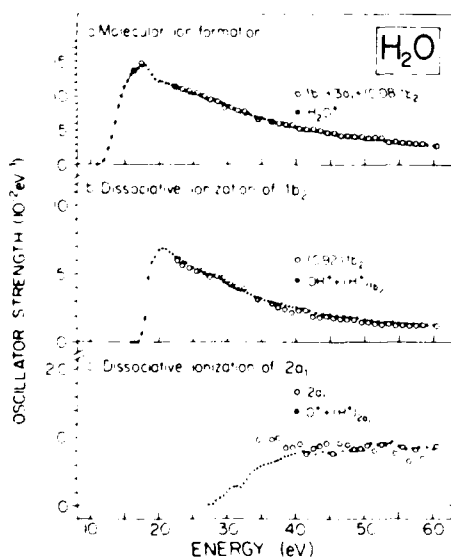
Energy (eV)	Oscillator strength (10^{-2} eV^{-1}) ^{a)}			
	1b ₁	3a ₁	1b ₂	2a ₁
16.5	6.06	7.40		
17.5	7.13	7.73	6.44	
22.5	5.67	5.29		
23.5	5.50	4.96	6.05	
24.5	5.15	4.83	6.01	
25.5	5.07	4.65	5.64	
26.5	4.43	4.60	5.63	
27.5	4.61	4.56	5.23	
28.5	4.37	4.35	5.19	
29.5	3.94	3.73	5.22	
30.5	3.98	3.79	4.73	
31.5	3.89	3.41	4.60	
32.5	3.92	3.31	4.34	
34.5	3.26	2.96	3.43	0.96
36.5	3.37	2.70	3.00	0.96
37.5	3.13	2.59	2.73	1.02
38.5	2.91	2.60	2.72	0.86
39.5	3.05	2.55	2.44	0.86
40.5	2.78	2.32	2.58	0.92
41.5	2.68	2.15	2.61	0.77
42.5	2.71	2.30	2.09	0.85
43.5	2.61	2.18	1.97	0.90
44.5	2.35	2.04	2.12	0.93
45.5	2.27	2.09	1.98	0.75
46.5	2.18	1.87	1.88	0.91
47.5	2.12	1.86	1.78	0.88
48.5	2.14	1.76	1.81	0.81
49.5	2.08	1.80	1.61	0.82
50.5	1.91	1.80	1.62	0.87
51.5	2.01	1.82	1.46	0.79
52.5	2.11	1.66	1.42	0.89
53.5	1.72	1.34	1.50	0.87
54.5	1.62	1.62	1.37	0.76
55.5	1.57	1.49	1.34	0.84
56.5	1.60	1.47	1.32	0.64
57.5	1.45	1.49	1.24	0.80
58.5	1.51	1.40	1.31	0.68
60.5	1.32	1.22	1.31	0.87

a) $\sigma(10^{-18} \text{ cm}^2) = 109.75 \text{ df/dE (eV}^{-1}\text{)}$.

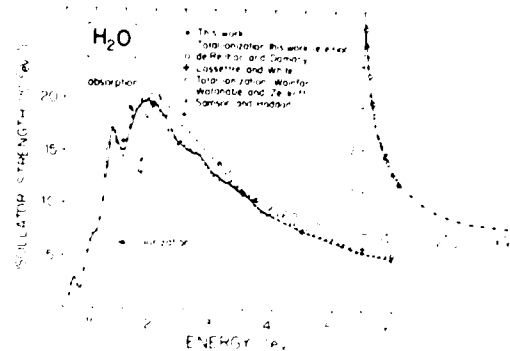
Reference: These data were taken from K. Tan, C.E. Brion, Ph. E. Van der Leeuw, and M.J. Van der Wiel, Chem. Phys. 29, 299 (1978).



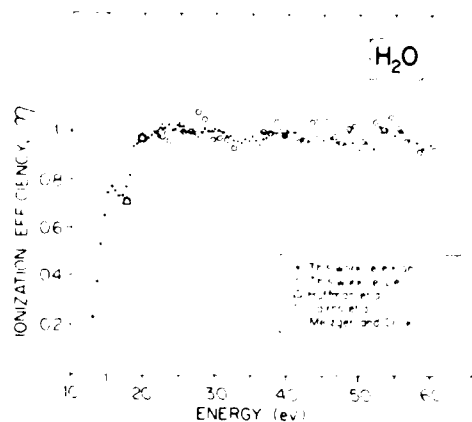
Photoionisation branching ratios for formation of electronic states of H_2O^+ .



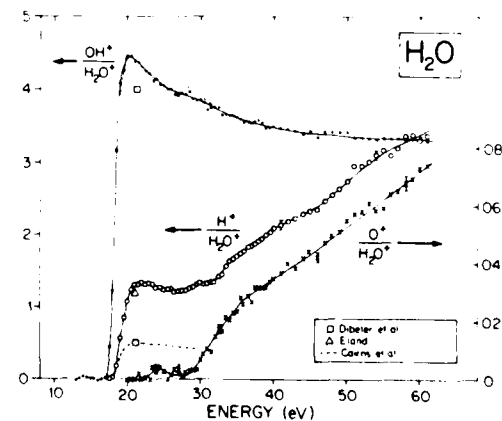
Partition of total oscillator strength for ionisation over molecular and fragment ion formation. (a) Contributions to H_2O^+ . (b) Contributions to OH^+ and H^+ from the $1b_2$ state. (c) Contributions to H^+ and O^+ from the $2a_1$ state.



Photoabsorption oscillator strength of H_2O .



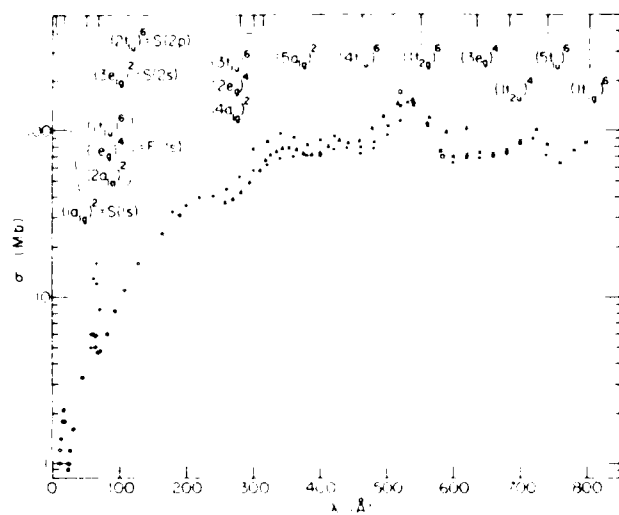
Photoionisation efficiency of H_2O .



Fractional abundances of the fragment ions formed from photoionisation of H_2O .

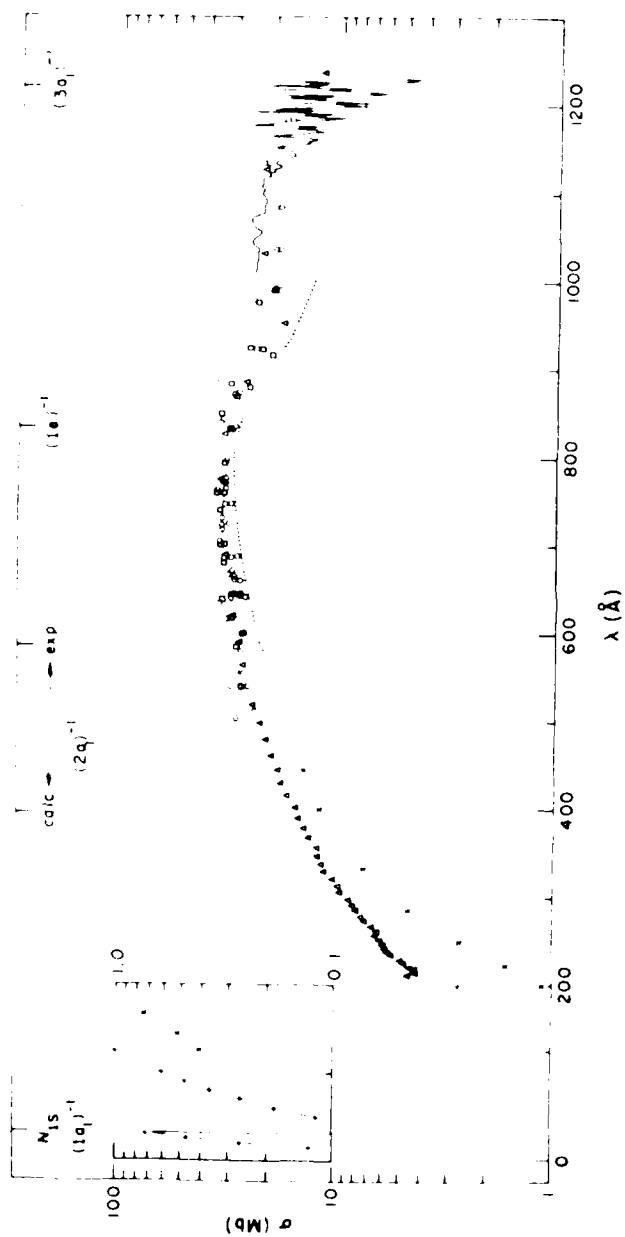
Graphical Data D-2.31 Spectroscopy Data for H_2O .

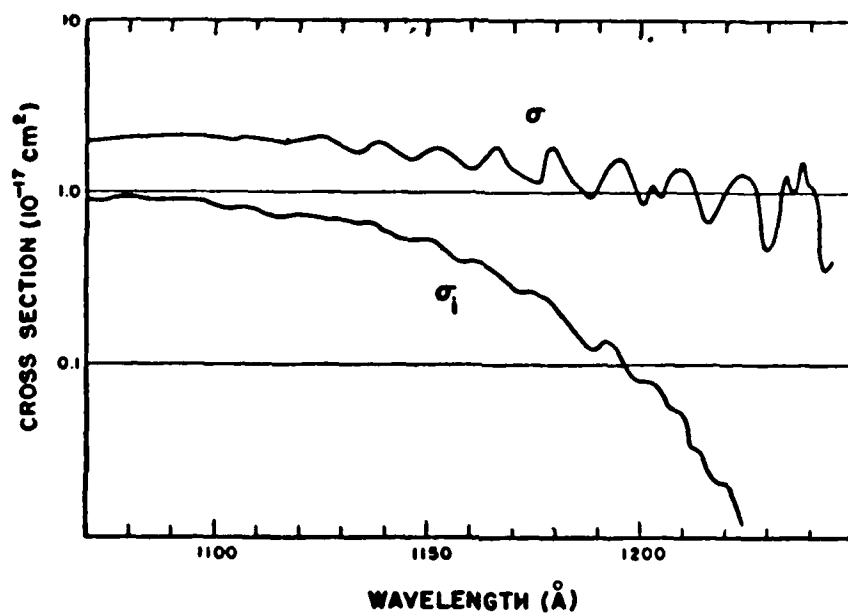
Reference: These data were taken from K. Tan, C.E. Brion, Ph. E. Van der Leeuw, and M.J. Van der Wiel, Chem. Phys. **29**, 229 (1978).



Graphical Data D-2.32

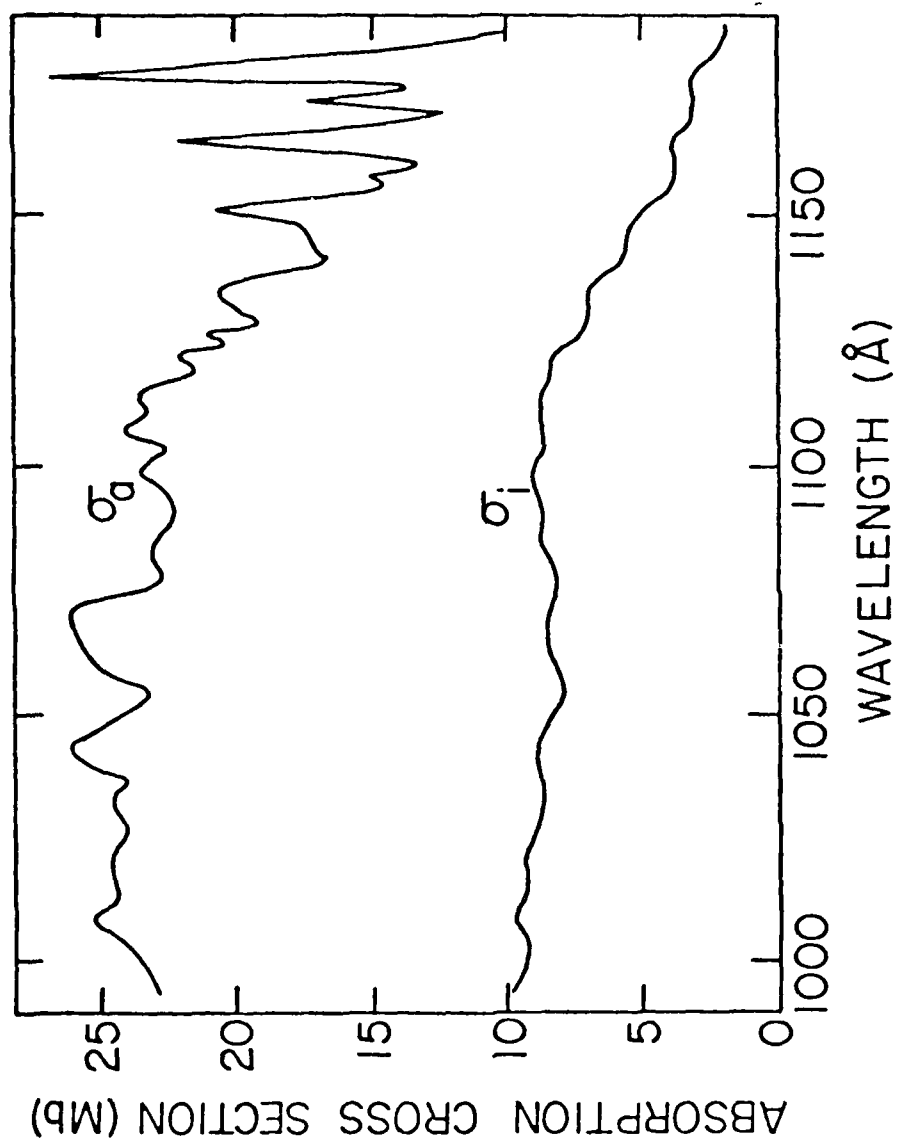
Photoabsorption cross section of SF_6 .





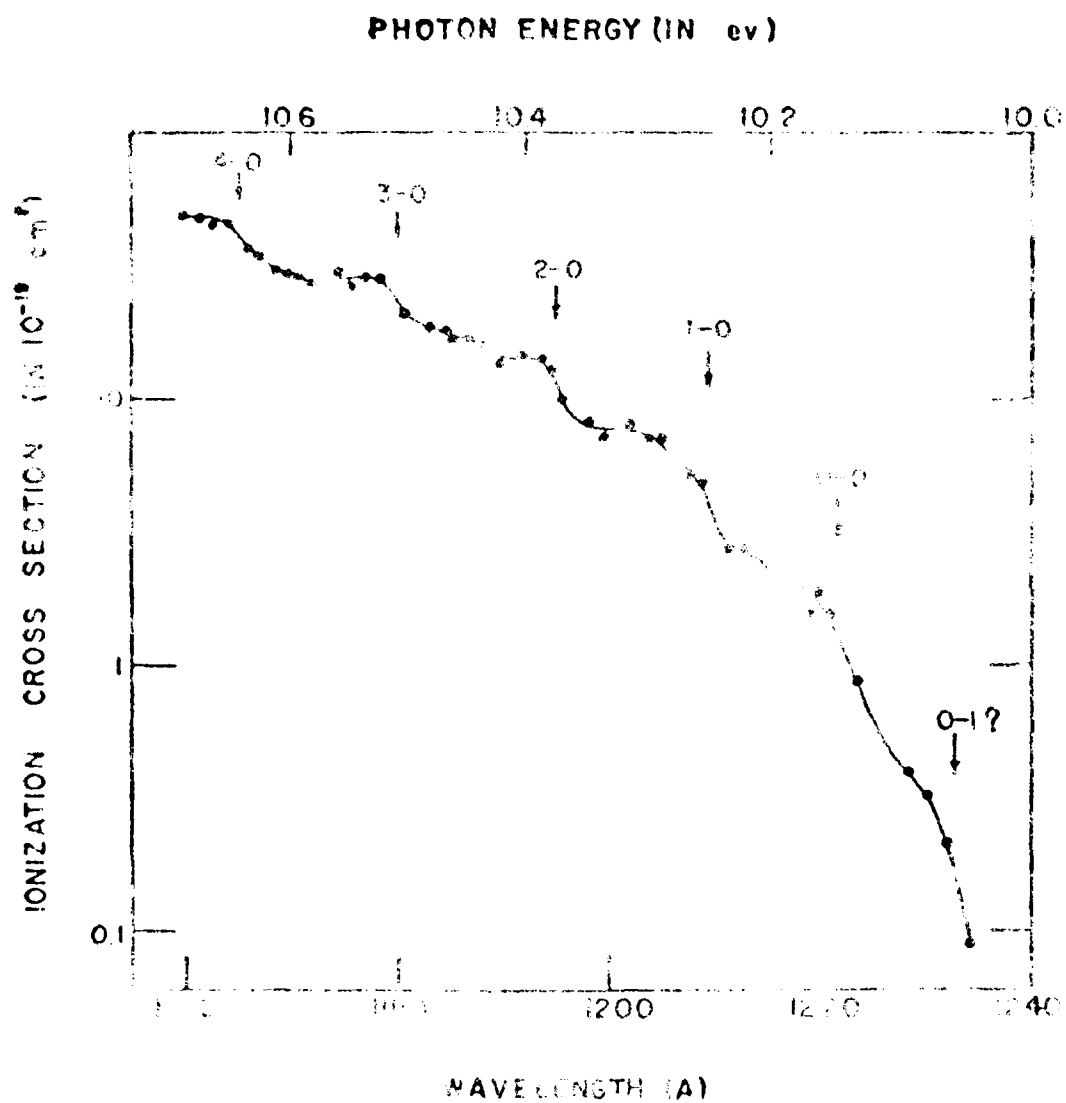
Graphical Data D-2.34

Photoabsorption (σ) and photoionization (σ_i) cross sections for NH_3 .



Graphical Data D-2.35

Photoabsorption (σ_a) and photoionization (σ_i) cross sections for NH_3 .



Continued from page 10

For a complete description of the data, see page 10

Tabular Data D-2.37

Photoabsorption and Partial Photoionization Oscillator Strengths
(Cross Sections) for NH_3

Energy(eV)	Oscillator Strengths (eV^{-1}) ^c			Absorption df_0/dE^a
	$3a_1$	$1e$	$2a_1$	
15	0.065 ^b	—	—	0.303
16	0.065	—	—	0.315
17	0.063	—	—	0.312
18	0.062	0.193	—	0.298
19	0.057	0.234	—	0.291
20	0.055	0.217	—	0.272
21	0.052	0.209	—	0.261
22	0.049	0.196	—	0.245
23	0.050	0.183	—	0.233
24	0.048	0.170	—	0.220
25	0.048	0.154	—	0.202
27	0.047	0.133	—	0.180
30	0.041	0.111	—	0.152
32	0.036	0.093	0.006	0.135
34	0.029	0.079	0.013	0.121
36	0.026	0.070	0.013	0.109
38	0.026	0.065	0.011	0.102
40	0.021	0.056	0.010	0.087
42	0.021	0.047	0.011	0.079
44	0.017	0.044	0.011	0.072
46	0.017	0.039	0.010	0.066
48	0.017	0.032	0.010	0.058
50	0.017	0.028	0.011	0.056

^a Total absorption oscillator strength

^b df_0/dE for $3a_1$ can still be computed below 19 eV where $\eta_i < 1$ from the raw data using df_0/dE and the ionization data where $\eta_i = 1$.

^c Partial ionization cross-sections (Mb) may be obtained by multiplying by a factor 109.75.

Reference: These data are from C.E. Brion, A. Hamnett, G.R. Wight, and M.J. Van der Wiel, J. Electron Spectrosc. 12, 323 (1977).

Tabular Data D-2.38

Photoabsorption and Photoionization Oscillator Strengths
(Cross Sections) for NH_3

Energy (eV)	$f^{(0)}$ (eV^{-1})		Fragment NH_3^+ (%)			
	Absorption	NH_3	NH_2^+	NH^+	N^+	H^+
10	0.110					
11	0.204	0.060				
12	0.204	0.087				
13	0.162	0.083				
14	0.240	0.091				
15	0.303	0.121	0			
16	0.315	0.137	24			
17	0.312	0.114	100			
18	0.298	0.107	158			
19	0.291	0.106	170			
20	0.272	0.099	173	0.2		
21	0.261	0.096	168	0.2		
22	0.245	0.090	164	0.2		
23	0.233	0.085	164	0.2		0.5
24	0.220	0.081	163	0.3		0.8
25	0.202	0.075	167	1.2		1.2
26	0.192	0.074	161	2.1		1.5
27	0.180	0.068	154	4.4		3.0
28	0.166	0.065	156	5.5	0.2	5.2
29	0.162	0.059	158	6.8	0.3	6.9
30	0.152	0.057	153	7.7	0.5	9.4
31	0.141	0.054	154	8.5	0.7	9.7
32	0.135	0.050	148	8.8	1.0	11.6
33	0.127	0.048	146	9.2	1.2	12.6
34	0.121	0.045	145	9.8	1.7	14.4
35	0.110	0.042	146	11.1	2.2	17.1
36	0.109	0.039	141	10.6	1.9	17.5
37	0.105	0.038	142	11.8	2.1	17.9
38	0.102	0.036	141	12.0	2.2	19.4
39	0.092	0.033	141	12.8	2.0	20.5
40	0.087	0.032	139	13.3	2.5	24.7
41	0.086	0.030	141	14.0	3.1	23.5
42	0.079	0.029	138	14.4	2.7	25.8
43	0.075	0.027	137	14.9	2.8	27.3
44	0.072	0.026	136	15.4	3.0	28.8
45	0.069	0.024	139	16.7	3.2	32.4
46	0.066	0.023	136	16.3	3.4	30.7
47	0.062	0.022	137	17.1	3.4	31.4
48	0.058	0.021	135	17.7	3.3	33.3
49	0.059	0.020	136	19.4	3.8	34.2
50	0.056	0.019	135	19.2	3.8	40.0
51	0.054	0.018	133	19.8	4.4	37.8
52	0.053	0.017	133	19.3	4.0	38.2
53	0.051	0.017	134	21.3	4.3	41.0
54	0.049	0.016	132	21.4	4.2	38.9
55	0.045	0.015	136	23.1	4.5	46.6
56	0.044	0.015	132	21.8	5.0	43.0
57	0.041	0.014	130	24.1	5.3	44.9
58	0.039	0.013	129	22.5	5.5	46.0
59	0.039	0.013	131	25.0	5.6	49.4
60	0.042	0.012	129	23.2	5.7	49.2

Note: $\sigma(10^{-18} \text{ cm}^2) = 109.75 \times f^0 (\text{eV}^{-1})$

Reference: These data are from G.R. Wight, M.J. Van der Wiel, and C.E. Brion, J. Phys. B 10, 1863 (1977).

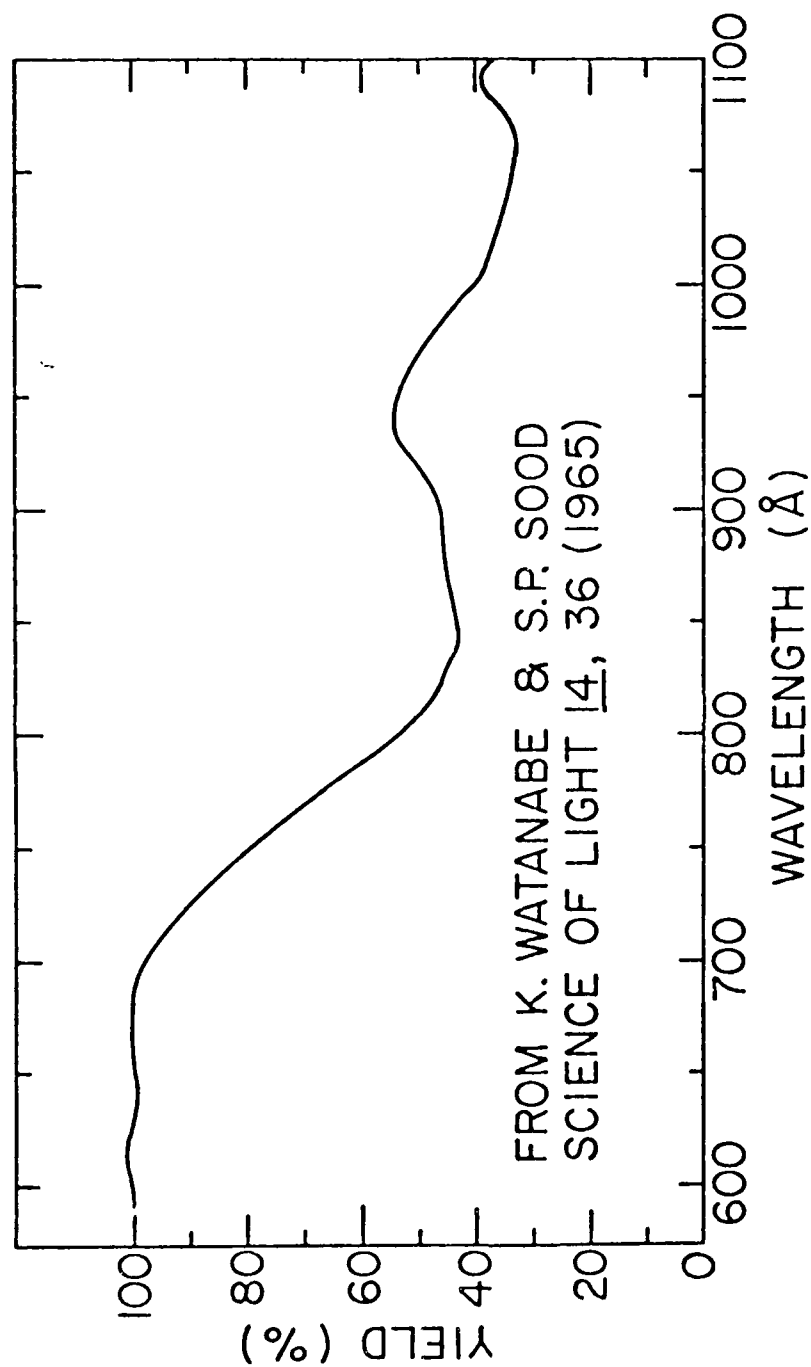
Tabular Data D-2.39

Photoelectron Branching Ratios () for Various
States in the Photoionization of NF_3

<i>Energy (eV)</i>	<i>3a₁</i>	<i>1e</i>	<i>2a₁</i>
18	24.3	75.7	—
19	19.5	80.5	—
20	20.1	79.9	—
21	20.1	79.9	—
22	20.2	79.8	—
23	21.3	78.7	—
24	22.6	77.4	—
25	23.6	76.4	^a
27	26.3	73.7	^a
30	27.2	72.8	^a
32	26.4	69.2	4.4
34	23.6	65.3	11.0
36	24.0	64.2	11.7
38	25.4	64.2	10.4
40	24.1	64.1	11.7
42	27.2	59.0	13.8
44	23.4	61.4	15.2
46	25.6	59.0	15.5
48	28.8	54.4	16.7
50	29.6	50.8	19.6

^a Not measured.

Reference: These data are from C.E. Brion, A. Hamnett, G.P. Wight, and M.J. Van der Wiel, J. Electron Spectrosc. 12, 323 (1977).



Graphical Data D-2.40

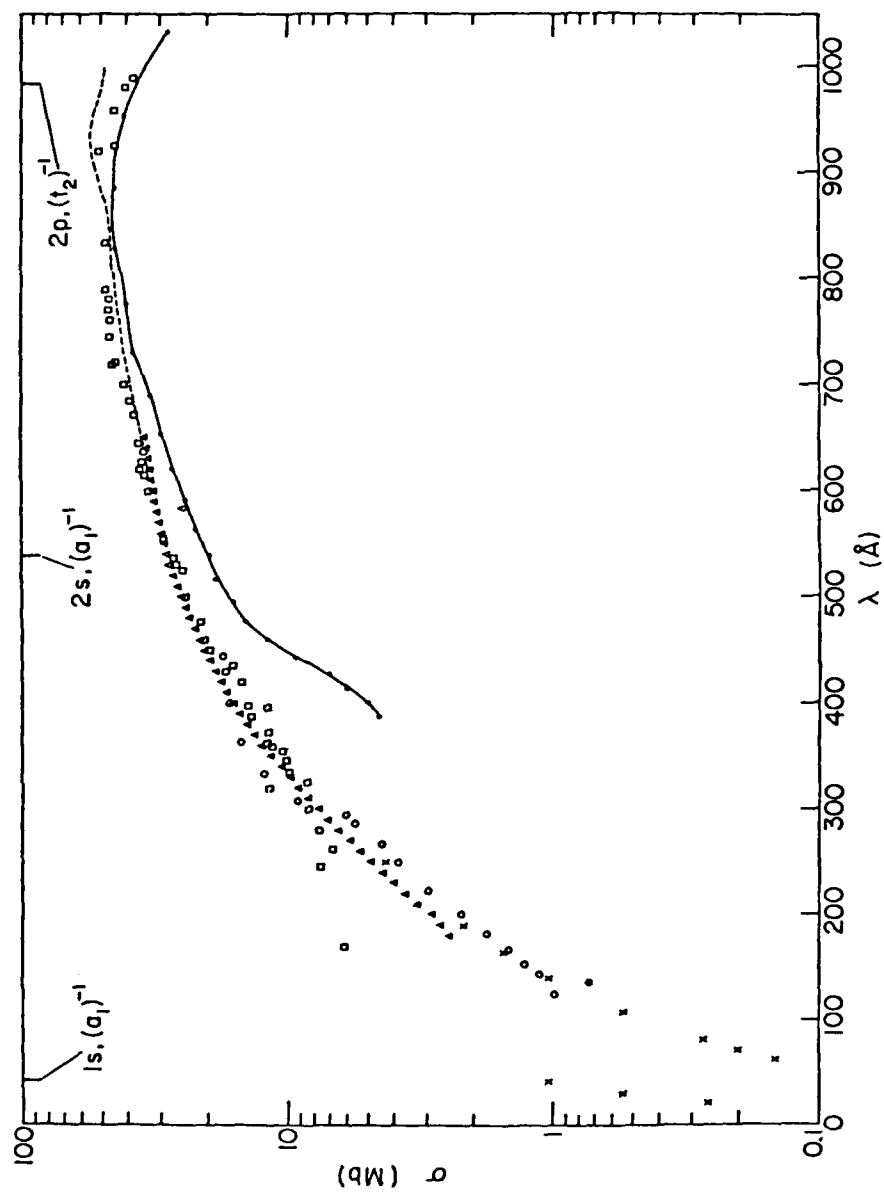
Quantum yield of ionization for NH_3 .

Tabular Data D-2.41

Photoabsorption Cross Section for CH₄

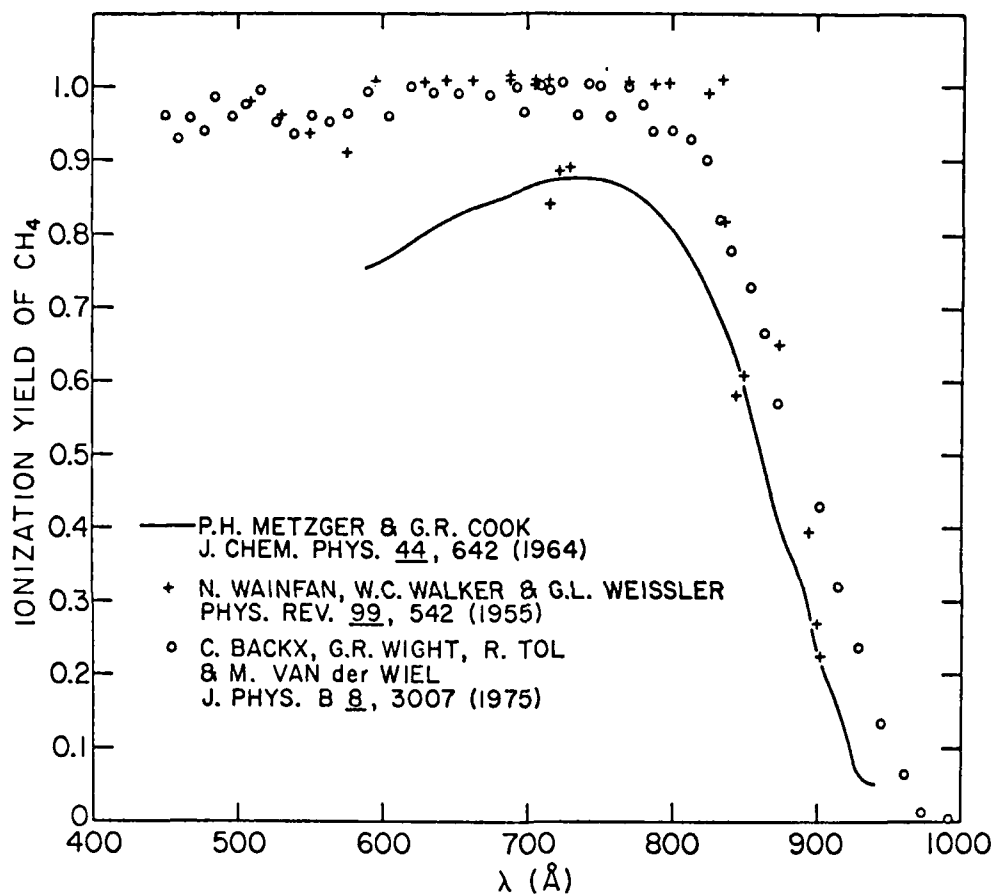
$\lambda(\text{\AA})$	$\sigma(\text{Mb})$	$\lambda(\text{\AA})$	$\sigma(\text{Mb})$
23.6	0.263	425.0	16.0
31.4	0.550	450.0	18.6
40.96	1.06	475.0	21.0
44.4	0.0617	500.0	23.2
64.35	0.148	525.0	25.2
72.20	0.205	550.0	28.0
81.98	0.276	575.0	30.3
108.65	0.550	600.0	33.0
139.50	1.049	625.0	34.5
164.60	1.532	650.0	36.0
190.3	2.19	675.0	38.0
250.5	4.31	700.0	40.0
		725.0	44.0
275.0	5.35	750.0	46.0
300.0	6.9	775.0	46.5
325.0	8.5	800.0	48.0
350.0	10.2	825.0	48.0
375.0	12.0	850.0	49.0
400.0	14.0	875.0	51.0
		900.0	51.0
		925.0	51.0
		950.0	47.5
		975.0	41.5
		985.0	39.0

Reference: These data were taken from J. Berkowitz, Photoabsorption, Photoionization, and Photoelectron Spectroscopy (Academic Press, New York, 1979).



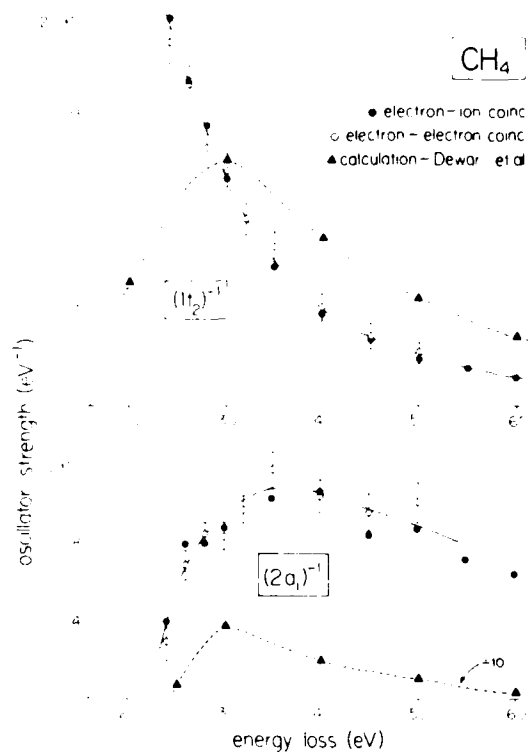
Graphical Data D-2.42

Photoabsorption cross section of CH_4 .



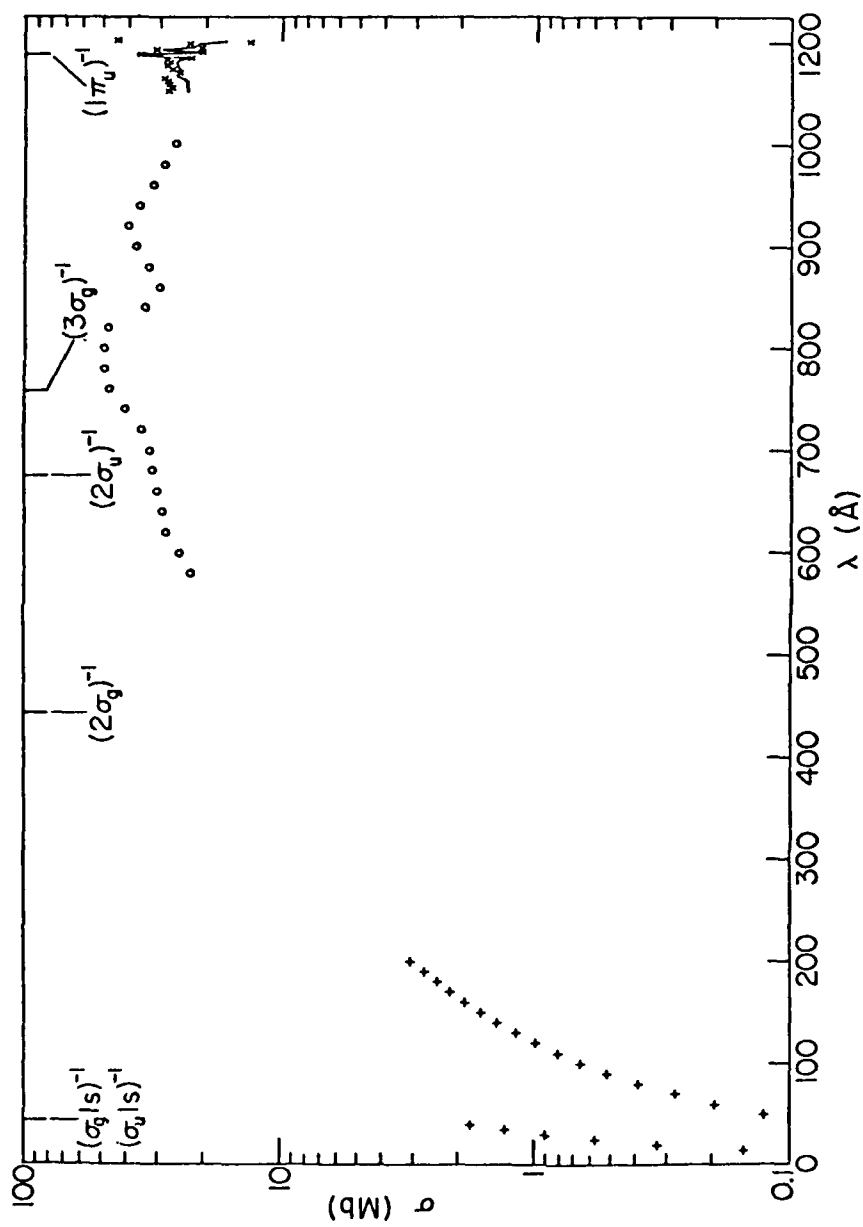
Graphical Data D-2.43

Quantum yield of ionization for CH₄.



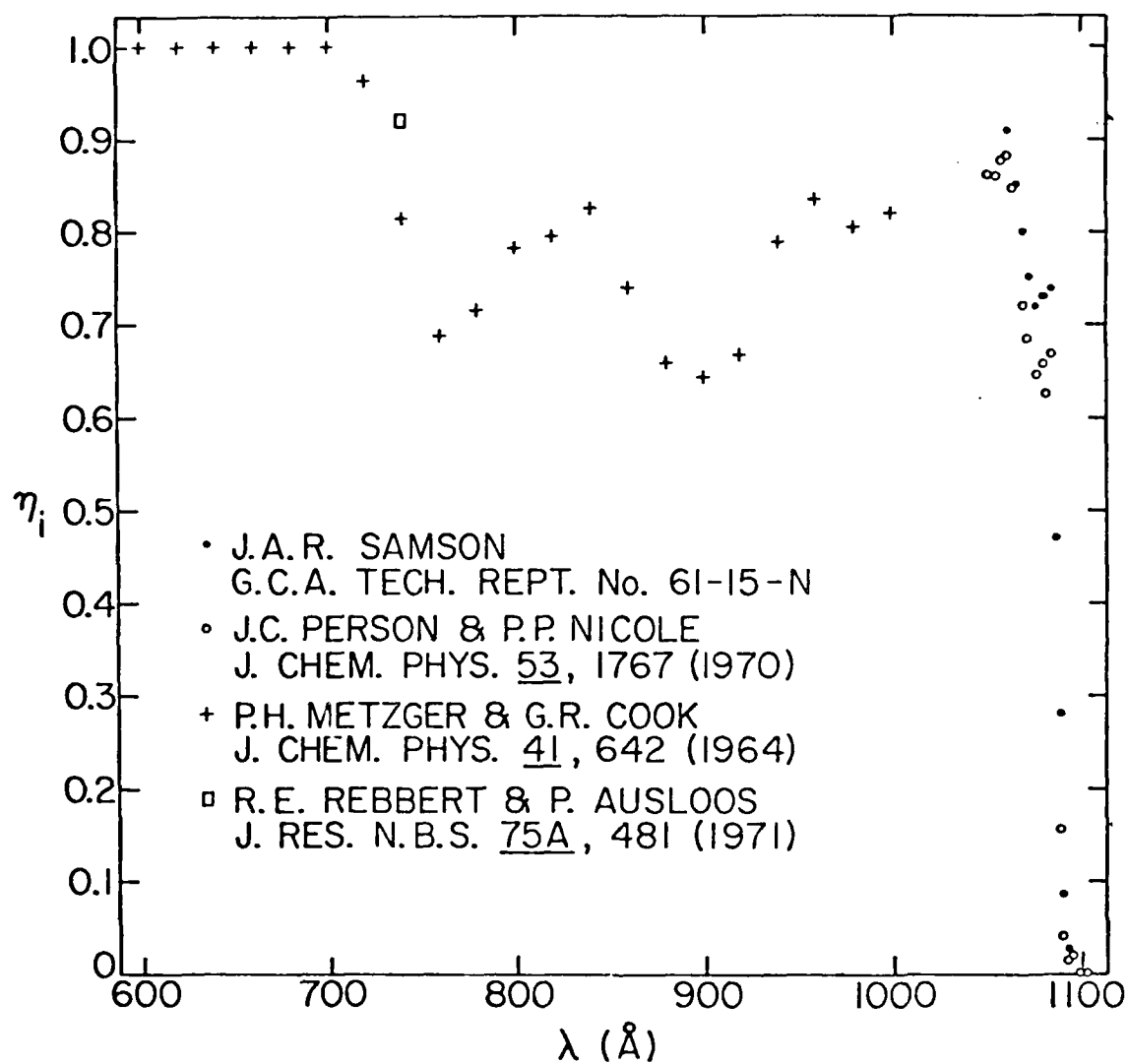
Graphical Data D-2.44

Partial photoionization oscillator strengths (cross sections) for production of the $(1t_2)^{-1}$ and $(2a_1)^{-1}$ states of CH_4^+ in the photoionization of CH_4 .



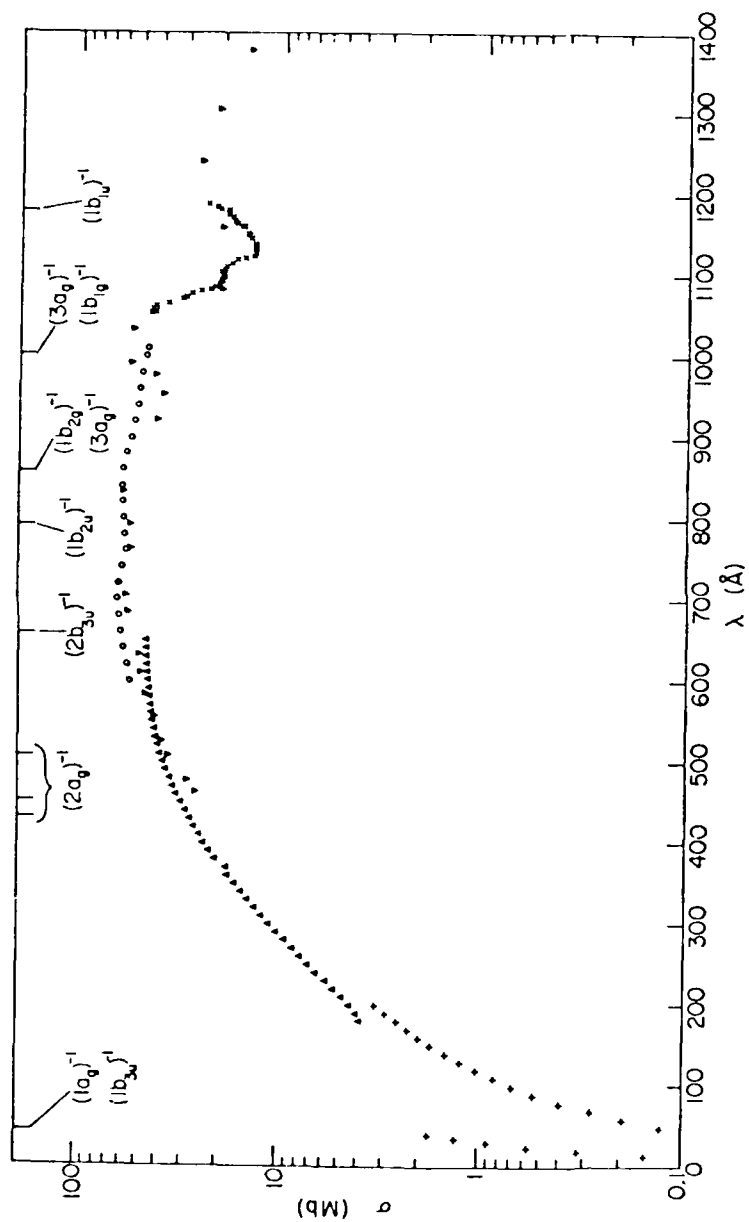
Graphical Data D-2.45

Photoabsorption cross section of C_2H_2 .



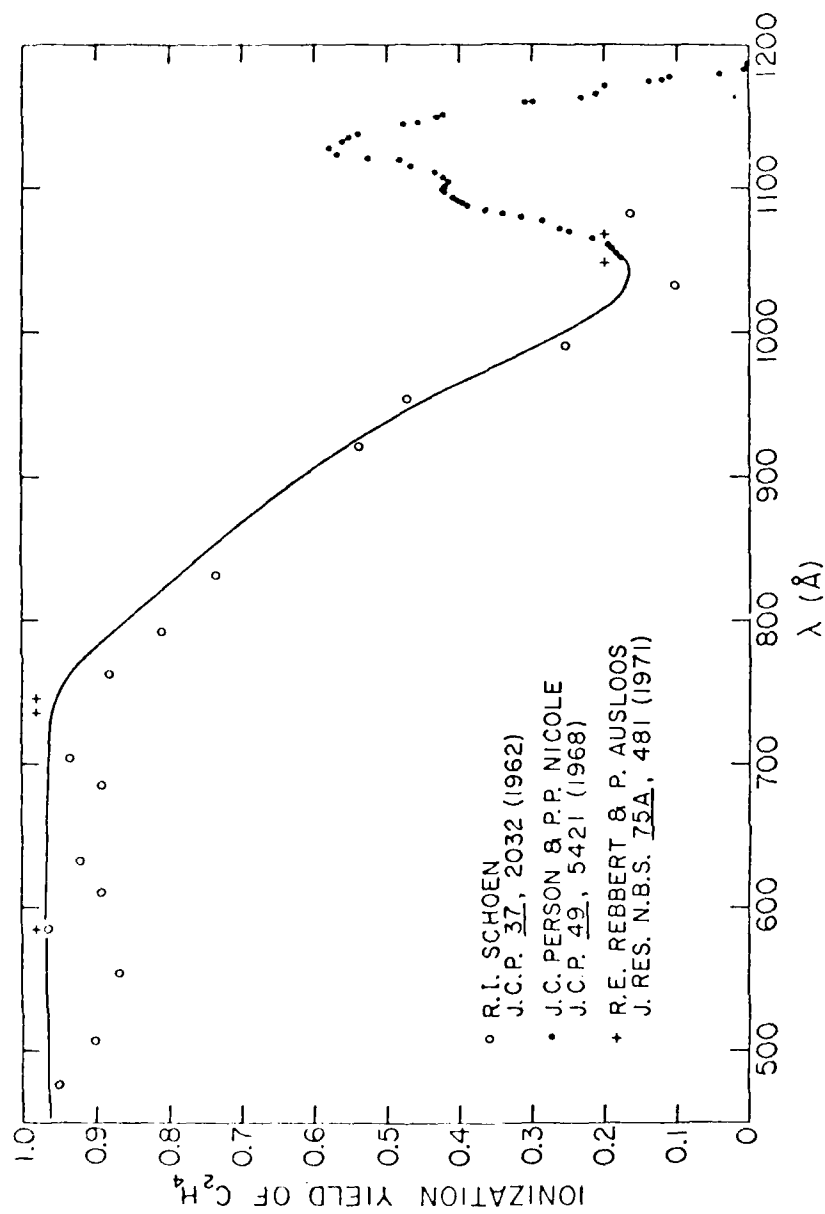
Graphical Data D-2.45

Quantum yield of ionization, η_i , for C_6H_6 .



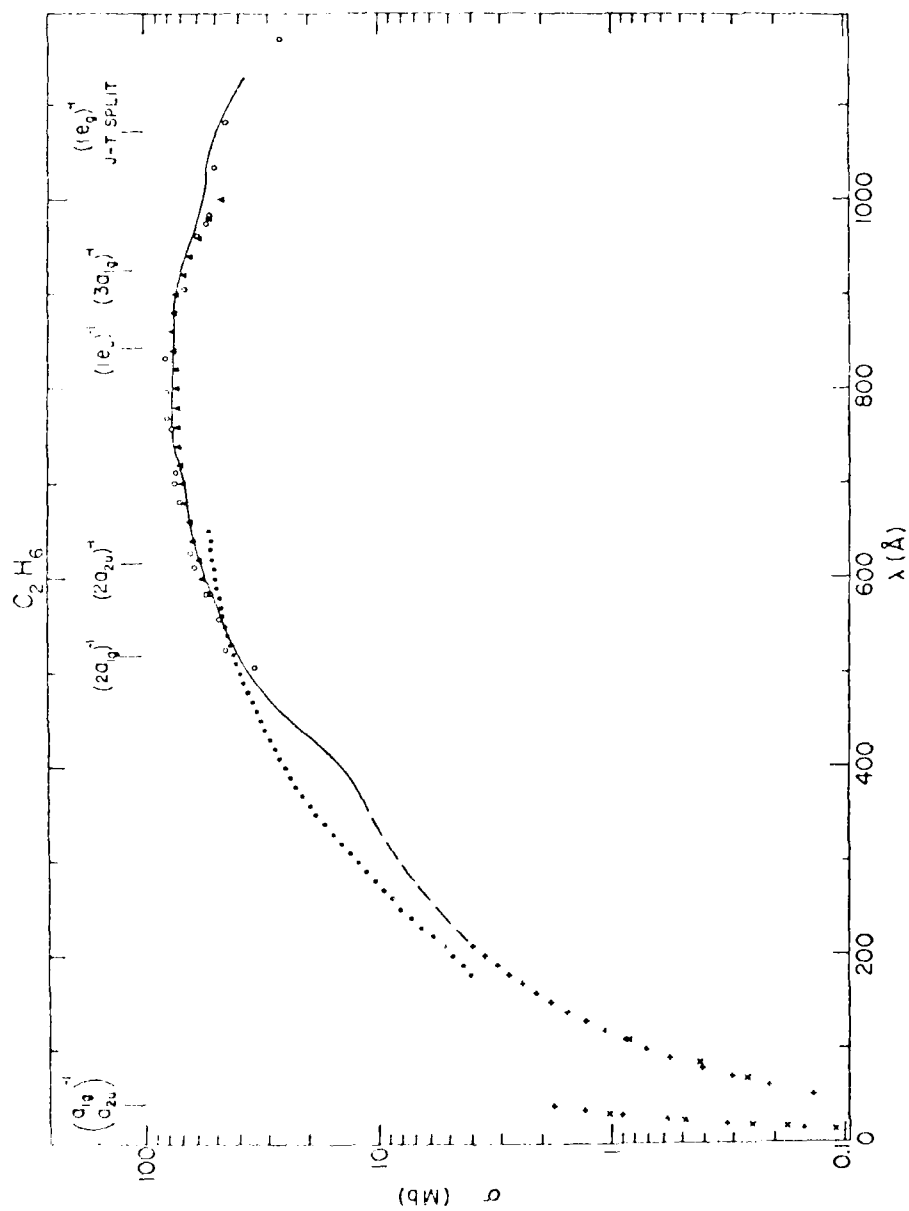
Graphical Data D-2.7

Photabsorption cross section of C_2H_4 .



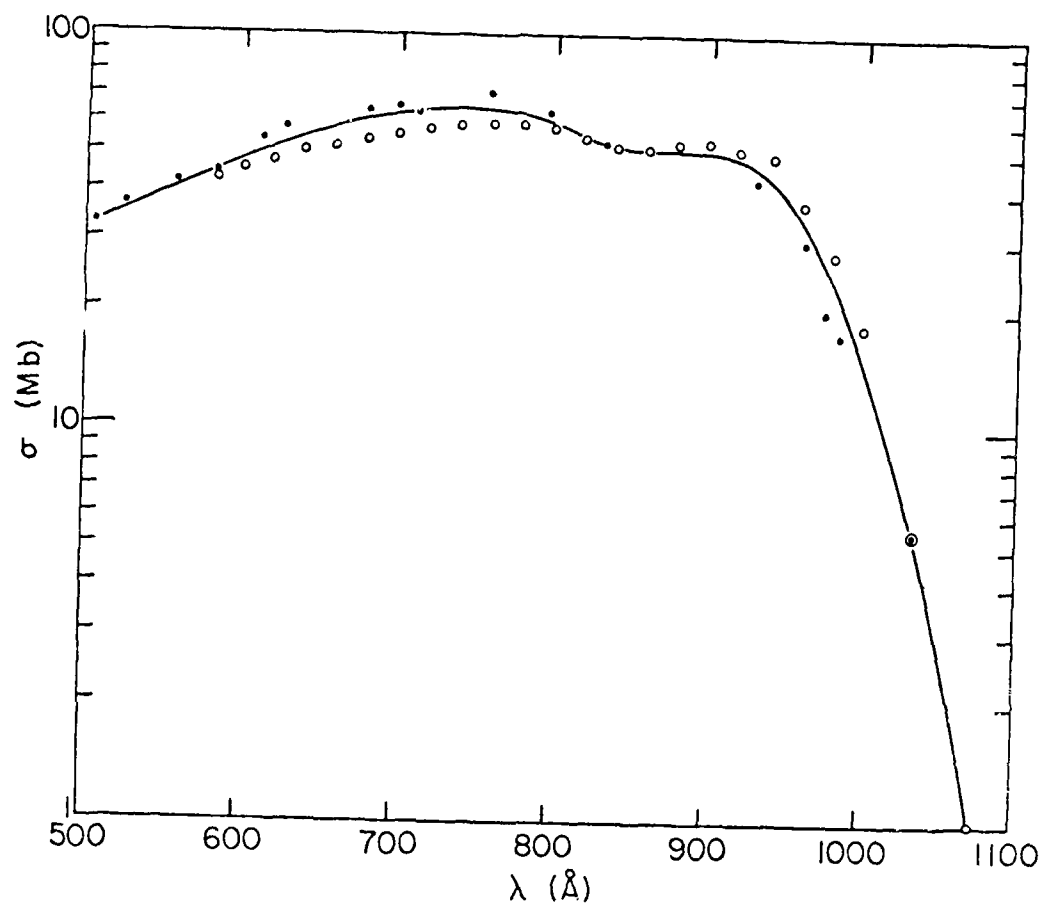
Graphical Data D-2.43

Quantum yield of ionization for C_2H_4 .



Graphical Data D-2.49

Photoabsorption cross section of C_2H_6 .



Graphical Data D-2.50

Photoabsorption cross section for C_2H , in the threshold region.

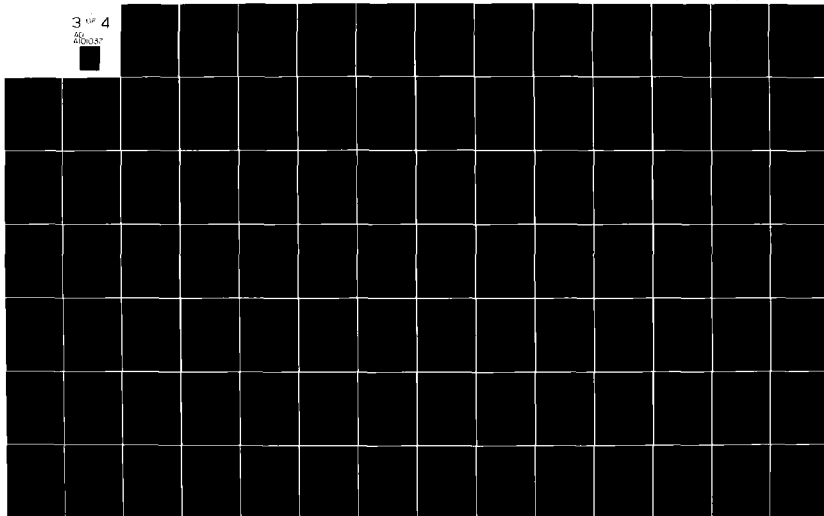
AD-A101 037

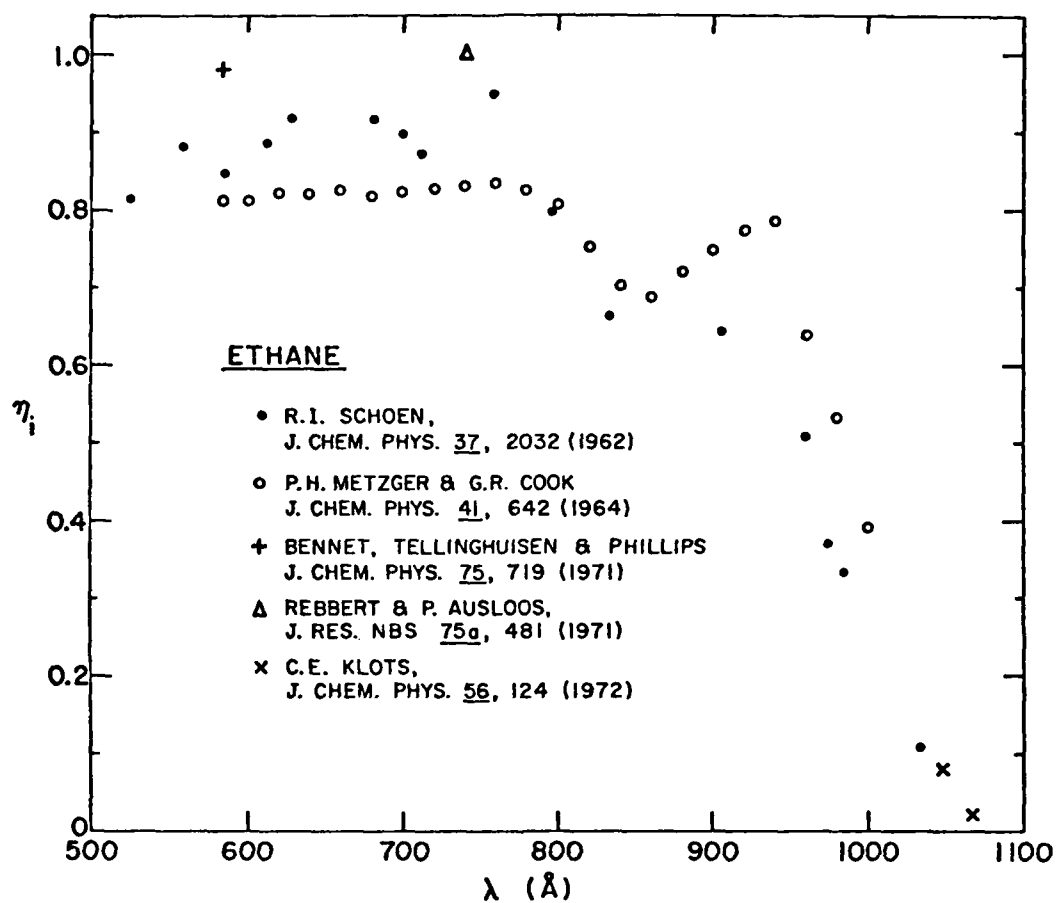
ARMY MISSILE COMMAND REDSTONE ARSENAL AL DIRECTED E--ETC F/G 20/5
COMPILATION OF ATOMIC AND MOLECULAR DATA REVELANT TO GAS LASERS--ETC (11)
DEC 80 E W MCDANIEL, W R FLANNERY, E W THOMAS
DRSMI-RM-81-4-VOL-8

UNCLASSIFIED

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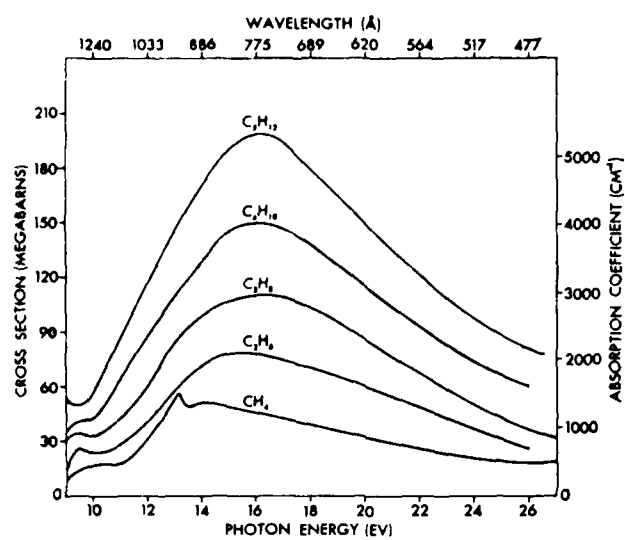
3 of 4
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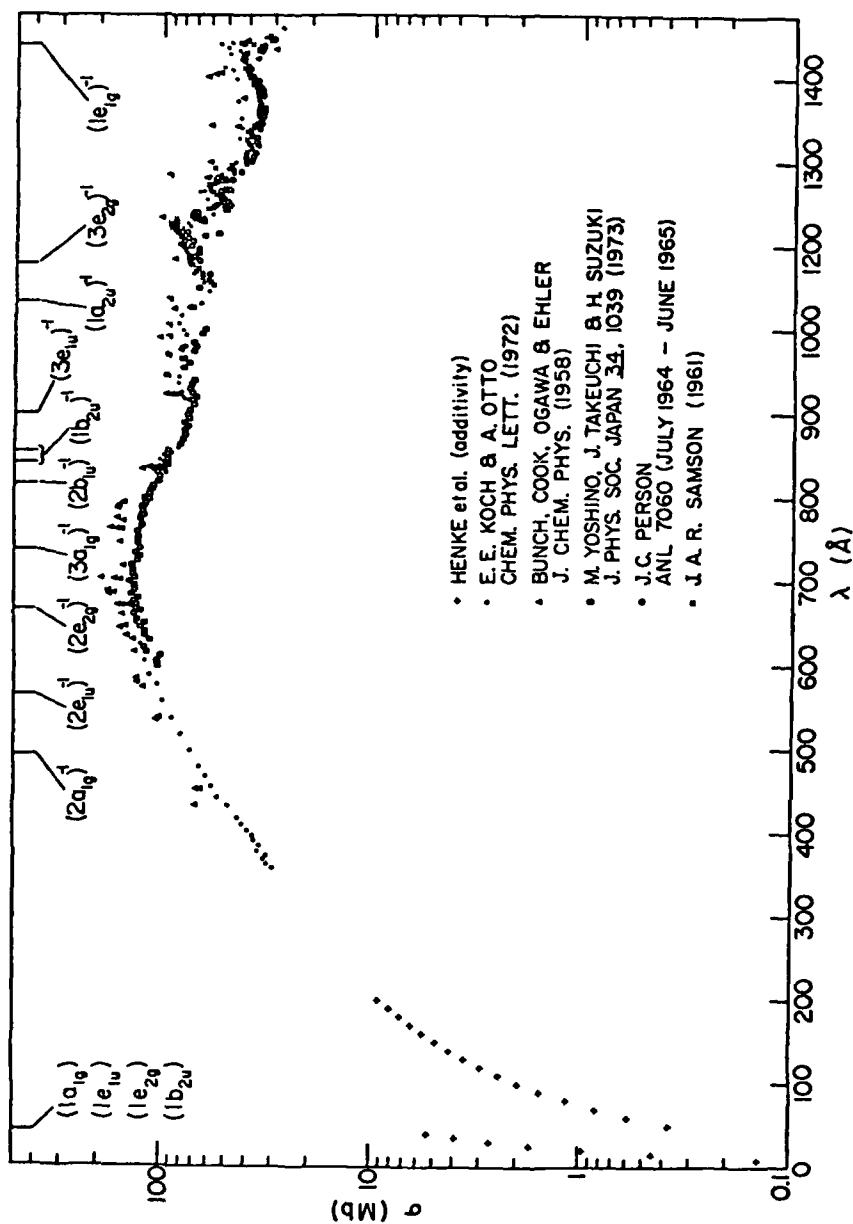
Graphical Data D-2.51

Quantum yield of ionization, η_i , for C_2H_6 .



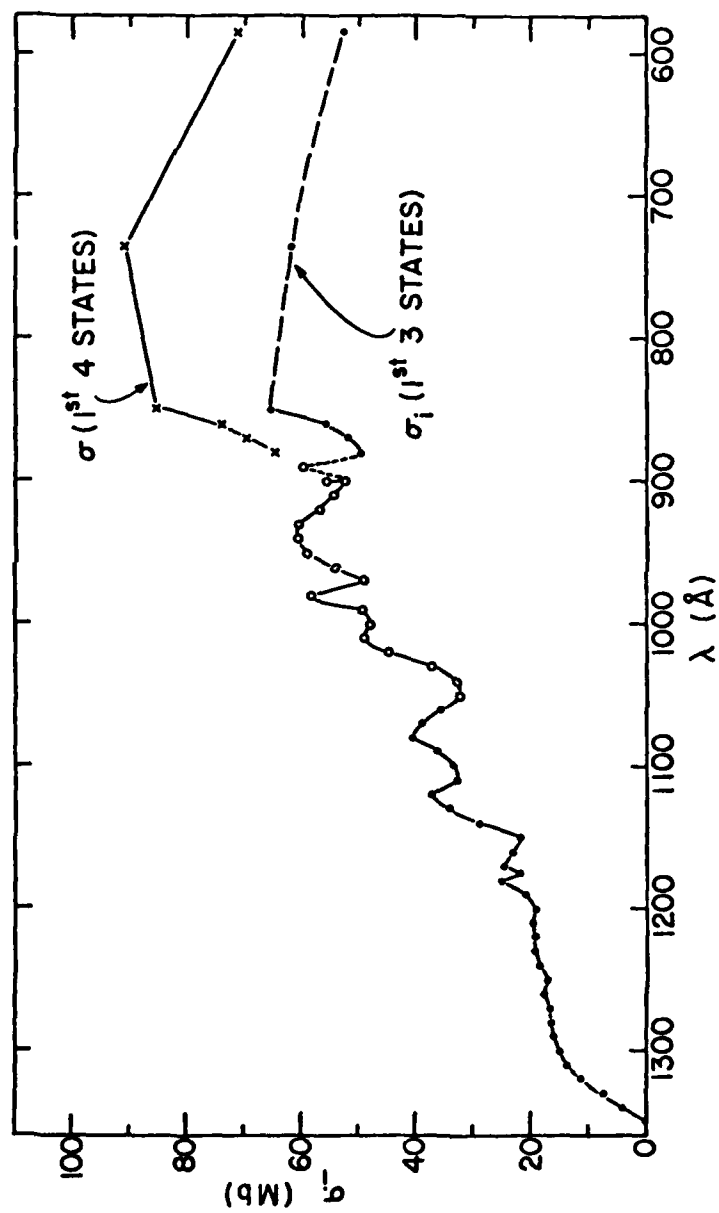
Graphical Data D-2.52

Photoabsorption cross sections of the alkanes



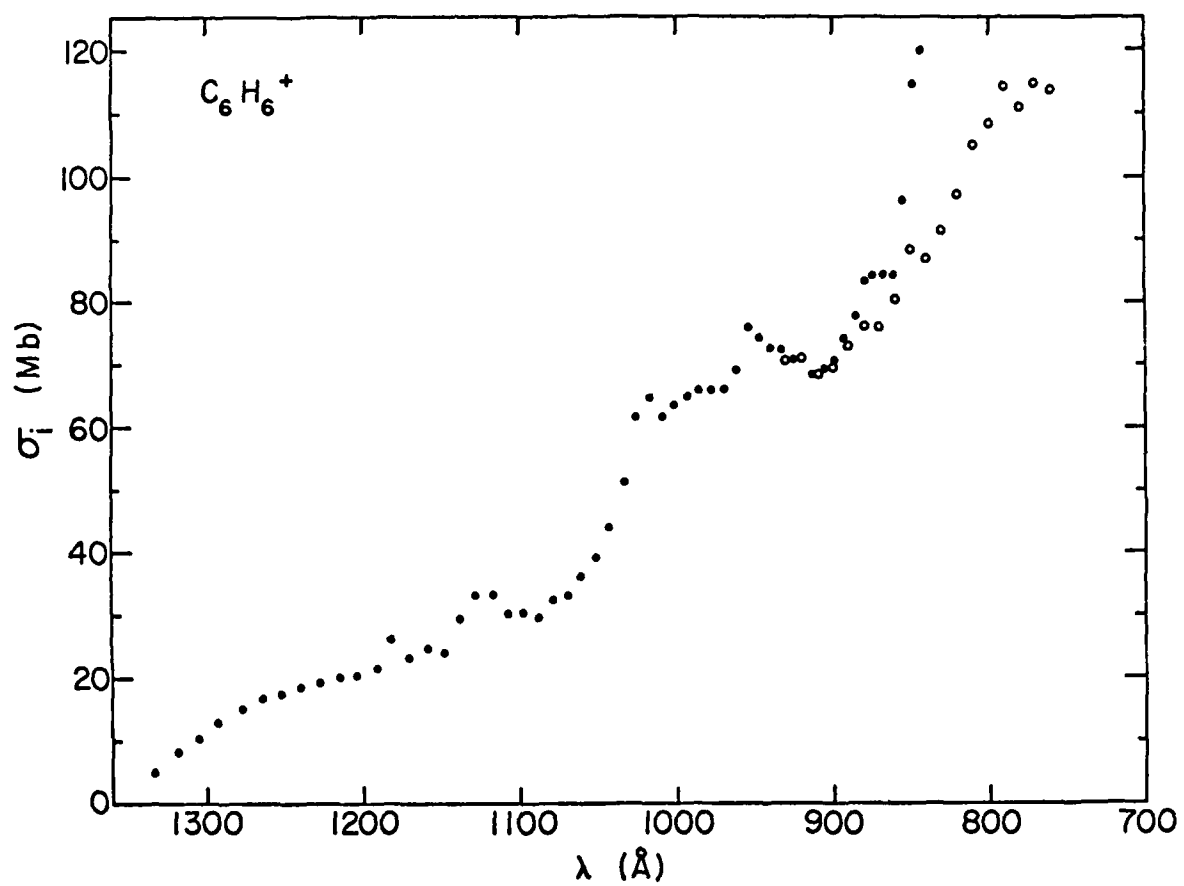
Graphical Data D-2.53

Photoabsorption cross section of C_6H_6 .



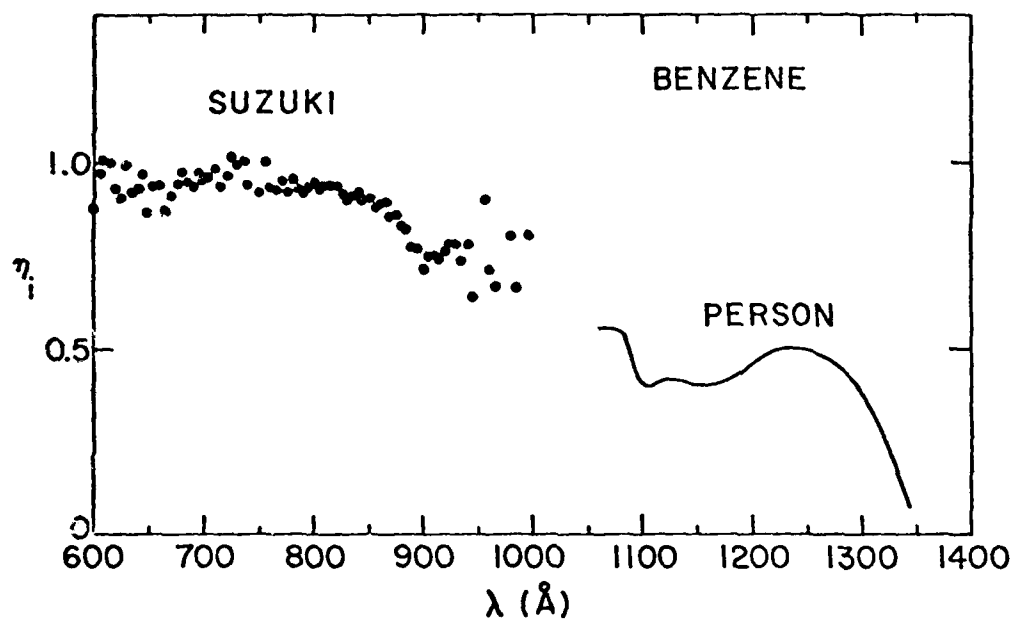
Graphical Data D-2.54

Photoionization cross section of C_6H_6 for production of the lowest three and four states of C_6H_6^+ .



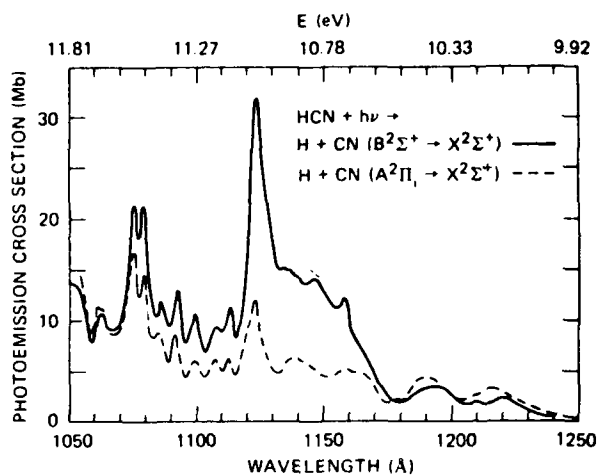
Graphical Data D-2.55

Photoionization cross section of C_6H_6 for production of $C_6H_6^+$.



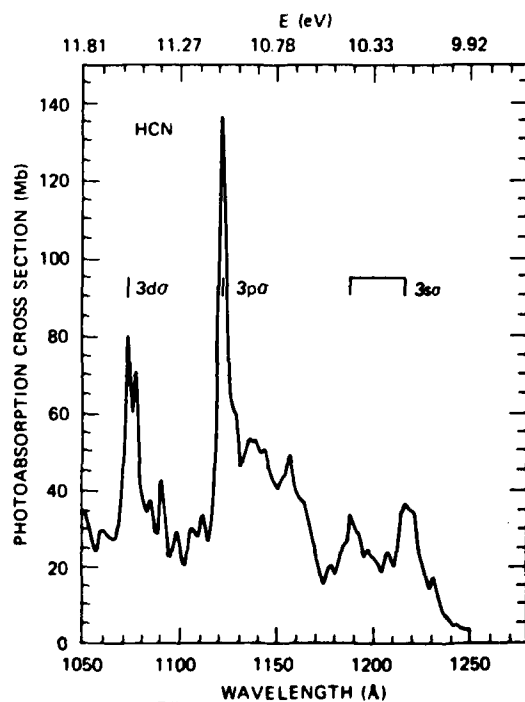
Graphical Data D-2.56

Quantum yield of ionization, η_i , for C_6H_6 .



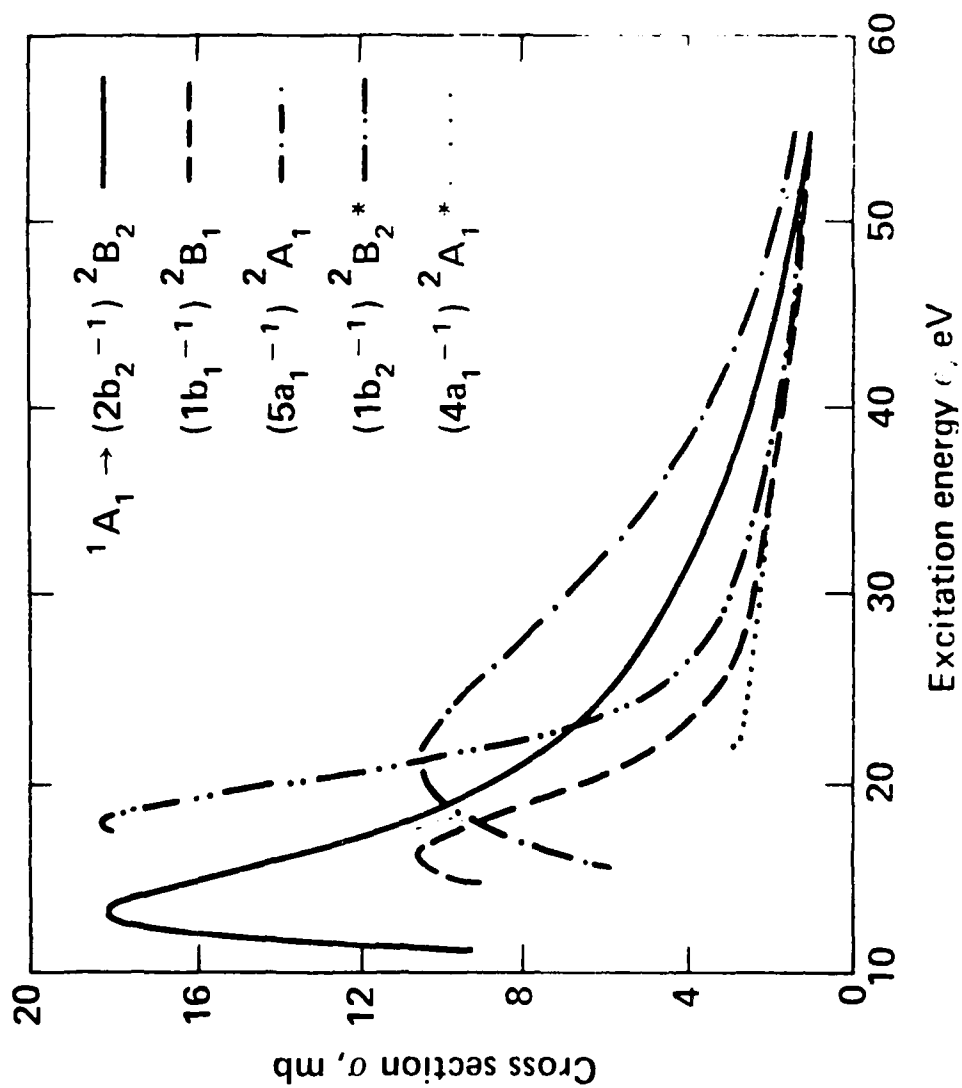
Graphical Data D-2.57

Cross sections for various emissions from HCN photodissociation.



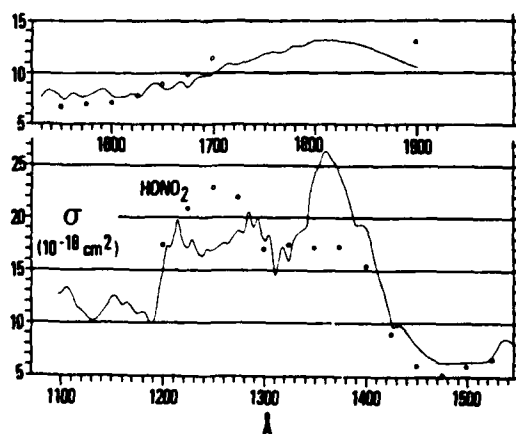
Graphical Data D-2.58

Photoabsorption cross section of HCN with the positions of various electronic states indicated.



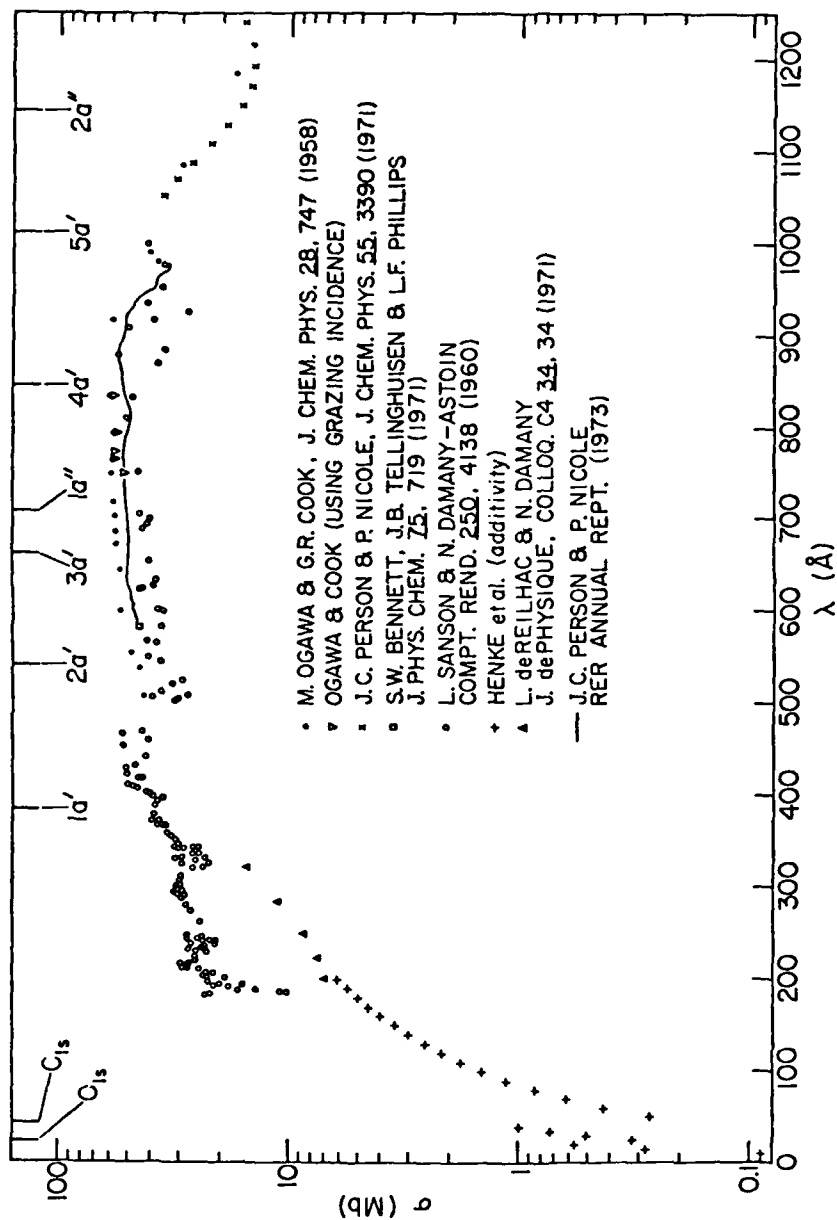
Graphical Data D-2.59

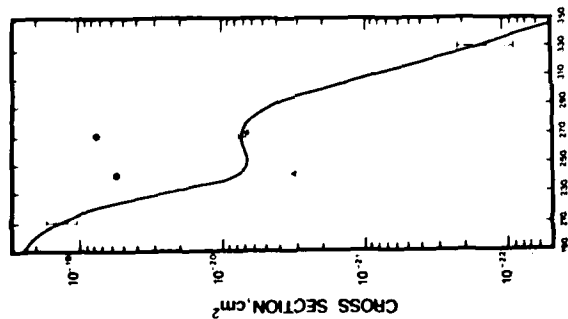
Photoionization cross sections for various channels in $\text{H}_2\text{CO} \rightarrow \text{H}_2\text{CO}^+$ vertical transitions.



Graphical Data D-2.60

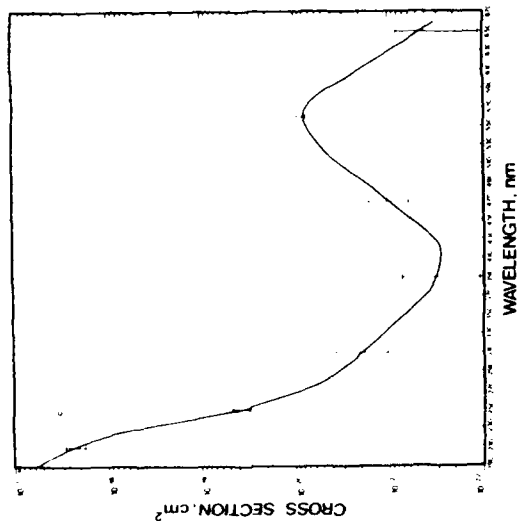
Photoabsorption cross section of HONO_2 (nitric acid)





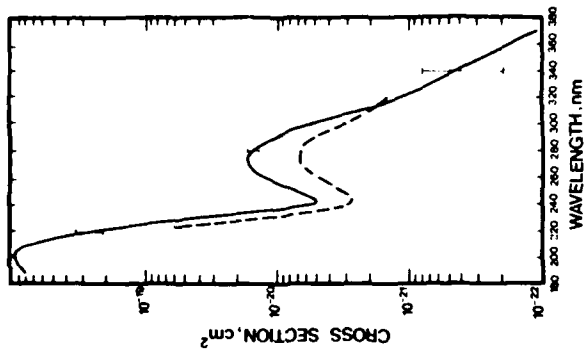
Ultraviolet absorption spectrum of $\text{CF}_3\text{NO}(\text{g})$ at 295 K.

(a)



Absorption spectrum of $\text{CCl}_3\text{NO}(\text{g})$ at 295 K.

(b)

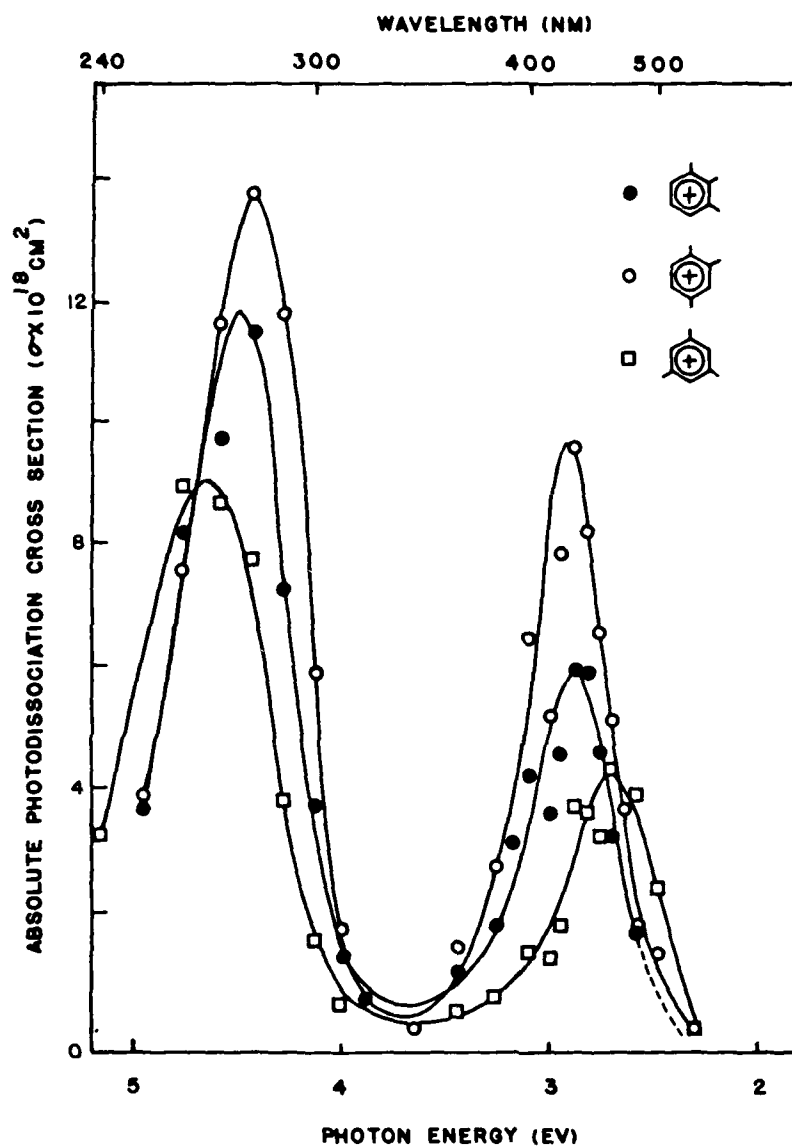


Absorption spectrum of $\text{CCl}_3\text{NO}_2(\text{g})$ at 295 K.

(c)

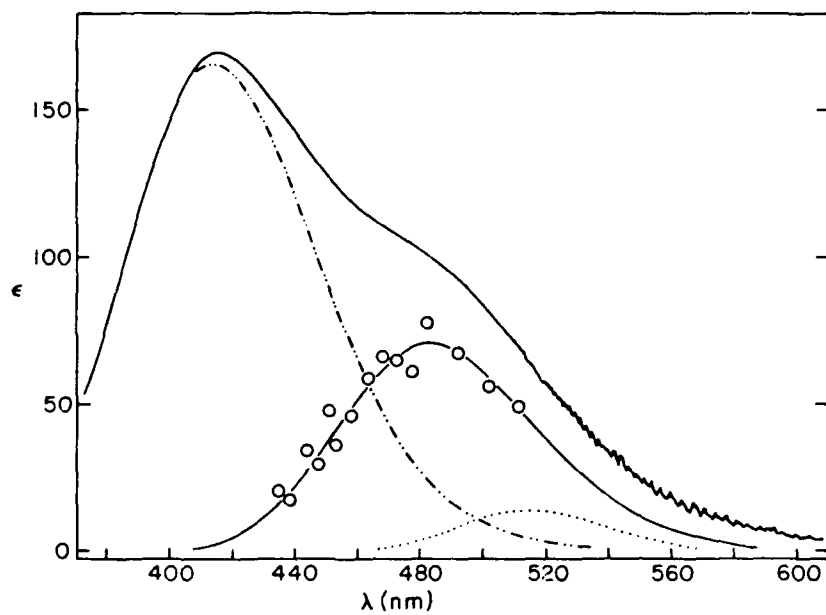
Graphical Data D-2.62 Absorption of spectrum of $\text{CF}_3\text{NO}(\text{g})$, $\text{CCl}_3\text{NO}(\text{g})$ and $\text{CCl}_3\text{NO}_2(\text{g})$ at 295°K.

Reference: These data are from T.D. Allston, M.L. Fedyk, and G.A. Takacs, Chem. Phys. Lett. 60, 97 (1978).



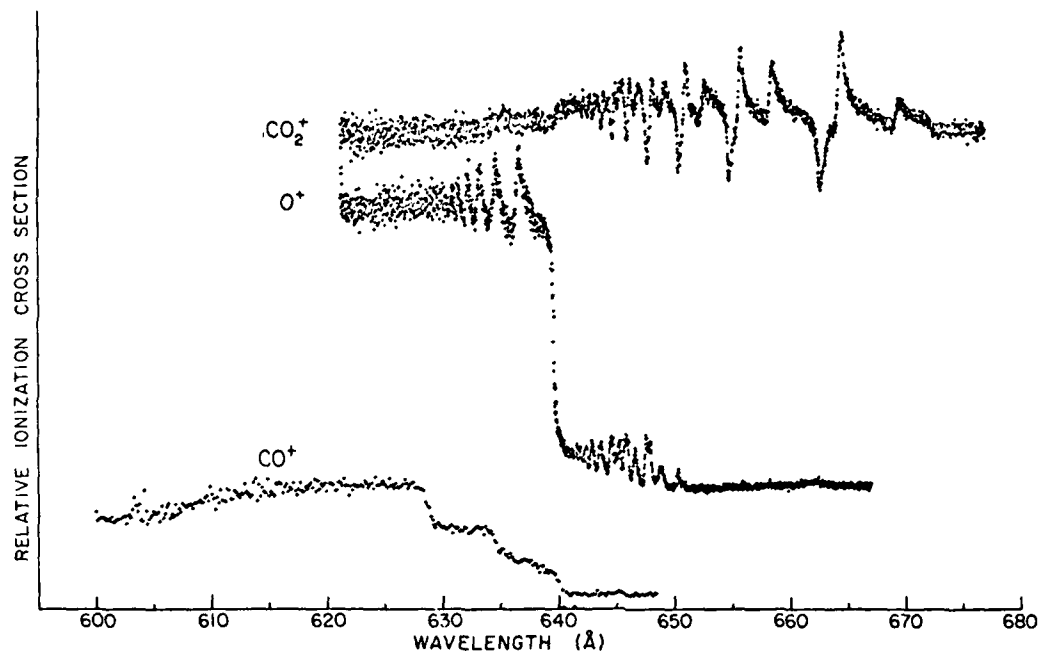
Graphical Data D-2.63

Photodissociation cross sections for three trimethylbenzene isomers



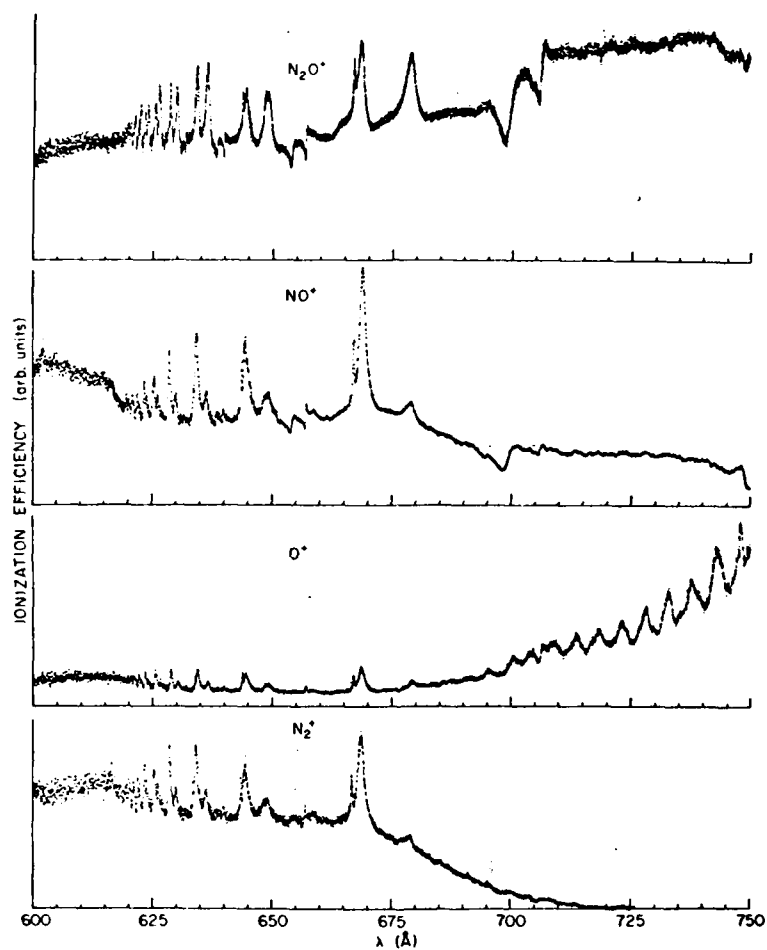
Graphical Data D-2.64

Relative photoabsorption cross section of Br_2 decomposed into $X \cdot B$, $X \cdot A$, and $X \cdot ^1HBr$



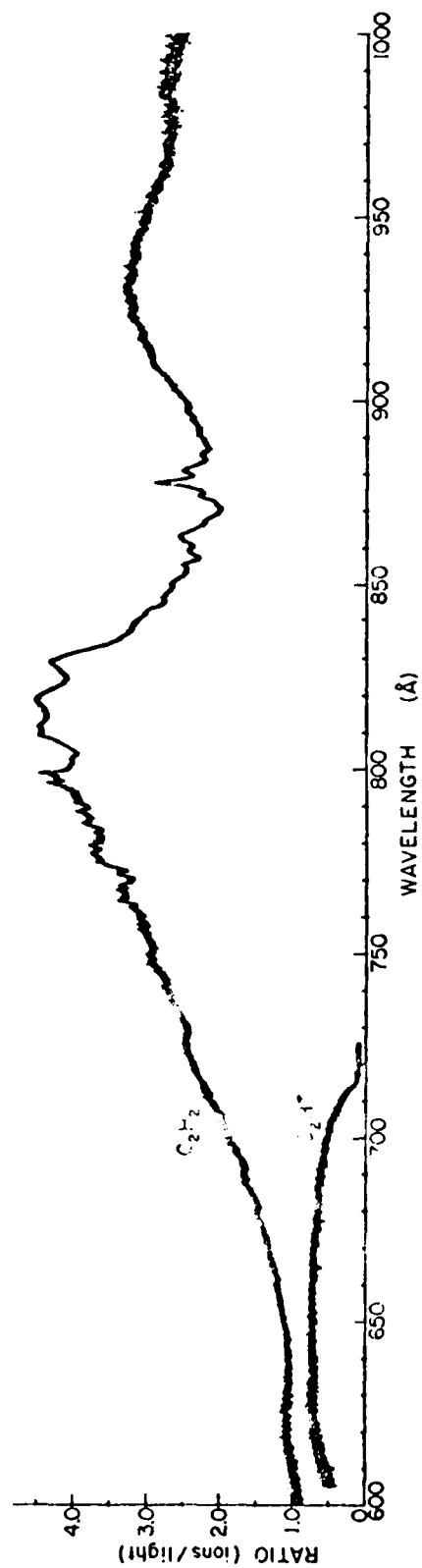
Graphical Data D-2.65

Relative photoionization cross sections of CO_2 for production of various ions.

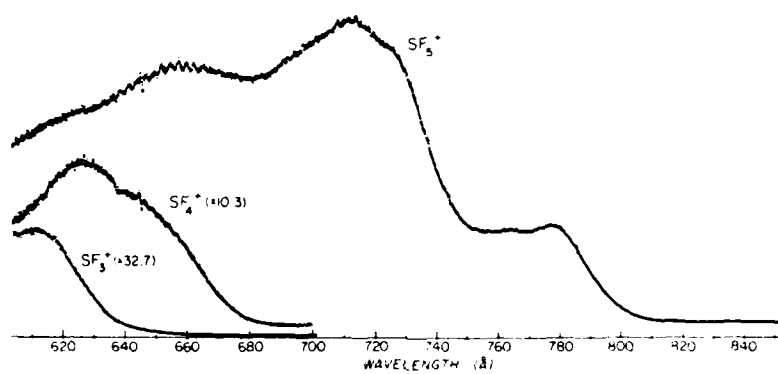


Graphical Data D-2.66

Relative photoionization cross sections of N_2O for production of various ions.

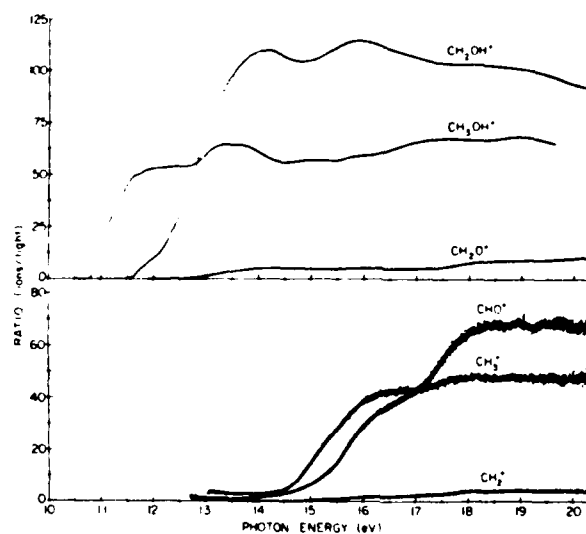


Graphical Data D-2.67
Relative photoionization cross sections of C_2H_2 for production of various ions.



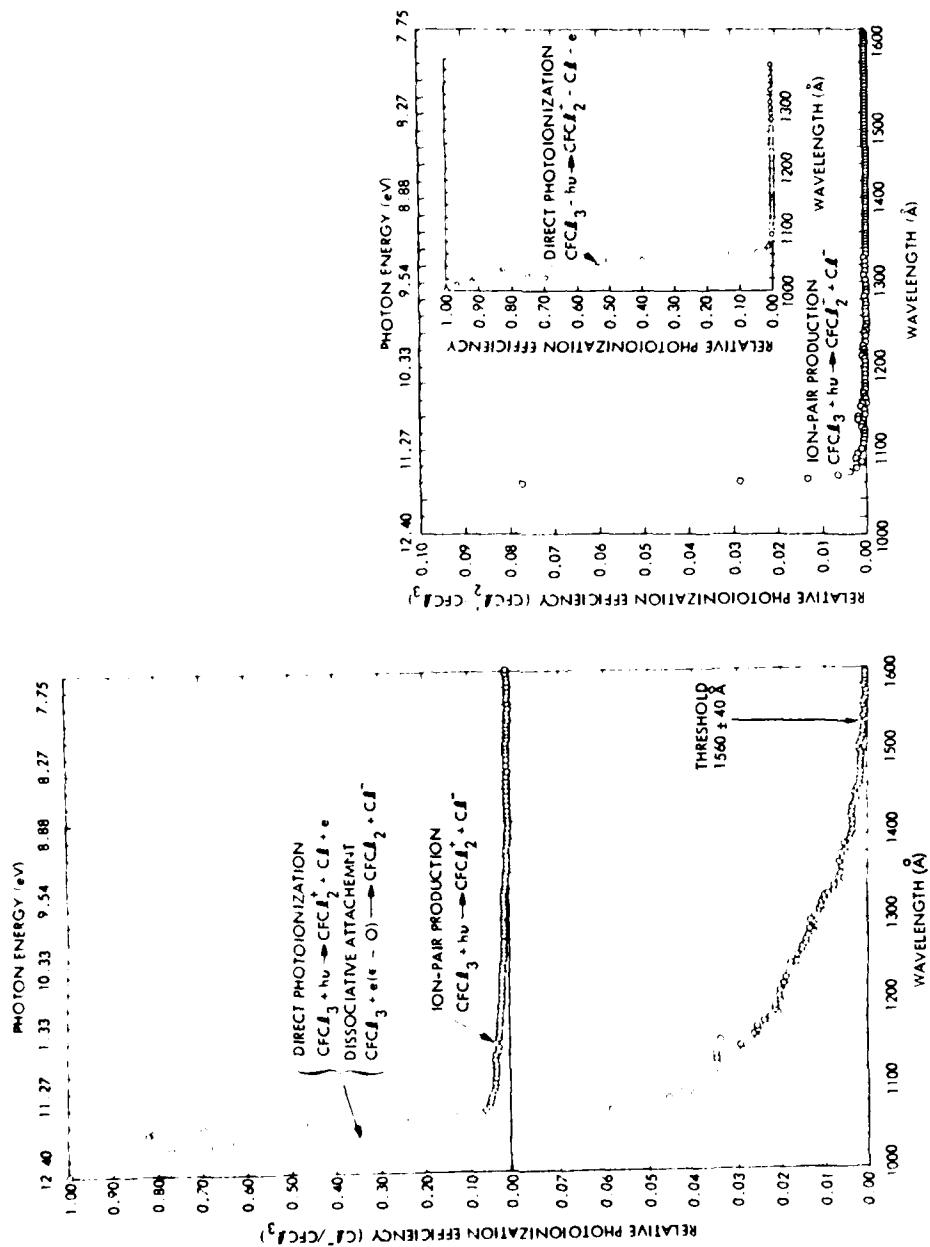
Graphical Data D-2.68

Relative photoionization cross section of SF₆ for production of various ions.



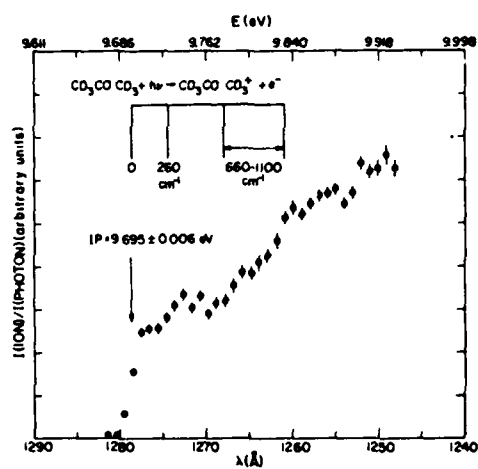
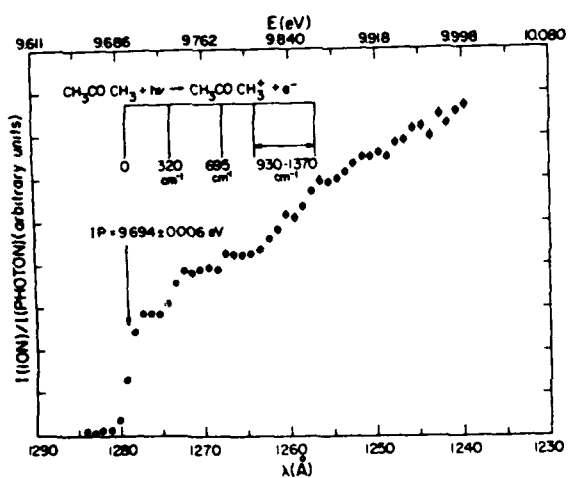
Graphical Data D-2.69

Relative photoionization cross section of CH_3OH for production of various ions.



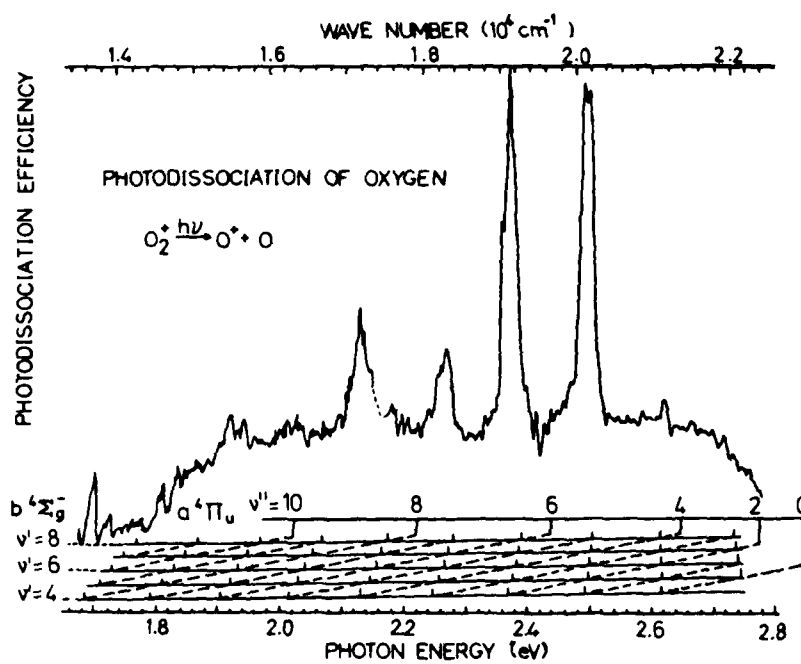
Graphical Data D-2.70

Relative photoionization efficiency for CFC₁₃ production and CFC₁₂⁺ in the photoionization of CFC₁₃.



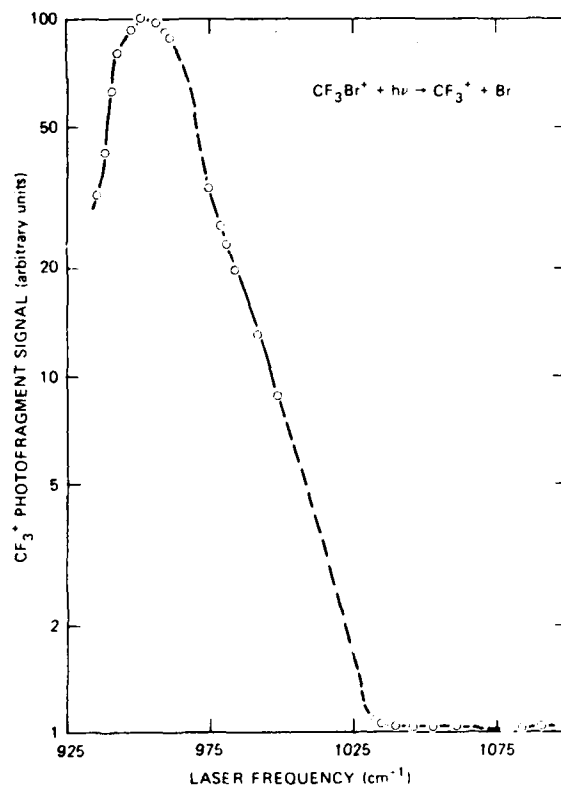
Graphical Data D-2.71

Relative photoionization cross sections for CH_3COCH_3 and CD_3COCd_3 .



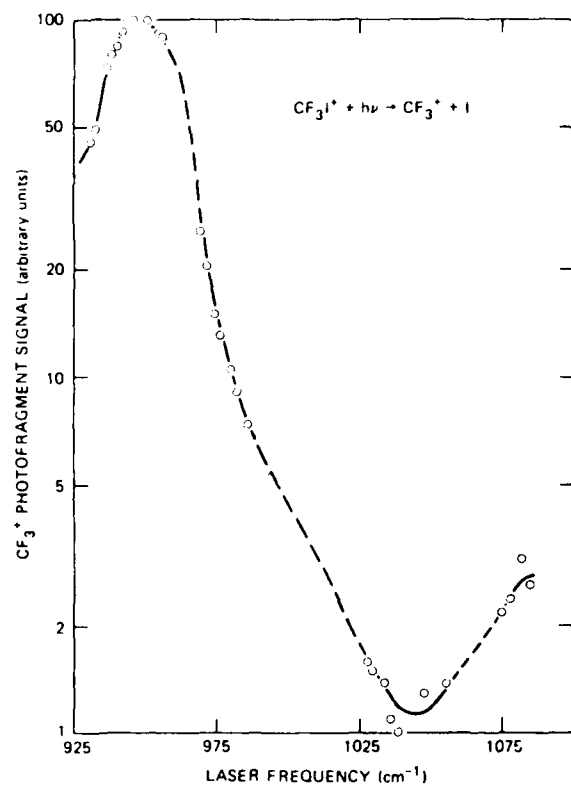
Graphical Data D-2.72

Relative photodissociation cross section of O_2^+ .



Graphical Data D-2.73

Relative photodissociation cross section of $\text{CF}_3\text{Br}^+ \rightarrow \text{CF}_3^+ + \text{Br}$.

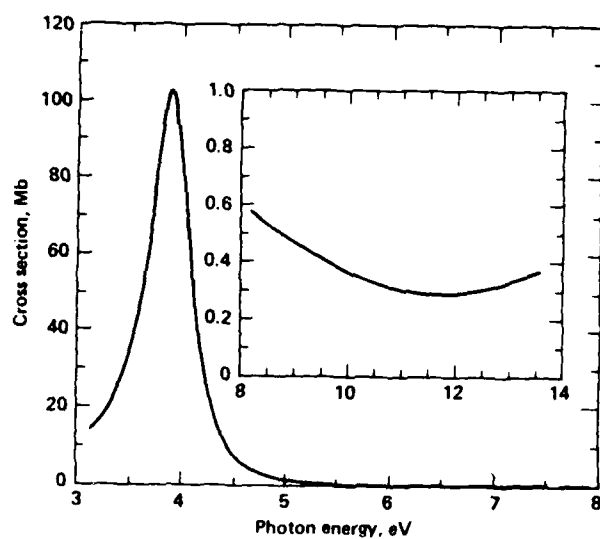


Graphical Data D-2.74

Relative photodissociation cross section of $\text{CF}_3\text{I}^+ \rightarrow \text{CF}_3^+ + \text{I}$.

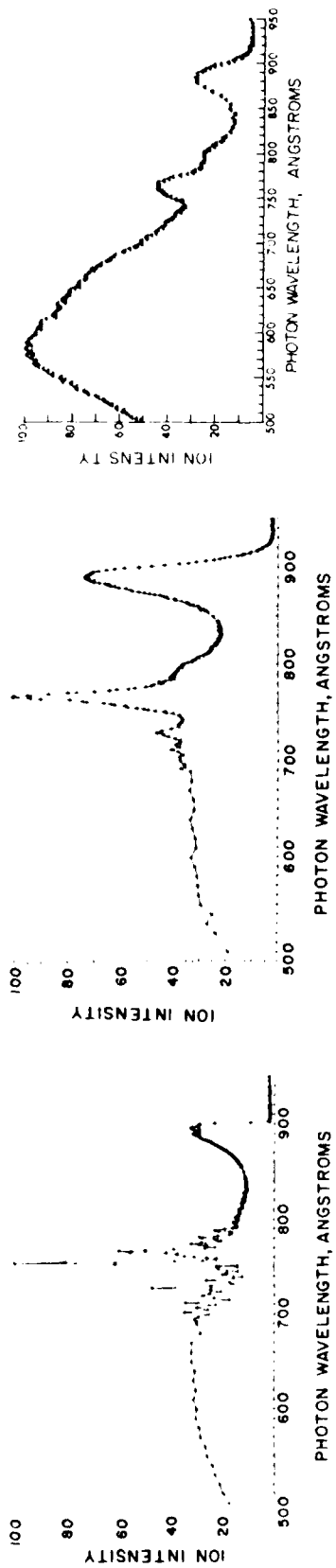
D-3. PHOTOABSORPTION, PHOTOIONIZATION, AND PHOTODISSOCIATION CROSS SECTION
OF MOLECULES AND POSITIVE MOLECULAR IONS (DIMERS AND EXCIMERS)

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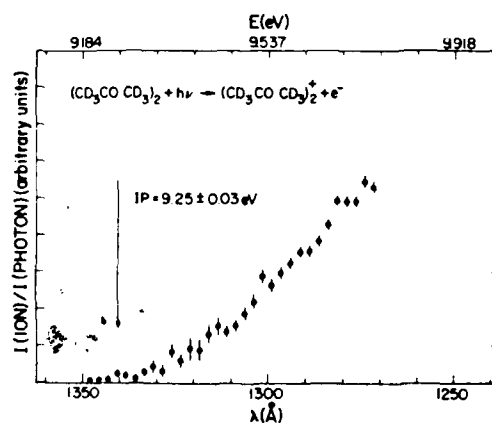
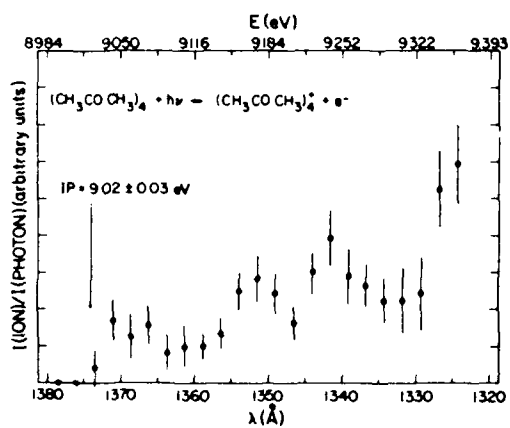
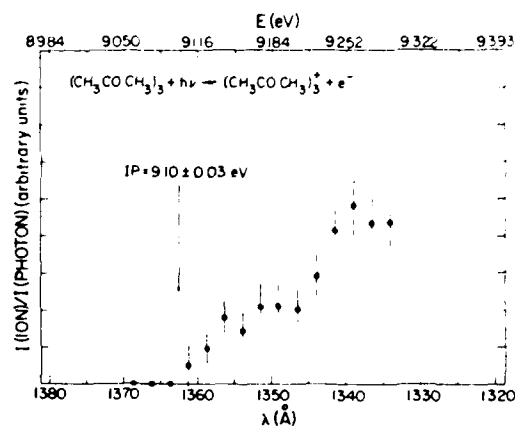
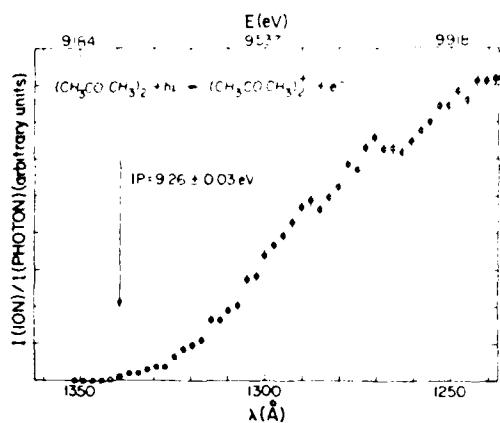
Graphical Data 3.1

Vertical photoionization cross section of the $1\Sigma_g^+$ excimer state of Ar_2 calculated at $R = 4.8a_0$.



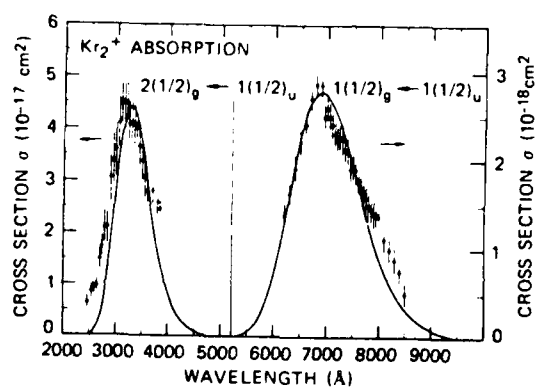
Graphical Data D-3.2

Relative photoionization cross sections for CO, monomers and dimers at 245K.



Graphical Data D-3.3

Relative photoionization cross sections for various dimers (clusters) of CH_3COCH_3 and CD_3COCD_3



Graphical Data D-3.4

Photoabsorption cross section for Kr₂⁺.

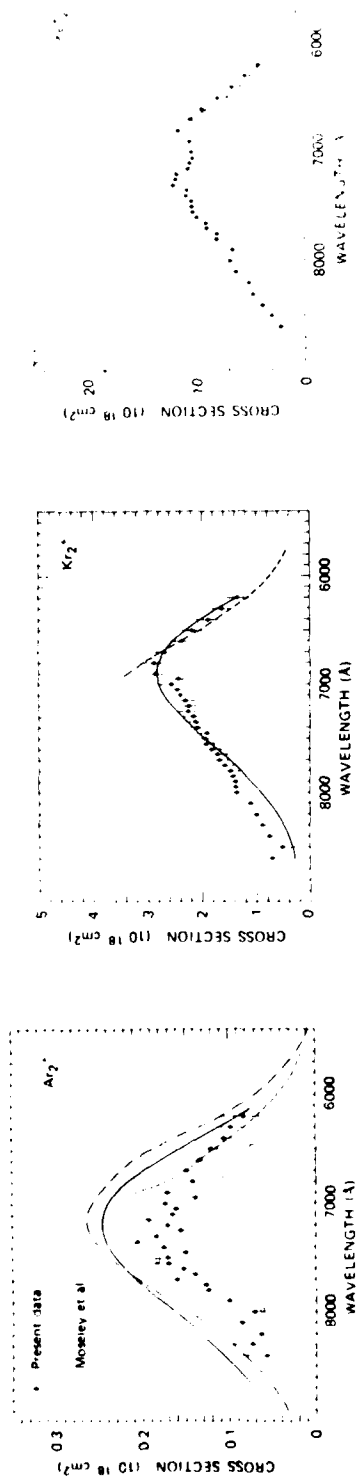
Tabular Data D-3.5

Photodissociation Cross Sections of Ne_2^+ , Ar_2^+ , Kr_2^+ ,
and Xe_2^+ Dimer Ions

$\lambda (\text{\AA})$	Ne_2^+	Ar_2^+	Kr_2^+	Xe_2^+
5309		< 0.012	0.033 ± 0.004	0.028 ± 0.005
5208		< 0.028	0.026 ± 0.006	0.022 ± 0.004
4825		0.024 ± 0.012	0.100 ± 0.014	0.102 ± 0.011
4762	< 0.11	0.035 ± 0.016	0.122 ± 0.013	0.147 ± 0.016
4680		0.082 ± 0.022	0.203 ± 0.022	0.244 ± 0.024
4579		0.130 ± 0.012	0.39 ± 0.03	0.64 ± 0.03
4131	< 0.09	1.05 ± 0.10	3.18 ± 0.28	5.50 ± 0.55
4067		1.60 ± 0.17	3.63 ± 0.37	7.32 ± 1.33
{3569	1.93 ± 0.20	13.3 ± 1.1	24.8 ± 1.9	29.6 ± 2.0
{3507				

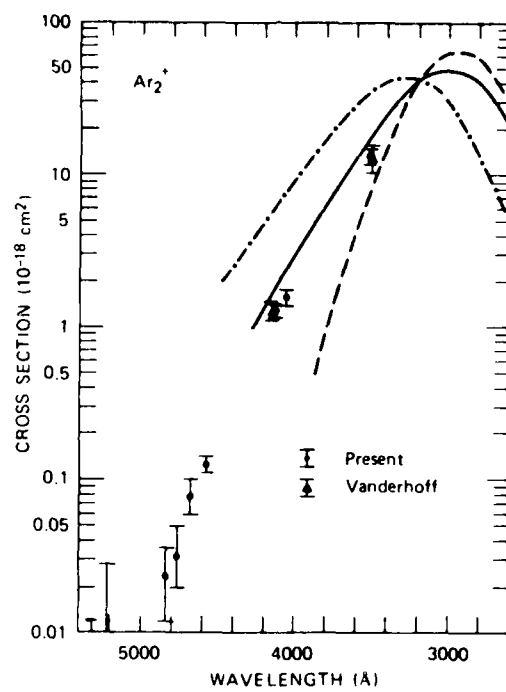
Note: Ne_2^+ and Ar_2^+ were measured at 10 Td, and Kr_2^+ and Xe_2^+ at 20 Td.

Reference: These data were taken from L. C. Lee and G. P. Smith, Phys. Rev. A 19, 2329 (1979).



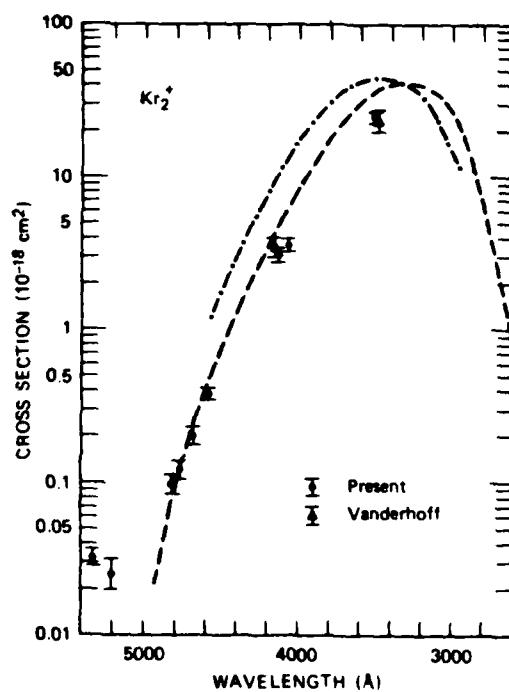
Graphical Data D-3.6

Photodissociation cross sections of Ar_2^+ , Kr_2^+ , and Xe_2^+ dimer ions.



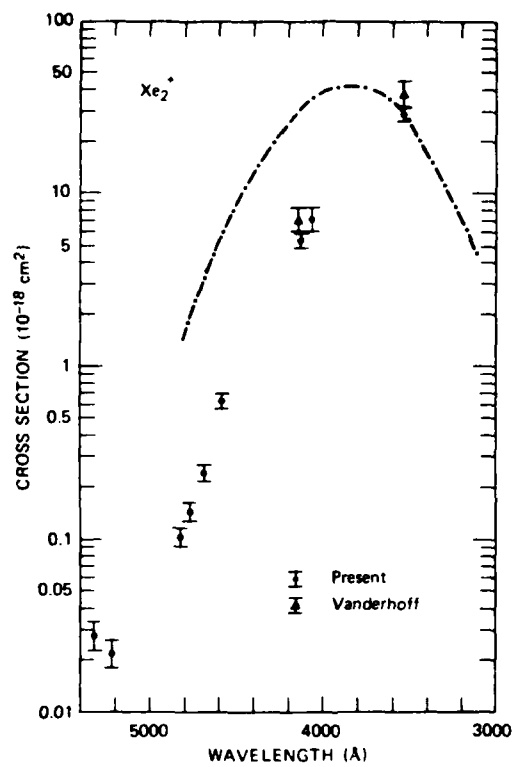
Graphical Data D-3.7

Photodissociation cross sections for Ar_2^+ .



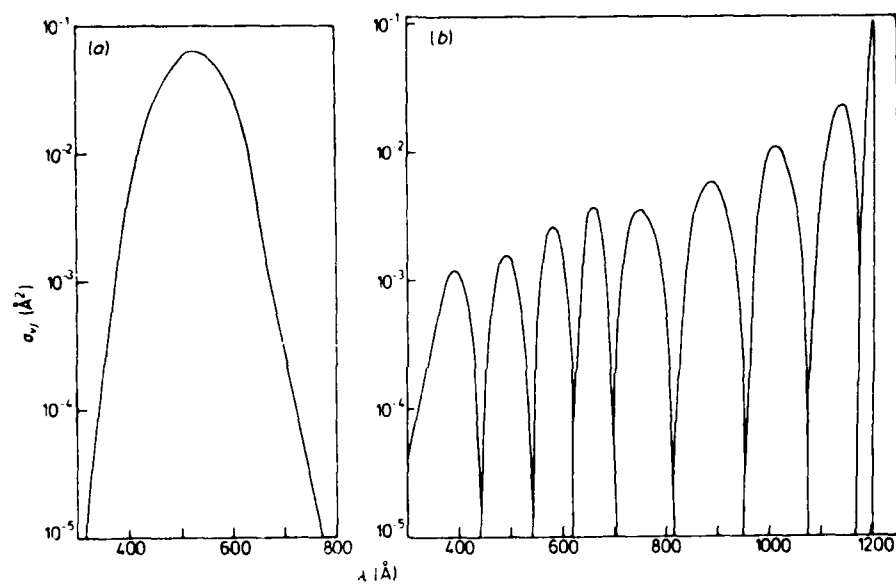
Graphical Data D-3.8

Photodissociation cross section for Kr_2^+ .



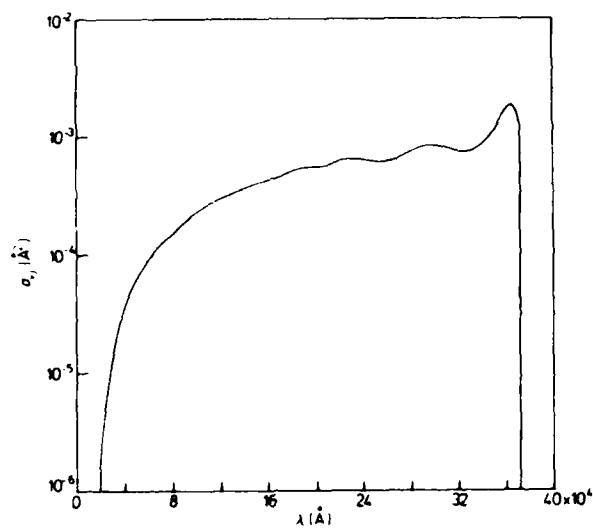
Graphical Data D-3.9

Photodissociation cross sections for Xe_2^+ .



Graphical Data D-3.10

Partial photodissociation cross section for HeH^+ by electronic excitation from the vibrational level (a) $v = 0$ and (b) $v = 8$ (with $j = 1$).

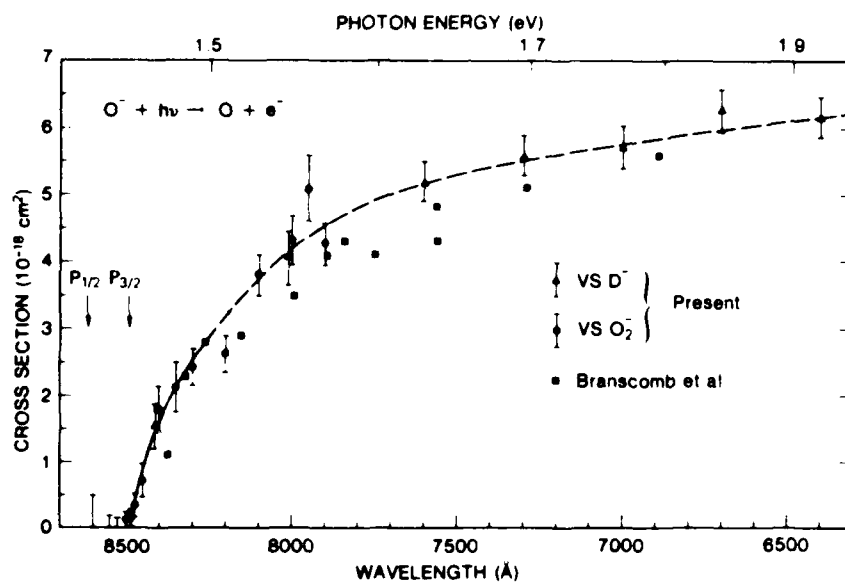


Graphical Data D-3.11

Partial photodissociation cross section for HeH^+ by vibrational excitation from the vibrational level $v = 8$ (with $j = 1$).

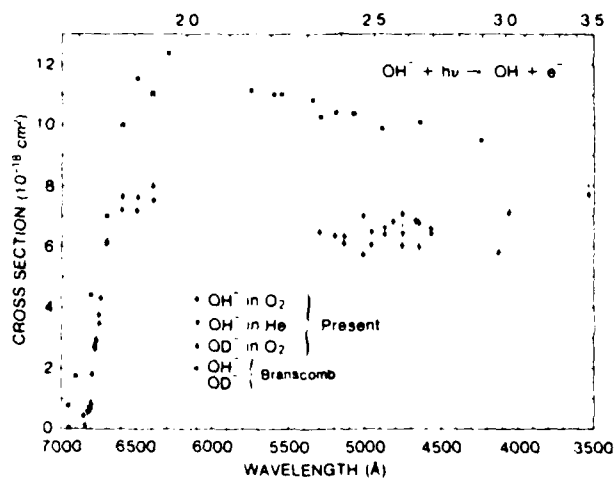
D-4. PHOTODETACHMENT, PHOTODISSOCIATION, AND PHOTODESTRUCTION OF NEGATIVE IONS

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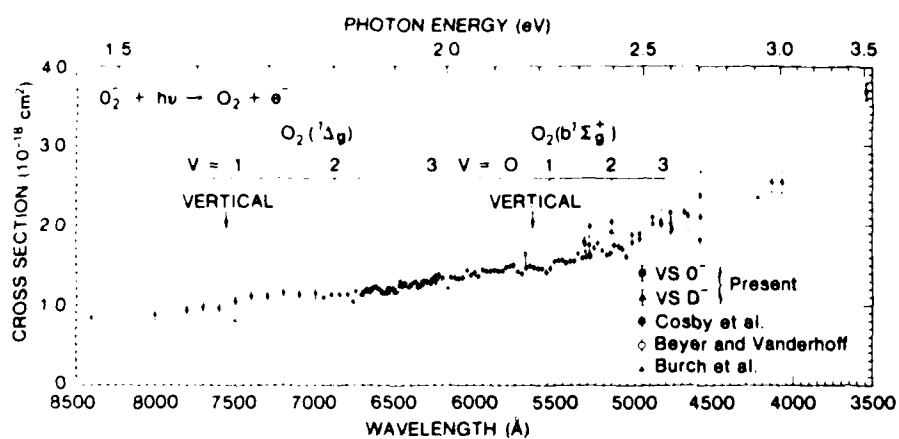
Graphical Data D-4.1

Photodetachment cross section for O^- placed on an absolute scale by normalization to D^- and O_2^- .



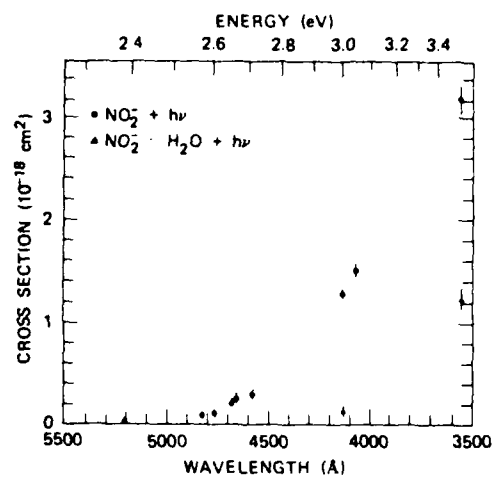
Graphical D-4.2

Photodetachment cross sections for OH^- and OD^- measured in O_2 and He .



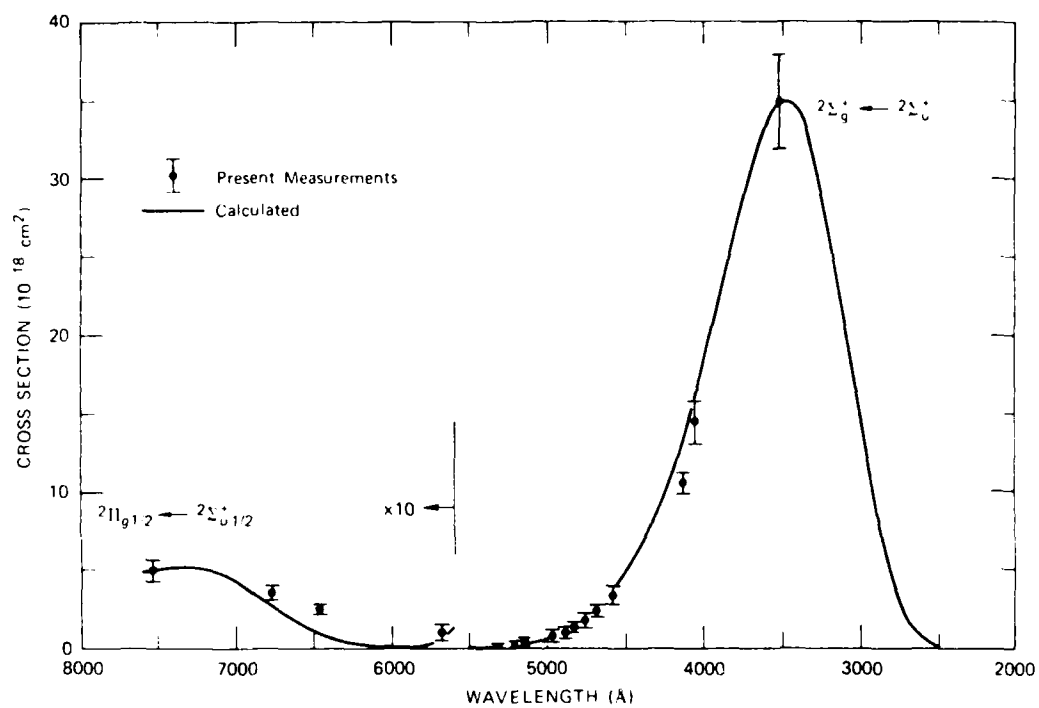
Graphical Data D-4.3

Photodetachment cross section for O_2^- placed on an absolute scale by normalizing to D^- and O^- .



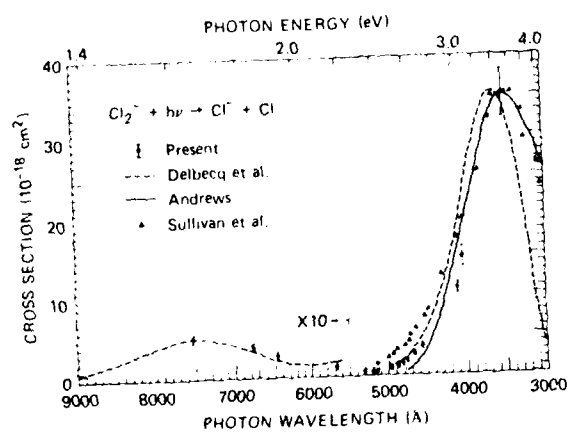
Graphical Data D-4.4

Photodetachment cross section for thermal NO_2^- and $\text{NO}_2^-\cdot\text{H}_2\text{O}$.



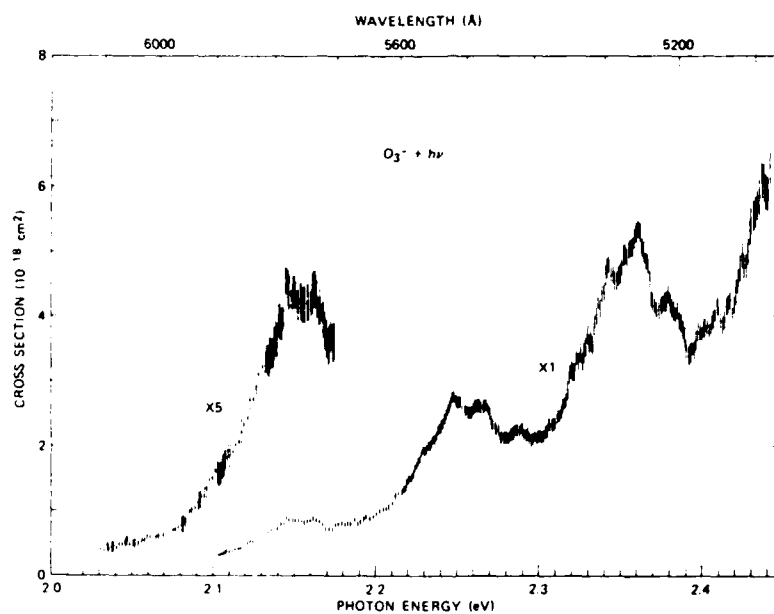
Graphical Data D-4.5

Measured and calculated photodissociation cross section for C_2 .



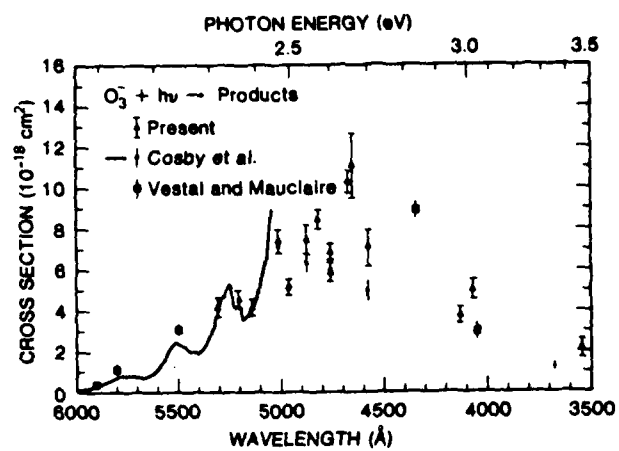
Graphical Data D-4.6

Photodissociation cross section for Cl_2^- .



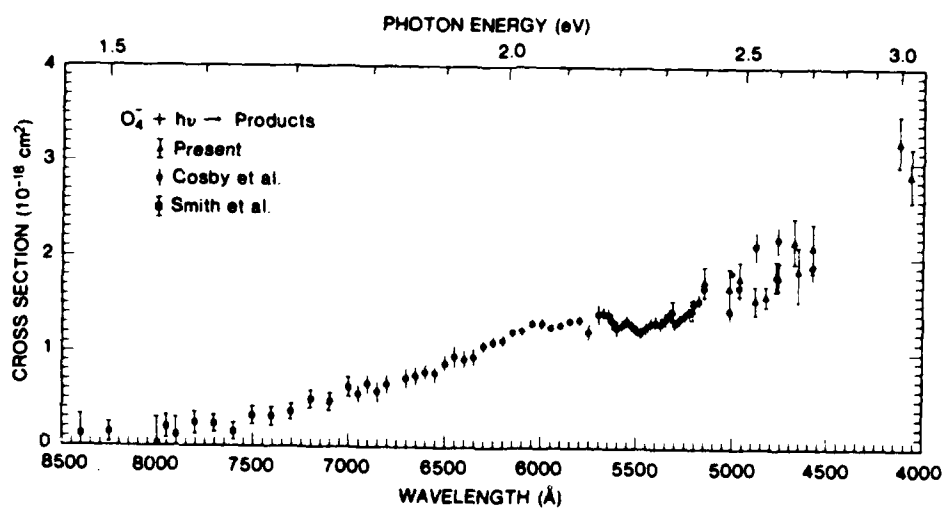
Graphical Data D-4./

Photodestruction cross section of O_3^- . The predominant process observed is photodissociation into $O^- + O_2$.



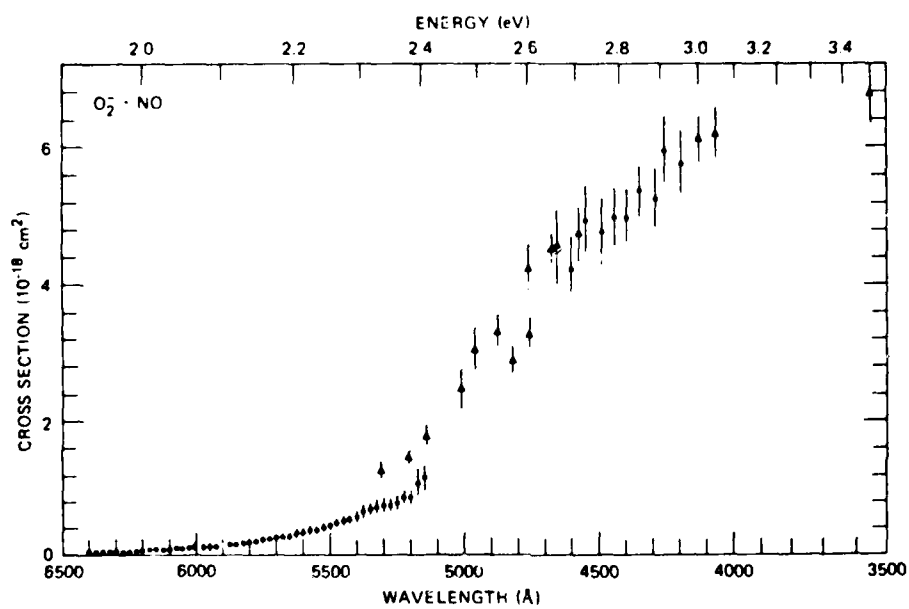
Graphical Data D-4.8

Photodestruction cross section for O_3^- .



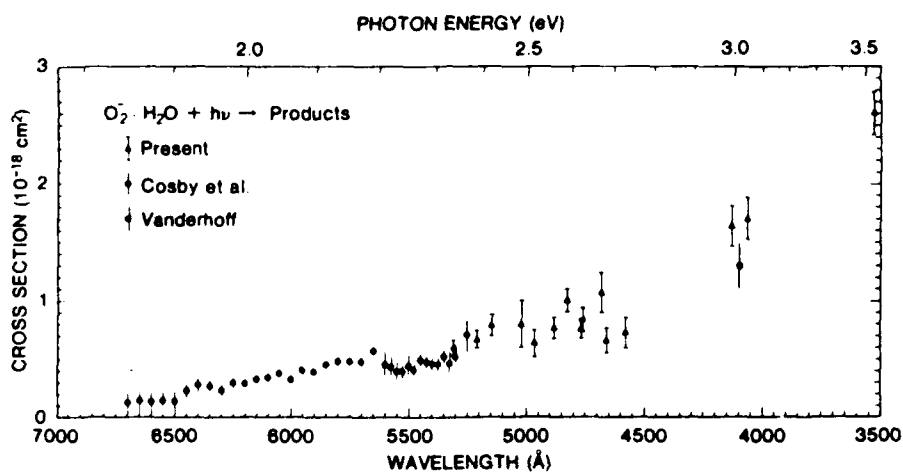
Graphical Data D-4.9

Photodestruction cross section for O_4^- .



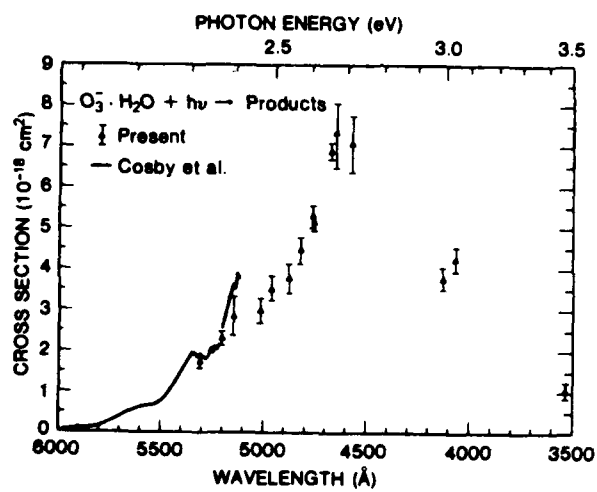
Graphical Data D-4.10

Photodestruction cross section for $O_2^- \cdot NO$.



Graphical Data D-4.11

Photodestruction cross section for $O_2^- + H_2O$.



Graphical Data D-4.12

Photodestruction cross section for $O_3 \cdot H_2O$.

Tabular Data D-4.13

Photodestruction Cross Section Upper Limits
for NO_2^- , NO_3^- , $\text{O}_3^- \cdot \text{NO}$, and Hydrates

E (eV)	λ (Å)	NO_2^-	$\text{NO}_2^- \cdot \text{H}_2\text{O}$	NO_3^-	$\text{NO}_3^- \cdot \text{H}_2\text{O}$	$\text{O}_3^- \cdot \text{NO}$	$\text{O}_3^- \cdot \text{NO} \cdot \text{H}_2\text{O}$
1.503	8250					< 0.067	< 0.072
1.653	7500		< 0.041				< 0.036
1.746	7100					< 0.054	
2.335	5309	< 0.029	< 0.014	< 0.011	< 0.10		
2.381	5208	0.023 ± 0.009	< 0.026	< 0.054	< 0.10		
2.569	4825		< 0.046	< 0.075	< 0.22		
2.603	4762		< 0.095	< 0.092	< 0.12		
2.649	4680		< 0.071	< 0.106	< 0.16		
2.708	4579		< 0.045				
3.001	4131		0.16 ± 0.03	< 0.074	< 0.10		
3.479	3564		1.23 ± 0.16	0.10 ± 0.03	0.41 ± 0.08		
3.535	3507						

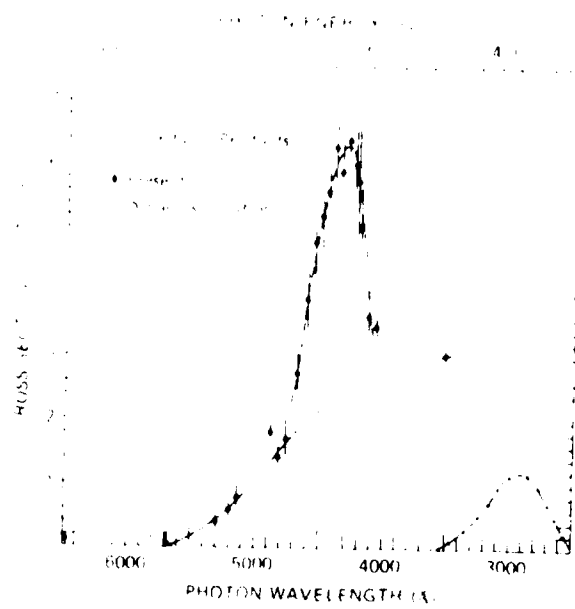
Reference: The above data were taken from G. P. Smith, L. C. Lee,
and P. C. Cosby, J. Chem. Phys. 71, 4468 (1979).

Tabular Data D-4.14

Photodestruction Cross Sections for Cl_2^- ,
 ClO^- , Cl_2^- , and Br Cl^-

λ (Å)	Cl_2^-	ClO^-	Cl_2^-	BrCl^-
3507				
3569	35.1 ± 3.0	3.06 ± 0.32	6.72 ± 0.65	3.29 ± 0.28
4067	14.9 ± 1.1	3.49 ± 0.82		
4131	10.8 ± 0.4	3.64 ± 0.36	0.79 ± 0.07	0.94 ± 0.20
4579	3.55 ± 0.42			
4680	2.41 ± 0.28	2.77 ± 0.32	< 0.3	0.30 ± 0.12
4762	1.64 ± 0.19	1.75 ± 0.25		< 0.24
4765	1.85 ± 0.20			
4825	1.46 ± 0.18	1.44 ± 0.24		
4880	1.19 ± 0.13	1.82 ± 0.27		
4965	0.99 ± 0.12			
5145	0.43 ± 0.06	0.82 ± 0.32		
5208	0.39 ± 0.05	0.60 ± 0.11		< 0.14
5309	0.28 ± 0.03	0.44 ± 0.06	< 0.15	< 0.08
5682	0.11 ± 0.05	0.23		
6471	0.25 ± 0.03	0.11 ± 0.05	< 0.02	< 0.03
6764	0.37 ± 0.04			
7525	0.51 ± 0.06	0.01		

Reference: The above data were taken from L. C. Lee, G. P. Smith, J. T. Moseley, P. C. Cosby, and J. A. Guest, J. Chem. Phys. 70, 3237 (1979).



Graphical Data p-4, 15

For the theoretical section for $\lambda = 4000$.

E. TRANSPORT PROPERTIES OF ELECTRONS, IONS, AND NEUTRALS IN GASES

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(No new entries here.)	

See Chapter E in Vol. II and Chapter E in Vol. V for data.

I. TRANSPORT PROPERTIES OF ELECTRONS, IONS,
AND NEUTRALS IN GASES

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Acknowledgement: It is a pleasure to acknowledge the expert help of Professor Harry W. Ellis in the preparation of this chapter.

E-1. TRANSPORT PROPERTIES OF ELECTRONS

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Definitions and Relationships

- v_d Drift velocity of electrons — average velocity along the field direction in a gas exposed to a constant, uniform electric field E . v_d is usually expressed in units of cm/sec.
- K Mobility of the electrons, defined by $K = v_d/E$. K is usually expressed in $\text{cm}^2/\text{V-sec}$.
- E/N Electron energy parameter = ratio of the electric field intensity to the gas number density. E/N is usually expressed in units of V-cm^2 , or in Townsends, where $1 \text{ Td} = 10^{-17} \text{ V-cm}^2$.
- Td Unit of E/N , the "Townsend" = 10^{-17} V-cm^2 .
- D Diffusion coefficient of the electrons. A scalar at low E/N , D is then related to the mobility by the Einstein (or Nernst-Townsend) relation $K = eD/kT$, where T is the gas temperature, e the electronic charge, and k the Boltzmann constant. At higher E/N , D is a tensor quantity.
- $D_\perp$ The component of the diffusion tensor perpendicular to the electric field.
- $D_\parallel$ The component of the diffusion tensor parallel to the electric field.

 $D_\perp/K$ and $D_\parallel/K$ are measures of the average electron energy at a given E/N . In the limit $E/N \rightarrow 0$, $D_\parallel = D_\perp = D$, the scalar diffusion coefficient.
- α The first Townsend ionization coefficient. Usually it is expressed as α/N , which then has units of cm^2 .
- a The electron attachment coefficient, usually expressed as a/N , which has units of cm^2 .

For electrons in a given gas at a given temperature, v_d , NK , $ND_\perp$, $ND_\parallel$, α/N , a/N , and the average electron energy are functions of E/N alone, N being the gas number density.

Before about 1970, the energy parameter was usually expressed in terms of E/p , where p is the gas pressure in units of torr. To convert from E/p to E/N , one may use the relation

$$E/N \text{ (in Td)} = (1.0354 \cdot T \cdot 10^{-2}) (E/p) \text{ where } T \text{ is the gas temperature.}$$

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Table E-1.1. Sources of Electron Mobility Data Not Presented in Previous Volumes.*

ELECTRON MOBILITY

Gas	Est. Gas Temp ($^{\circ}$ K)	Approx. E/N (Td)	Ref.
Cs	310 - 414 566 - 725	4 - 110 1 - 50	10 2
Hg	468 - 608	0.07 - 14	8
Na	724 - 862	2 - 70	8
Tl	1213 - 1273	0.6 - 15	8
Xe (high density)	163 - 288	10^{-4} - 0.2	6
CH ₄ (high density)	91 - 193	10^{-4} - 0.2	5
BF ₃	300	5 - 70	1
C _n H _{2n+2} (n = 2, 3, 4)	298 - 673	0.01 - 10	7
C(CH ₃) ₄	298 - 673	0.01 - 10	7
CClF ₃	300	2 - 30	3
CH ₂ Cl ₂	300	2 - 30	3
CH ₃ Cl	300	2 - 30	3
C ₂ H ₅ Cl	300	2 - 30	3
C ₄ H ₉ Br	300	2 - 30	3
He - F ₂ MIX	300	3 - 15	9
N ₂ - CO ₂ MIX	300	3 - 93	11
Xe - CF ₄ MIX	296	0.3 - 10	4
Xe-C ₂ H ₂ MIX	296	0.3 - 10	4
Xe-CF ₄ -C ₂ H ₂ MIX	296	0.3 - 10	4

* A substantial amount of electron swarm data has already been presented in previous volumes: see pages 717-732 of Vol. II and pages 2081-2108 of Vol. V.

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Table E-1.2. Sources of Electron Diffusion Data Not Presented in Previous Volumes.

ELECTRON DIFFUSION

Gas	Data	Approx. E/N (Td)	Ref.
He	$D_T/K, \alpha/N$	3 - 847	1
Ar	$D_T/K, \alpha/N$	8 - 1271	1
CO	D_T/K	5 - 350	3
NO	$D_T/K, \alpha/N$	56 - 1412	1
CO ₂	D_T/K	5 - 180	3
	$D_T/K,$	30 - 5000	2
CH ₄	$D_T/K, \alpha/N$	14 - 5650	1

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E-2. TRANSPORT PROPERTIES OF IONS

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Definitions and Relationships

$\vec{v}_d$ = Drift velocity of ion = average velocity of drift of ion along field lines in a gas exposed to a constant, uniform electric field $\vec{E}$. v_d is usually expressed in cm/sec.

K = Mobility of ion, defined by the equation $\vec{v}_d = K \vec{E}$. K is usually expressed in $\text{cm}^2/\text{V-sec}$.

K_o = Reduced mobility of ion = mobility of ion reduced to S.T.P., defined by the equation

$$K_o = \frac{p}{760} \frac{273.16}{T} K,$$

where p is the gas pressure in torr and T is the gas temperature in degrees Kelvin at which K was measured.

P_o = Reduced pressure = $\frac{273.16}{T} p$.

E/N = Ionic energy parameter = ratio of electric field intensity to gas number density. E/N is usually expressed in units of (volts/cm) / ($1/\text{cm}^3$) = $\text{V} - \text{cm}^2$.

$K_o(0)$ = Zero-field reduced mobility = K_o in the limit $E/N \rightarrow 0$.

Td = Unit of E/N , the "Townsend" = $10^{-17} \text{ V} - \text{cm}^2$.

$v_d = 0.0269 \cdot (E/N) \cdot K_o$, where v_d is in 10^4 cm/sec , E/N is in Td , and K_o is in $\text{cm}^2/\text{V} - \text{sec}$.

$$\vec{\vec{D}} = \begin{bmatrix} D_L & 0 & 0 \\ 0 & D_T & 0 \\ 0 & 0 & D_L \end{bmatrix} = \text{ionic diffusion tensor.}$$

D_L = (Scalar) longitudinal diffusion coefficient = coefficient of diffusion along electric field.

D_T = (Scalar) transverse diffusion coefficient = coefficient of diffusion transverse to electric field.

In the limit $E/N \rightarrow 0$, $D_L = D_T = D$, the scalar diffusion coefficient.

For a particular ionic species in a given gas at a given temperature, v_d , NK , ND_L , ND_T , and the average ionic energy are functions of E/N alone.

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- III. A summary of the data for the various ions and gases collected in Atomic Data and Nuclear Data Tables 17 and 22.

Table E-2.1. Sources of Ion Mobility Data Not Presented in Previous Volumes.*

<u>Ions in Helium</u>			
<u>Ion</u>	<u>Gas Temp (°K)</u>	<u>Approx. E/N (Td)</u>	<u>Ref(s).</u>
H ⁺	300	5 - 60	17
D ⁺	300	5 - 70	17
He ⁺	77 - 700	low-field	5, 28, 30
Li ⁺	20 - 500	low field	16
C ⁺	297	5 - 110	9
Na ⁺	90 - 480	low field	32
S ⁺	297	5 - 110	9
Ti ⁺	300	low field	18
Cd ⁺	526	low field	20
Cs ⁺	80 - 490	low field	32
Th ⁺	300	low field	18
Ne ⁺⁺	300	15 - 90	19
Ar ⁺⁺	305	10 - 100	19
Kr ⁺⁺	305	10 - 90	19
Xe ⁺⁺	302	10 - 90	19
He ₂ ⁺	293 120 - 700	5 - 40 low field	15 30
Ne ₂ ⁺	300 77 - 300	5 - 50 low field	3 8
HeNe ⁺	?	low field	33
(CH _n) ⁺ (n = 1, 2, 3, 4, 5)	300	5 - 100	36
SO ₂ F ₂ ⁻	300	low field	31
SO ₂ F ₂ ⁻ · (SO ₂)	300	low field	31
F ⁻ · (SF ₆)	300	low field	31

* A substantial amount of ion swarm data has already been presented in previous volumes: see pages 733-748 of Vol. II and pages 2109-2116 of Vol. V.

Table E-2.1 (continued)

ION MOBILITYIons in Neon

Ion	Gas Temp ($^{\circ}$ K)	Approx. E/N (Td)	Ref(s) .
He ⁺	77, 200, 300	low field	8
Ne ⁺	77, 200, 300	low field	5
Ar ⁺	77	10 - 20	15
Xe ⁺	303	20 - 140	19
Xe ⁺⁺	303	20 - 140	19
He ₂ ⁺	300	5 - 50	3
Ne ₂ ⁺	77 77, 200, 300	10 - 55 low field	15 5
N ₂ ⁺	300	low field	24

Ions in Argon

Ar ⁺	77 77, 200, 300	15 - 90 low field	15 5
K ⁺	291, 400, 460	low field	16
Ar ⁺⁺	77, 200, 300	low field	5
H ₃ ⁺	300	10 - 110	25
ArH ⁺	300	30 - 110	25
Ar ₂ ⁺	77	50 - 100	15
H ₃ O ⁺ · (H ₂ O) _n n = 1, 2, 3	337	low field	38
ReO ₃ ⁻	295	15 - 200	4
ReO ₄ ⁻	295	15 - 200	4
(WO ₃) _n ⁻ n = 1, 2, 3	295	25 - 160	4
H ₃ O ⁺	337	30 - 110	38

Table E-2.1 (continued)

ION MOBILITYIons in Krypton

<u>Ion</u>	<u>Gas Temp ($^{\circ}$K)</u>	<u>Approx. E/N (Td)</u>	<u>Ref(s)</u>
Li^+	291	low field	26, 37
Na^+	300	5 - 500	35
Rb^+	195 - 455	low field	16
Kr_2^+	295	5 - 170	15

Ions in Xenon

Li^+	291	low field	26, 37
Na^+	300	5 - 500	35
Cs^+	200 - 450	low field	16
Xe_2^+	300	40 - 230	2
	200 - 300	low field	14
Xe_3^+	200 - 300	low field	14

Ions in Hydrogen

H_3^+	77, 195, 300	low field	6
H_3O^+	300	5 - 140	6, 13

Table E-2.1 (continued)

ION MOBILITYIons in Nitrogen

Ion	Gas Temp (°K)	Approx. E/N (Td)	Ref(s).
Na^+	291	low field	26, 37
O_2^+	300 - 640	low field	10
NO^+	300 - 640	low field	10
Cl^-	340 - 470	low field	12
Cl_2^-	300 - 470	low field	12
NO_2^-	215 - 675	low field	11
NO_3^-	215 - 675	low field	11
CO_3^-	215 - 675	low field	11

Ions in Oxygen

Na^+	304	5 - 500	34
Cl^-	330 - 470	low field	12
Cl_2^-	300 - 470	low field	12
O_4^-	300	low field	27
CO_3^-	300 - 470	low field	12
SF_5^-	300	low field	31
SF_6^-	300	low field	31

Ions in Carbon Dioxide

Li^+	?	35 - 350	21
	307	25 - 900	34
O_2^+	?	40 - 400	21
NO^+	?	30 - 400	21

Table E-2.1 (concluded)

ION MOBILITYIons in Methane

Ion	Gas Temp ($^{\circ}$ K)	Approx. E/N (Td)	Ref(s).
Li^+	304	10 - 600	34
SF_5^-	300	low field	31
SF_6^-	300	low field	31

Ions in Sulfur Hexafluoride

SF_5^-	300	5 - 140	29
SF_6^-	300	5 - 140	29
$\text{SF}_6^- \cdot (\text{SF}_6)$	300	20 - 120	29
$\text{SF}_6^- \cdot (\text{SF}_6)_2$	300	20 - 120	29

Ions in Metal Vapors

Ion - Gas	Gas Temp ($^{\circ}$ K)	Approx. E/N (Td)	Ref(s).
$\text{Rb}^+ - \text{Rb}$	621	85 - 340	23
$\text{Rb}_2^+ - \text{Rb}$	621	85 - 340	23
$\text{Cs}^+ - \text{Cs}$	580 - 650	20 - 480	7, 23
$\text{Cs}_2^+ - \text{Cs}$	580 - 650	60 - 480	7, 23
$\text{Hg}^+ - \text{Hg}$	350 500	low field 210 - 570	1 22
Hg_2^+	500	110 - 325	22

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ION MOBILITY

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Table E-2.2. Sources of Ion Diffusion Data Not Presented in Previous Volumes.

ION DIFFUSION				
Ion - Gas	Measured Quantity	Gas Temp (°K)	Approx. E/N (Td)	Ref.
He ⁺ - He	D	293	low field	2
He ₂ ⁺ - He	D	293	low field	2
(CH _n) ⁺ - He n = 0,1,2,3,4,5	D _L	300	5 - 100	4
Na ⁺ - Kr	D _L	300	5 - 500	3
Na ⁺ - Xe	D _L	300	5 - 400	3
N ₃ ⁺ - N ₂	D _T	300	10 - 100	1
N ₄ ⁺ - N ₂	D _T	300	10 - 100	1
O ₂ ⁺ - O ₂	D _T	300	10 - 500	1
CO ⁺ ·CO - CO	D _T	300	10 - 130	1

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Table 2.3. References to Ion Transport Data Appearing in Compilations
I and II. [Atomic Data and Nuclear Data Tables 17, 177
(1976); 22, 179 (1978)].*

Mobilities (K)

Ar ⁺ in Ar. I - pg. 196	Cs ⁺ in Ar. II - pg. 193	Hg ⁺ in Ne. I - pg. 195
in He. I - pg. 185	in CO. II - pg. 204	H ₂ O ⁺ in He. II - pg. 188
Ar ²⁺ in Ar. I - pg. 196	in CO ₂ . II - pg. 204	H ₃ O ⁺ in He. II - pg. 188
Ar ²⁺ (¹ F) in Ar. II - pg. 195	in H ₂ . II - pg. 200	in N ₂ . II - pg. 201
Ar ²⁺ (³ F) in Ar. II - pg. 195	in He. II - pg. 187	H ₃ O ⁺ ·H ₂ O in He. II - pg. 188
Ar ₂ ⁺ in Ar. I - pg. 198	in Kr. II - pg. 196	in N ₂ . II - pg. 201
ArH ⁺ in He. I - pg. 187	in N ₂ . II - pg. 201	H ₃ O ⁺ ·2H ₂ O in He. II - pg. 188
Br ⁻ in Ar. II - pg. 194	in Ne. II - pg. 191	in N ₂ . II - pg. 201
in He. II - pg. 190	in O ₂ . II - pg. 203	I ⁻ in Ar. II - pg. 194
C ⁺ in CO. I - pg. 208	in Xe. II - pg. 198	in He. II - pg. 190
CH ₅ ⁺ in He. I - pg. 190	D ⁺ in D ₂ . I - pg. 202	K ⁺ in Ar. I - pg. 197
C ₂ H ₂ ⁻ in He. I - pg. 192	in He. I - pg. 184	in CH ₄ . II - pg. 204
CH ₃ O ₂ ⁺ in He. I - pg. 190	in Ne. I - pg. 195	in CO. I - pg. 207
Cl ⁻ in Ar. II (2) - pg. 194	D ₃ ⁺ in D ₂ . I - pg. 202	in CO ₂ . I - pg. 209
in He. I - 190, II - 190	F ⁻ in Ar. II - pg. 194	in D ₂ . I - pg. 203
in Kr. II - pg. 196	in He. II - pg. 190	in H ₂ . I - pg. 202
in Ne. II - pg. 192	in Kr. II - pg. 196	in He. I - pg. 183
in Xe. II - pg. 198	in Xe. II - pg. 198	in N ₂ . I - pg. 204
CO ⁺ in CO. I - pg. 207	H ⁺ in H ₂ . I - pg. 200	in NO. I - pg. 208
in He. I - pg. 186	in He. I - pg. 184	in Ne. I - pg. 195
CO ₂ ⁺ in Ar. I - pg. 198	in Ne. I - pg. 194	in O ₂ . I - pg. 205
in He. I - pg. 188	H ⁻ in H ₂ . I - pg. 200	in Xe. II - pg. 197
in N ₂ . I - 204, II (2) 203	in He. I - pg. 190	Kr ⁺ in Ar. II - pg. 193
in Ne. II - pg. 193	H ₂ ⁺ in He. I - pg. 186	in He. II - pg. 187
CO ₃ ⁻ in Ar. I - pg. 199	H ₃ ⁺ in H ₂ . I - pg. 200	in Kr. I - pg. 199, II-195
in CO ₂ . I - pg. 209	in He. I - pg. 187	Kr ⁺ in Kr. II - pg. 196, 197
in He. I - pg. 192	He ⁺ in He. I - pg. 184	Kr ⁺ in N ₂ . II - pg. 202
in O ₂ . I - pg. 206	He ²⁺ in He. II - pg. 189	Kr ₂ ⁺ in Kr. I - pg. 200
CO ₄ ⁻ in O ₂ . I - pg. 207	He ₂ ⁺ in He. I - pg. 186	Kr ²⁺ (A) in Kr. II - pg. 197
C ₂ O ₂ ⁺ in CO. I - pg. 207	HeH ⁺ in He. II - pg. 189	Kr ²⁺ (E) in Kr. II - pg. 197
COH ⁺ in Ar. I - pg. 198	Hg ⁺ in Ar. I - pg. 197	Li ⁺ in Ar. I - pg. 196
in He. I - pg. 188	in He. I - pg. 185	in D ₂ . I - pg. 202

Note: * I refers to Ellis, et al., ADNDT 17, 177 (1976); II to Ellis, et. al., ADNDT 22, 179 (1978).

Table E-2.3 (continued)

Mobilities (K) (continued)

Li^+ in H_2 . I - pg. 201	NH_4^+ in He. I - pg. 189	O_3^- in He. I - pg. 191
in He. I - pg. 183	in N_2 . I - pg. 205	in O_2 . I - pg. 206
in N_2 . II - pg. 200	N_2H^+ in N_2 . I- 203, II-202	O_4^+ in O_2 . I - pg. 205
in Ne. I - pg. 194	NO^+ in He. I - pg. 187	OH^- in He. I - pg. 191
in O_2 . II - pg. 203	in NO. I - pg. 208	O_2H^+ in He. I - pg. 188
N^+ in N_2 . I - pg. 204	$\text{NO}^+\cdot\text{H}_2\text{O}$ in He. I - pg. 189	O_2H_2^+ in He. I - pg. 189
in He. I - pg. 184	N_2O^+ in Ar. II - pg. 193	Rb^+ in Ar. I - pg. 197
N_2^+ in He. I - pg. 186	in He. II - pg. 189	in CO_2 . I - pg. 208
in N_2 . I - pg. 202	N_2O^+ in N_2 . II - pg. 202	in H_2 . II - pg. 200
N_3^+ in N_2 . I - pg. 204	in Ne. II - pg. 191	in He. I - pg. 83
N_4^+ in N_2 . II - pg. 202	NO_2^- in He. I - pg. 192	in Kr. II - pg. 195
Na^+ in Ar. I - pg. 196	N_2O_2^+ in NO. I - pg. 201	in N_2 . II - pg. 200
in CO_2 . I - pg. 209	N_2OH^+ in Ar. I - pg. 199	in Ne. I - pg. 194
in D_2 . I - pg. 202	in He. I - pg. 189	in O_2 . II - pg. 203
in H_2 . I - pg. 201	O^+ in Ar. I - pg. 197	in Xe. II - pg. 198
in He. I - pg. 183	in He. I - pg. 185	SF_5^- in He. I - pg. 193
in Ne. I - pg. 194	O^- in CO_2 . I - pg. 209	SF_6^- in He. I - pg. 193
Ne^+ in He. II - pg. 187	in He. I - pg. 191	SO_3^- in He. I - pg. 193
in Ne. I - pg. 195	in O_2 . I - pg. 206	SO_2F^- in He. I - pg. 192
$\text{Ne}^+(^2\text{P}_{3/2})$ in Ne. II - pg. 191	O_2^+ in Ar. I - pg. 198	U^+ in He. I - pg. 185
$\text{Ne}^+(^2\text{P}_{3/2})$ in Ne. II - 191	in He. I - pg. 187	Xe^+ in He. II - pg. 187
$\text{Ne}^{2+}(^1\text{D})$ in Ne. II - pg. 192	in O_2 . I - pg. 205	$\text{Xe}^+(^2\text{P}_{1/2})$ in Xe. II - pg. 199
$\text{Ne}^{2+}(^1\text{S})$ in Ne. II - pg. 192	O_2^- in He. I - pg. 190	$\text{Xe}^+(^2\text{P}_{3/2})$ in Xe. II - pg. 199
$\text{Ne}^{2+}(^3\text{P})$ in Ne. II - pg. 192	in O_2 . I - pg. 206	$\text{Xe}^{2+}(\text{A})$ in Xe. II - pg. 199
Ne_2^+ in He. I - pg. 193	O_3^- in Ar. I - pg. 199	$\text{Xe}^{2+}(\text{B})$ in Xe. II - pg. 199
NH_3^+ in He. I - pg. 188		

Longitudinal Diffusion Coefficients (D_L)

Cl^- in Ar. II - pg. 208	Cs^+ in CO_2 . II - pg. 215	D^+ in D_2 . II - pg. 211
in Kr. II - pg. 209	in H_2 . II - pg. 211	D^- in D_2 . II - pg. 212
in Ne. II - pg. 207	in He. II - pg. 206	D_3^+ in D_2 . II - pg. 212
in Xe. II - pg. 210	in Kr. II - pg. 209	F^- in Kr. II - pg. 209
CO^+ in CO . II - pg. 214	in N_2 . II - pg. 212	in Xe. II - pg. 210
Cs^+ in Ar. II - pg. 208	in Ne. II - pg. 207	H^+ in H_2 . II - pg. 210
in CO . II - pg. 214	in O_2 . II - pg. 213	H^- in H_2 . II - pg. 211
	in Xe. II - pg. 210	

Table E-2.3 (concluded)

Longitudinal Diffusion Coefficients (D_L) (continued)

H_3^+ in H_2 II - pg. 211	Li^+ in Ar. II - pg. 207	O^- in O_2 . IV - pg. 213
K^+ in Ar. II - pg. 208	in H_2 . II - pg. 210	O_2^+ in O_2 . II - pg. 213
in CO . II - pg. 214	in He. II - pg. 206	O_2^- in O_2 . II - pg. 214
in CO_2 . II - pg. 215	in Ne. II - pg. 206	Rb^+ in Ar. II - pg. 208
in H_2 . II - pg. 211	N^+ in N_2 . II - pg. 212	in CO_2 . II - pg. 215
in He. II - pg. 206	N_2^+ in N_2 . II - pg. 213	in H_2 . II - pg. 211
in Kr. II - pg. 208	Na^+ in Ar. II - pg. 208	in He. II - pg. 206
in N_2 . II - pg. 212	in CO_2 . II - pg. 215	in Kr. II - pg. 209
in Ne. II - pg. 207	in H_2 . II - pg. 210	in N_2 . II - pg. 212
in NO . II - pg. 214	in He. II - pg. 206	in Ne. II - pg. 207
in O_2 . II - pg. 213	in Ne. II - pg. 207	in O_2 . II - pg. 213
in Xe. II - pg. 209	NO^+ in NO . II - pg. 214	in Xe. II - pg. 209

Transverse Diffusion Coefficients (D_T)

H_3^+ in H_2 . II - pg. 216	K^+ in N_2 . II - pg. 216	N_2^+ in N_2 . II - pg. 216
K^+ in H_2 . II - pg. 216	N^+ in N_2 . II - pg. 216	O_2^+ in O_2 . II - pg. 216

Update of References in Compilation II [ADNDT 22, 179 (1978)]

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References for Table II

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F. INTERACTIONS WITH STATIC ELECTRIC AND MAGNETIC FIELDS

(No new entries here. See Vol. II for data.)

G. PARTICLE PENETRATION IN GASES (IONS, NEUTRALS, AND ELECTRONS)

(No new entries here. See Vols. II and V for data.)

H. PARTICLE AND PHOTON INTERACTIONS WITH SOLIDS

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(No new entries here.)	
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(No new entries here.)	
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(No new entries here).	

H-1. SPUTTERING

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INTRODUCTION

Sputtering Yield. Is defined as the number of target atoms removed per projectile atom incident. It is measured as a function of incident projectile energy and as a function of projectile incidence angle on the surface. Unless otherwise indicated the yields presented are for normal incidence.

Angular Dependence. Yield increases rapidly with increasing incidence angle (measured between projectile impact direction and target surface normal). For heavy particles (mass > 4 amu) the yield Y increases as

$$Y(\theta) = Y(0) \frac{1}{\cos \theta}$$

where θ is incidence angle and $Y(0)$ is the yield at normal incidence ($\theta=0^\circ$).

State of the Sputtered Species. From a monatomic target the ejected particles are generally neutral atoms with small fractions of ions and multimers (10^{-2} to 10^{-3}) of each species may be excited. While there is much information on charge and quantum states this is poorly digested and is not reproduced here.

Energy of Sputtered Species. For heavy particle impact where ejection results from a collision cascade the yield of sputtered particles $Y(E)$ as a function of ejected particle energy E often follows the equation

$$Y(E) \propto \frac{1}{(E+E_B)^2} \cdot \frac{1}{(1+E/E_B)}$$

Here E_B is the binding energy of the ejected atom while it was in the lattice, and may be approximated by sublimation energy found in many standard reference books.

Topographical Changes. It is well known that sputtering of a surface cause changes to topography resulting in such features as cones, ripples and faceted pits. Cones are due sometimes to impurity particles having a low sputtering coefficient which protects the underlying substrate. These various features are also often associated with preferential etching of different crystallographic faces of polycrystalline materials. Selective etching at grain boundaries and defects is also known.

Sputtering of Alloys and Compounds. The data we present are for monatomic (generally metallic) polycrystalline materials. In the sputtering of alloys or compounds one anticipates first a depletion of the component(s) exhibiting the highest sputtering yield leading to a changed surface composition. Thus yield measured by weight loss is a function of projectile dose. Also for compounds there may be a very high yield of molecules. Thus one cannot necessarily estimate yield from an alloy or compound by some average of the yields measured for the constituents in monatomic targets.

Effects of Ambient Atmosphere. An ambient atmosphere of reactive species (e.g. oxygen or hydrocarbons) produces substantial changes to the fluxes of sputtered ions and sputtered excited species. However, such an atmosphere does not appear to substantially change the flux of sputtered atoms which represents the bulk of ejected material. Thus the sputtering yields are not greatly influenced by the ambient atmosphere. Non reactive gases do not appear to

influence yield nor distribution of quantum states.

Semi-empirical Formulae to Estimate Yield. Sputtering yield is related to energy transfer to the lattice which causes atoms from the surface to acquire sufficient energy to overcome their binding to the solid and thereby escape. Thus yield should be related to stopping power, particle masses, projectile impact energy and target binding energy. A number of semi-empirical formulations have been devised to represent sputtering yield. Not only are they more convenient than raw data for modelling purposes but they also permit estimates for projectile target combinations that have not yet been subjected to experimental study. We discuss such formulae in sections B-1.13 and B-1.32.

Data Presented Here. Due to space limitations we present only a small fraction of the available sputtering yield data. Further data in graphical and tabular form can be found in the reviews listed below. Also the semi-empirical formulations given in sections B-1.13 and B-1.32 can be utilized.

Reviews and Data Compilations.

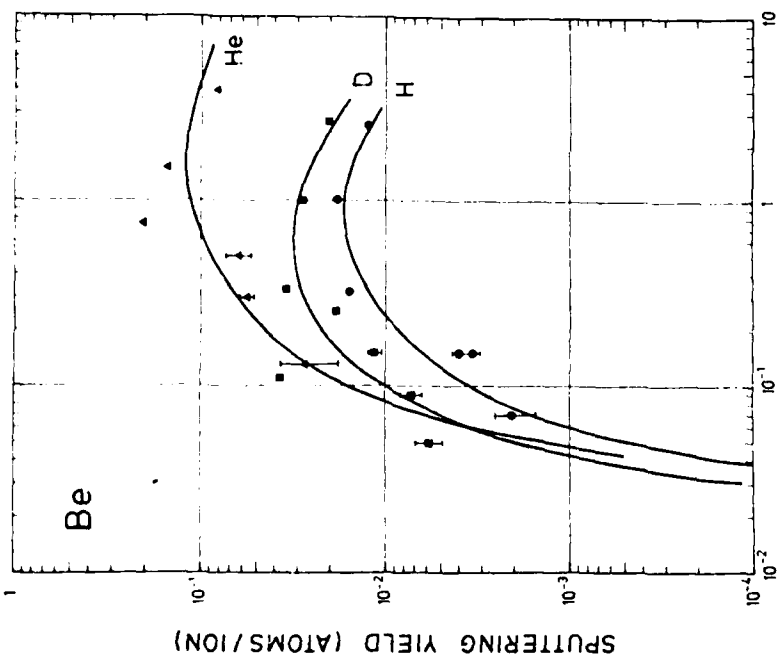
N. Matsunami, Y. Y. Yamamura, Y. Itikawa, N. Itoh, Y. Kazumata, S. Miyagawa, E. Morita and E. Shimizu. Energy Dependence of Sputtering Yields of Monatomic Solids. Institute of Plasma Physics, Nagoya University Nagoya Japan. Report No., IPPJ-MI-14 (June 1980).

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R. Behrisch (Ed). Sputtering by Ion Bombardment.. To be published as a book in the series "Topics of Applied Physics" by Springer-Verlag, Heidelberg.

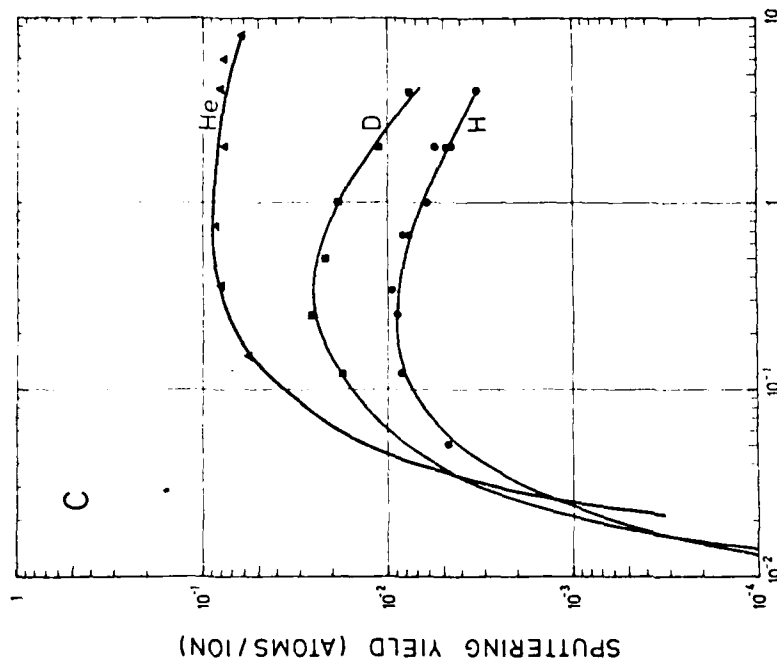
A. R. Krauss and R. B. Wright. J. Nucl. Mat. 89, 229 (1980). (A bibliography of data on kinetic energy and mass distribution of sputtered particles).



Graphical Data H-1.1

Energy dependence of the sputtering yield of Be with H, D, ^4He . The solid curves are a fit calculated from Eq. 1 of Table H-1.13

Reference: 162

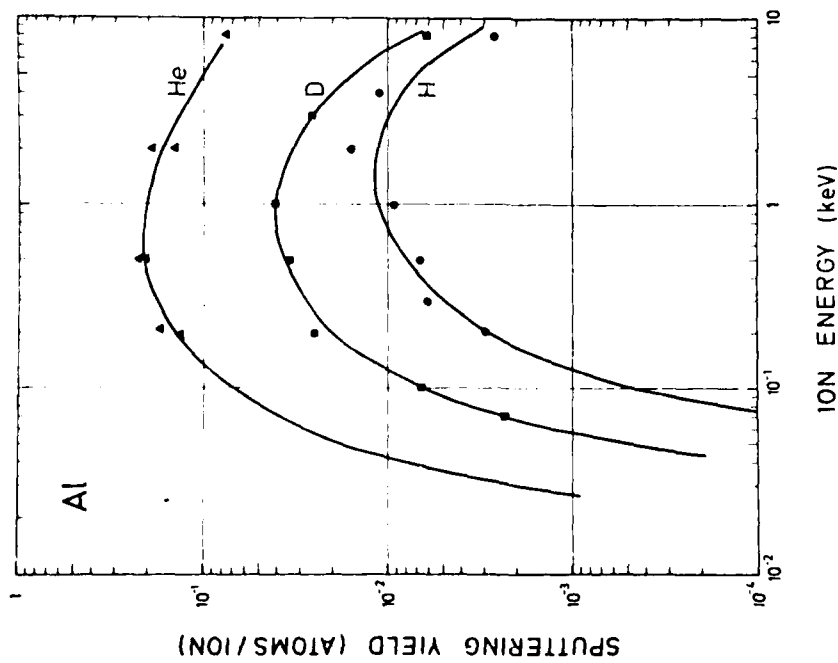


ION ENERGY (keV)

Graphical Data H-1.2

Energy dependence of the sputtering yield of C with H, D, He. The sputtering yields for different kinds of graphite vary widely according to their structure and orientation. Here only data for pyrolytic carbon (Union Graphite) are reported. For sputtering with D and H, an increase of the sputtering yield at elevated temperatures ($\sim 500^\circ\text{C}$) due to chemical effects has been observed. The solid curves are a fit calculated from Eq. 1. of Table H-1.13

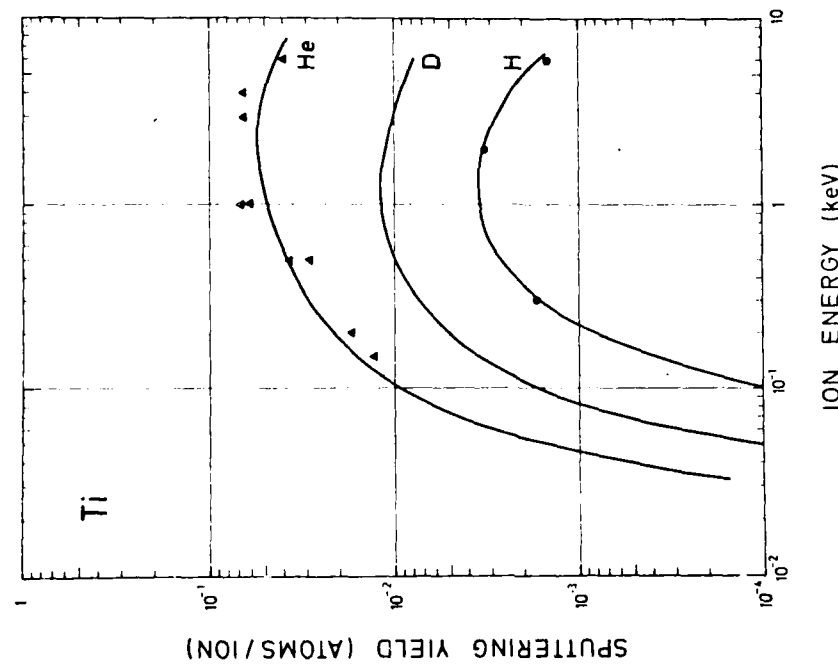
References: 1,3,465.



Graphical Data H-1.3

Energy dependence of the sputtering yield of Al with H, D, He. The solid curves are a fit calculated from Eq. 1 Table H-1.13

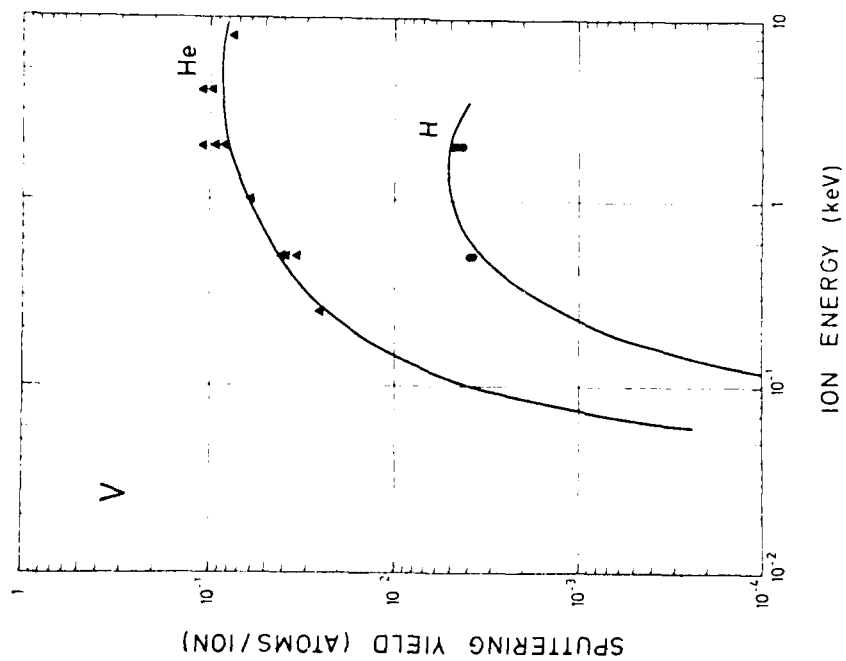
Reference: 1.6



Graphical Data H-1.4

Energy dependence of the sputtering yield of Ti with H and ^4He . The solid curves are a fit calculated from Eq. 1 of Table H-1.13; the line for D is interpolated by the same formulation.

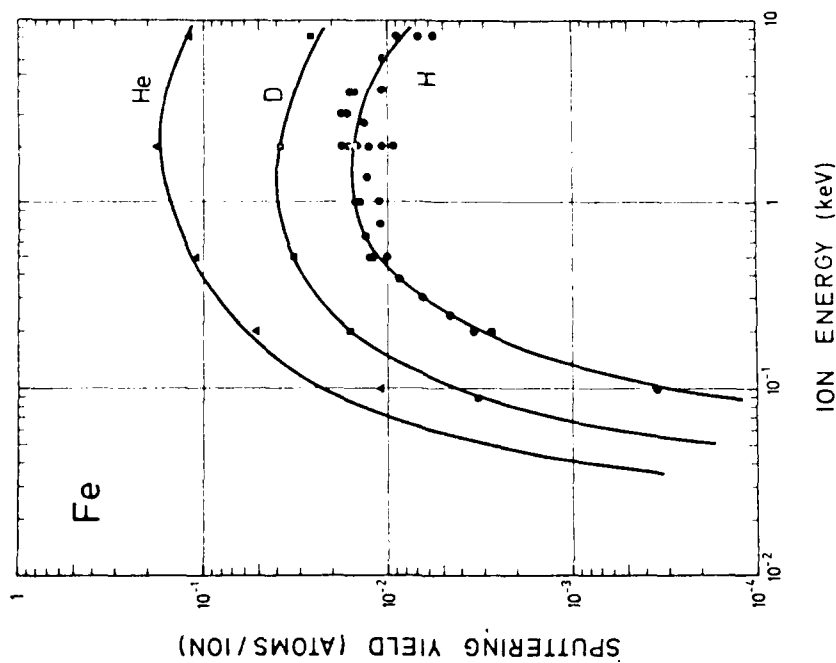
Reference: 1,567.



Graphical Data H-1.5

Energy dependence of the sputtering yield of V with H and ^4He . The solid curves are a fit calculated from Eq. 1 of Table H-1.13.

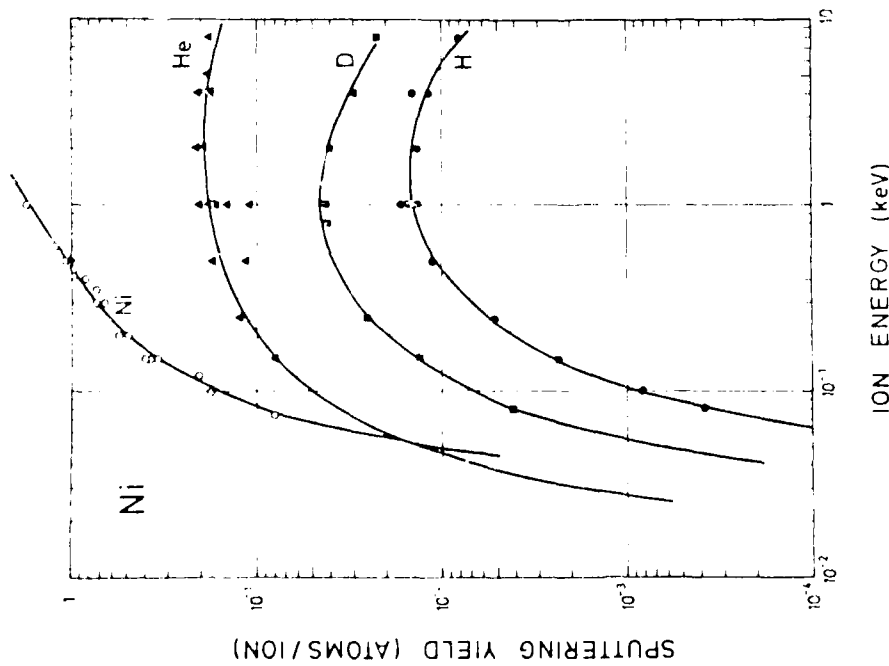
Reference: 1



Graphical Data H-1.6

Energy dependence of the sputtering yield of Fe with H, D and He. The solid curves are a fit calculated from Eq. 1 of Table H-1.13

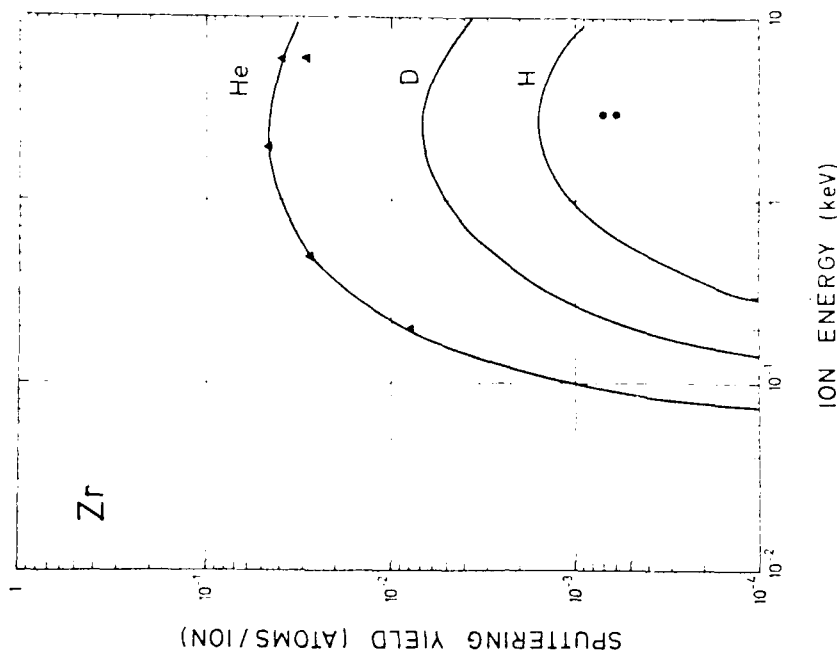
Reference: 1.



Graphical Data H-1.7

Energy dependence of the sputtering yield of Ni with H, D, He and Ni. The solid curves are a fit calculated from Eq. 1 of Table H-1.13

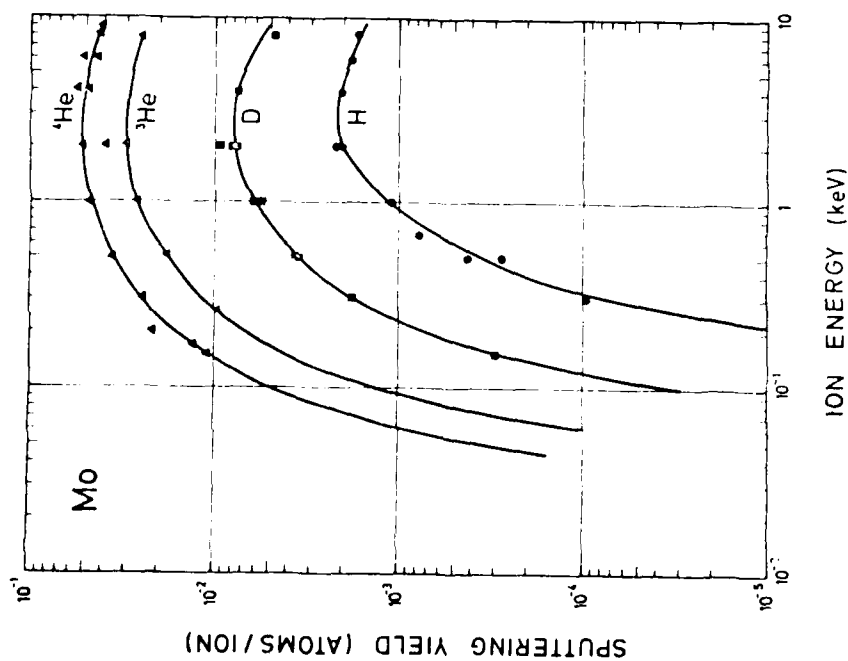
Reference: 1,8,9



Graphical Data H-1.8

Energy dependence of the sputtering yield of Zr with H and He. The solid curves are a fit calculated from Eq. 1 of Table H-1.13

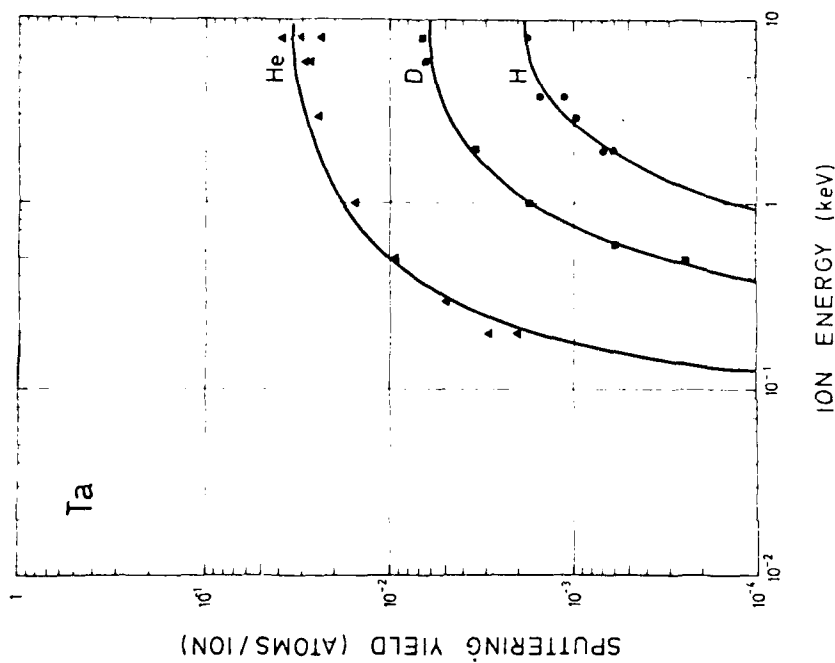
Reference: 1,10.



Graphical Data H-1.9

Energy dependence of the sputtering yield of Mo with H , D , ^3He , ^4He . The solid curves are a fit calculated from Eq. 1 of Table H-1.13

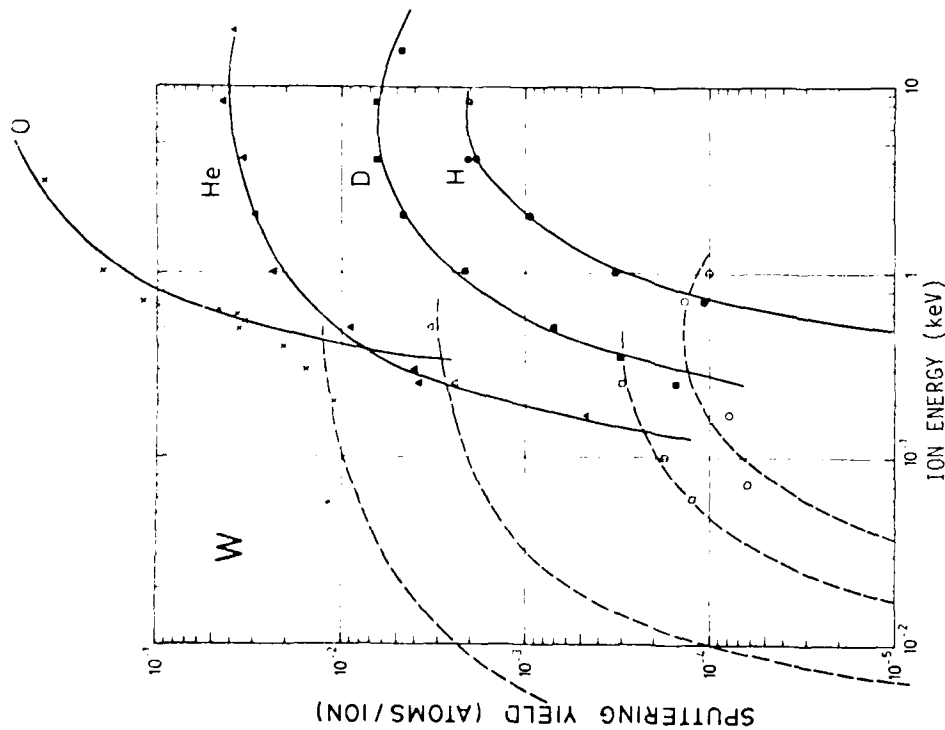
Reference: 1,11



Graphical Data H-1.10

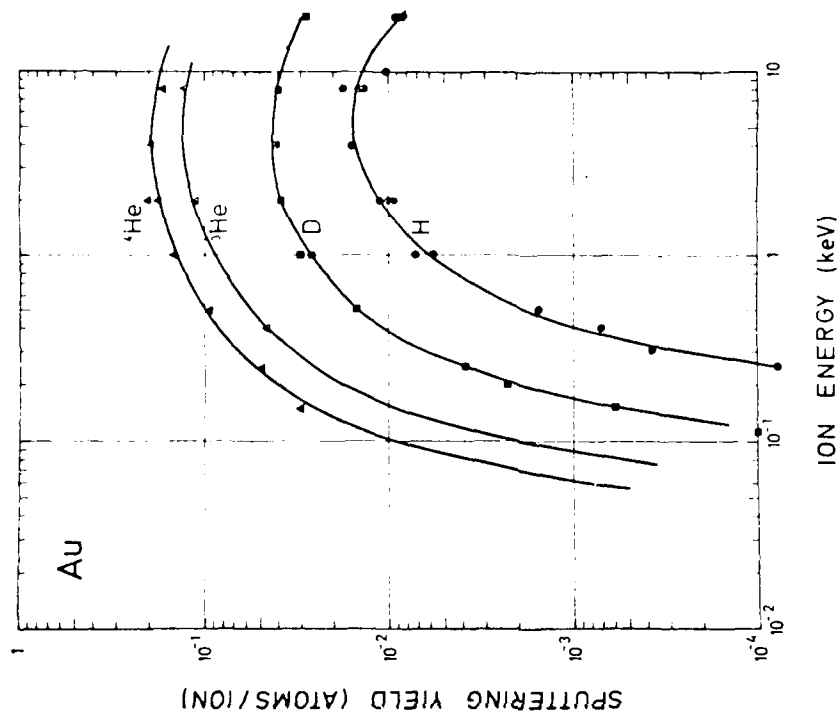
Energy dependence of the sputtering yield of Ta with H , D , and ^4He . The solid curves are a fit calculated from Eq. 1 of Table H-1.13.

Reference: 1,5



Graphical Data H-1.11

Energy dependence of the sputtering yield of W with H, D, He and O. The dashed curves indicate an additional sputtering mechanism at low energies in a background pressure of $8 \cdot 10^{-5}$ Torr O_2 . This mechanism might be due to sputtering of tungsten oxide molecules with lower surface binding energy than tungsten atoms. The solid and dashed curves are a fit calculated from Eq. 1 of Table H-1.13.
Reference: 1,5



Graphical Data H-1.12

Energy dependence of the sputtering yield of Au with H, D, ^3He and ^4He . The solid curves are a fit calculated from Eq. 1 of Table H-1.13.

Reference: 1,11.

Semi-Empirical Formulation for Sputtering

Yields Due to Light Particle Impact

<u>Symbols</u>	M_1	projectile mass (amu)
	M_2	target atom mass (amu)
	E	projectile energy (eV)
	E_{th}	threshold energy of the sputtering process (eV)
	E_B	surface binding energy of the target atom (eV).

Yield Expression

$$Y = 6.4 \cdot 10^{-3} M_2 \gamma^{5/3} \left(\frac{E}{E_{th}} \right)^{1/4} \left(1 - \frac{1}{E/E_{th}} \right)^{7/2} \text{ atoms/ion}$$

where

$$\gamma = \frac{4 M_1 M_2}{(M_1 + M_2)^2} \quad (1)$$

E_{th} - take from empirical values tabulated on the left or see note 4 below

Limitations

$$M_1/M_2 < 0.4$$

$$\frac{E}{E_{th}} < 20$$

Accuracy

Represents empirical data to within $\pm 25\%$.

Source

J. Roth, J. Bohdanský, W. Oltenberger
Data on Light Ion Sputtering, Max-Planck
Institut für Plasmaphysik, Garching May
1979 (Unpublished report).

H. L. Bay, J. Roth, J. Bohdanský. J. Appl.
Phys. 48, 4722 (1977).

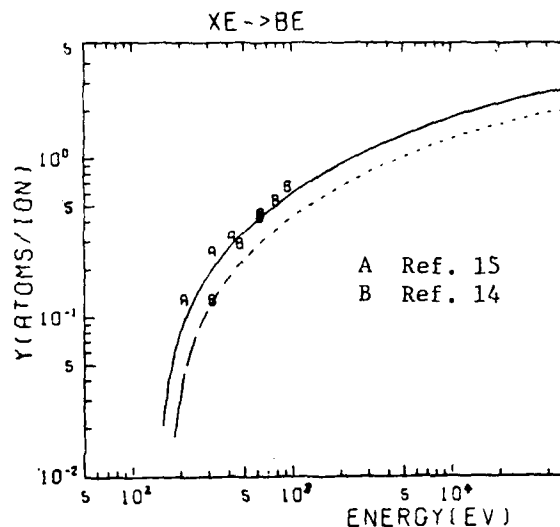
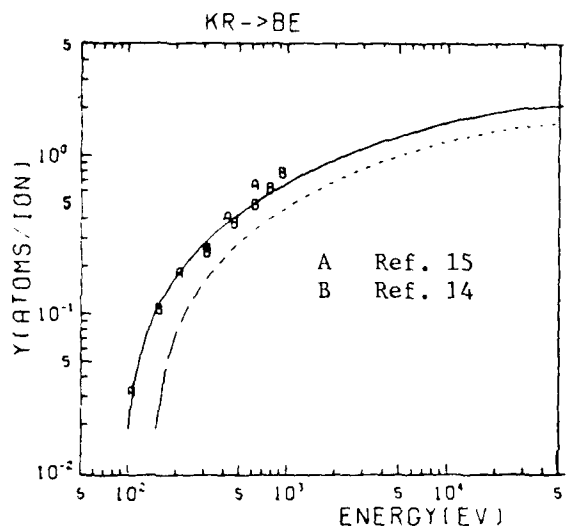
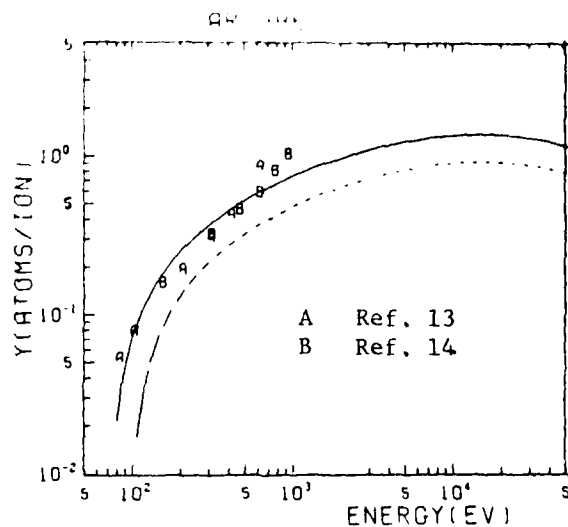
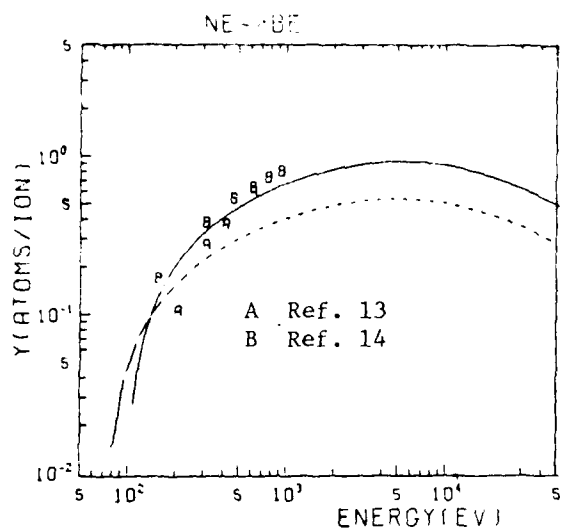
Threshold energy in eV

Ion \ Target	H	D	He ³	He ⁴
Al	53	34		20.5
Au	184	94	60	44
Be	27.5	24		33
C	9.9	11		16
Fe	64	40		35
Mo	164	86	45	39
Ni	47	32.5		20
Si	24.5	17.5		14
Ta	460	235		100
Ti	43.5			22
V	76			27
W	400	175		100
Zr				60

J. Bohdanský, J. Roth, M. L. Bay (to be published).

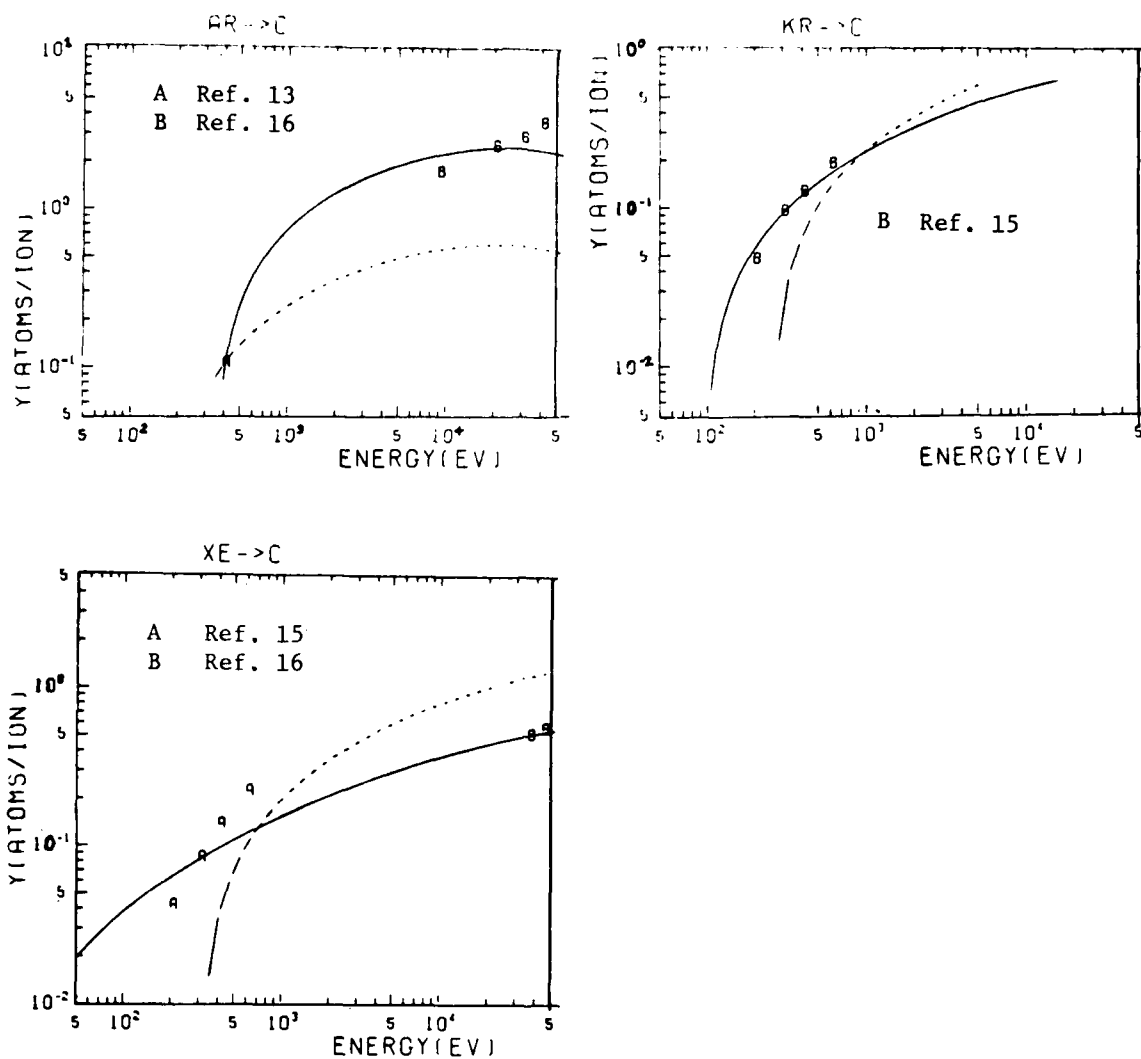
Notes

- 1) While this formulation has been developed primarily to represent yields for light ion impact it is also successful for sputtering of Ni by heavy ions within the limitations stated above.
- 2) The formulation is particularly accurate close to threshold.
- 3) At energies in excess of the range stated under "limitations" the empirical data scatters above the values given by the formulation.
- 4) Values of E_{th} for use in this formula should be taken from the table given above; these are values obtained by fitting to the data. For other cases one could estimate E_{th} as $E_B/\gamma(1-\gamma)$ where E_B is the binding energy. In turn E_B can be equated to the sublimation energy (e.g. JANAF Thermo-Chemical Tables, ed. D. R. Stull, H. Prophet, NSRDS-NBS 37), expressed in eV.



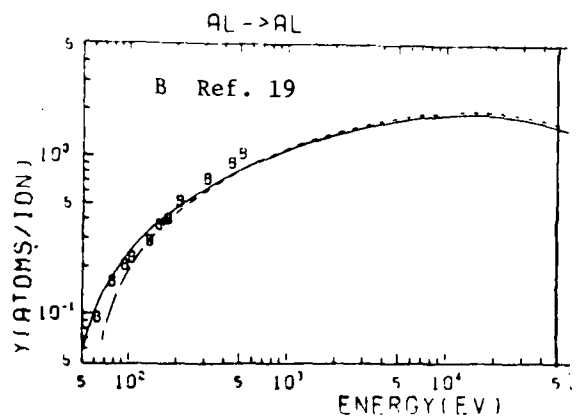
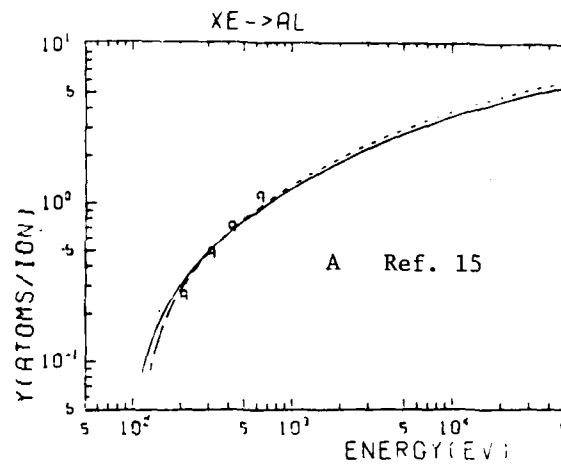
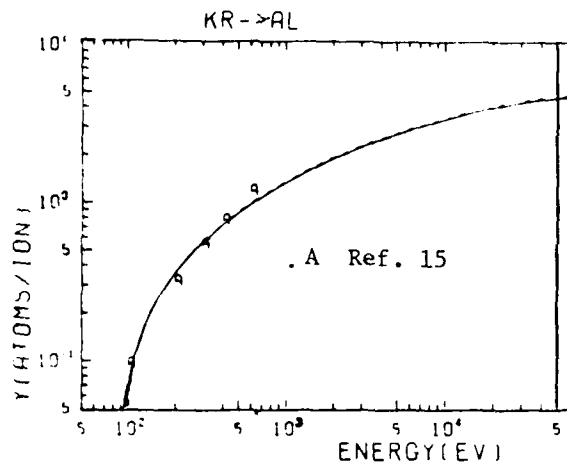
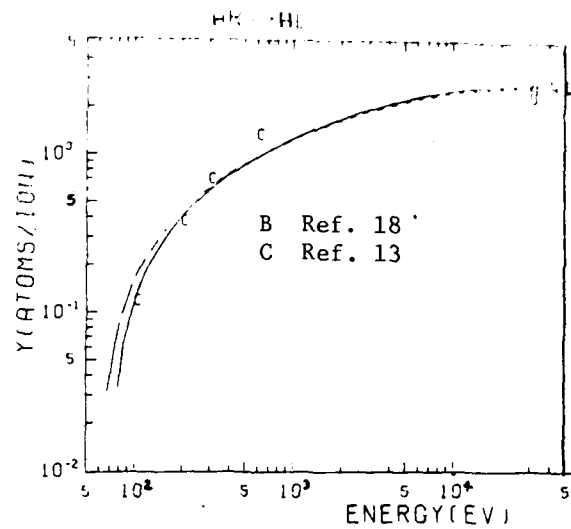
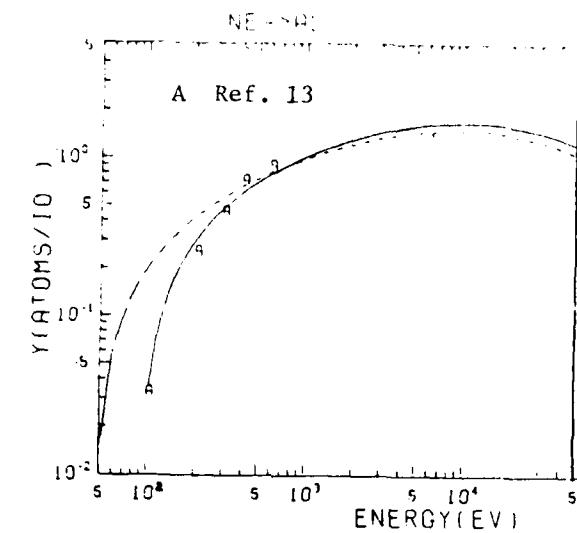
Graphical Data H-1.14

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ & Xe^+ impact on Be. The solid line is a best fit to the data by Eq. 1 of H-1.32. From Ref. 12.



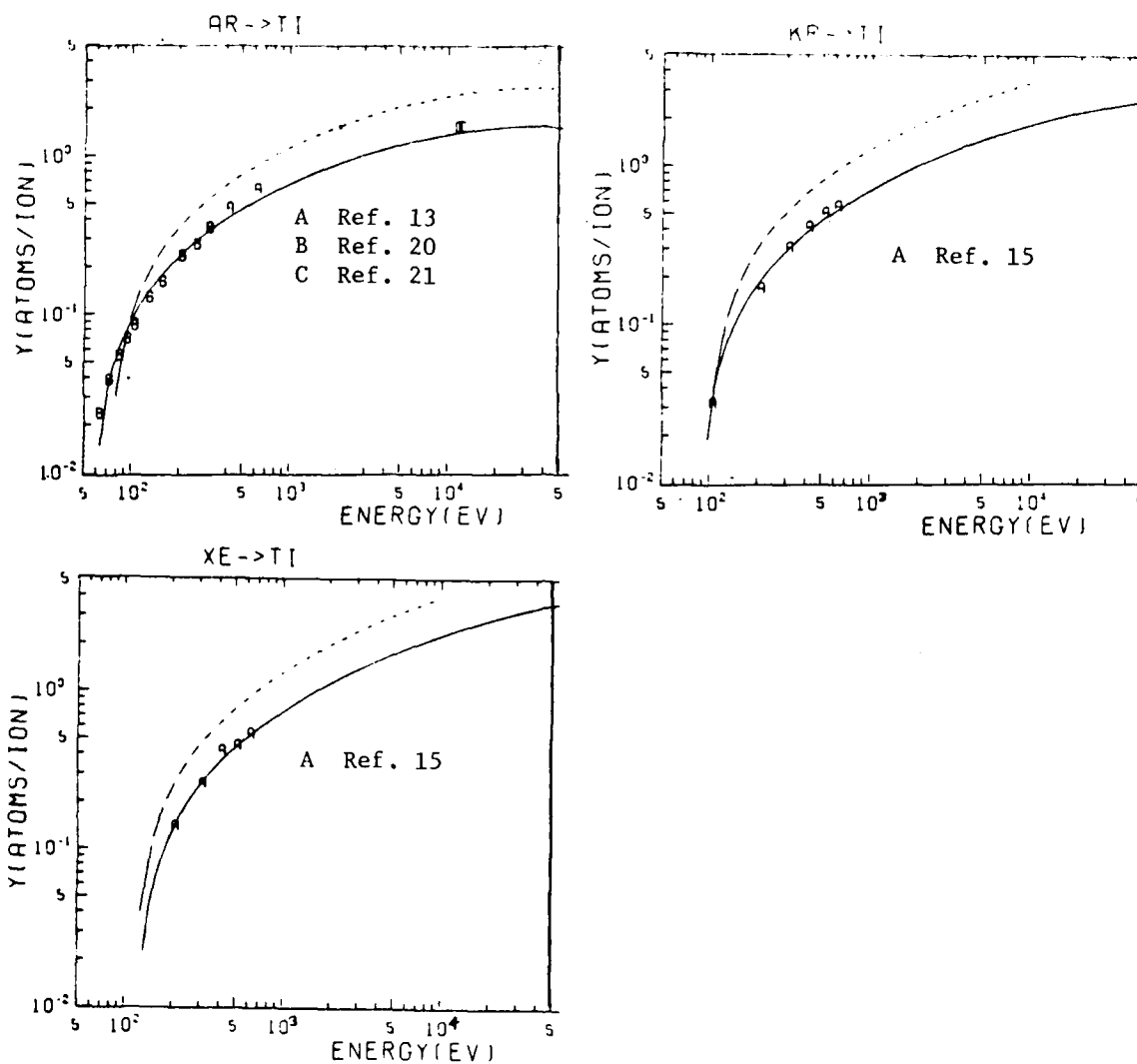
Graphical Data H-1.15

Energy dependence of the sputtering yield for Ar^+ , Kr^+ and Xe^+ impact on C. The solid line is a best fit of Eq. 1 (section H-1.32) to the data. From Ref. 12.



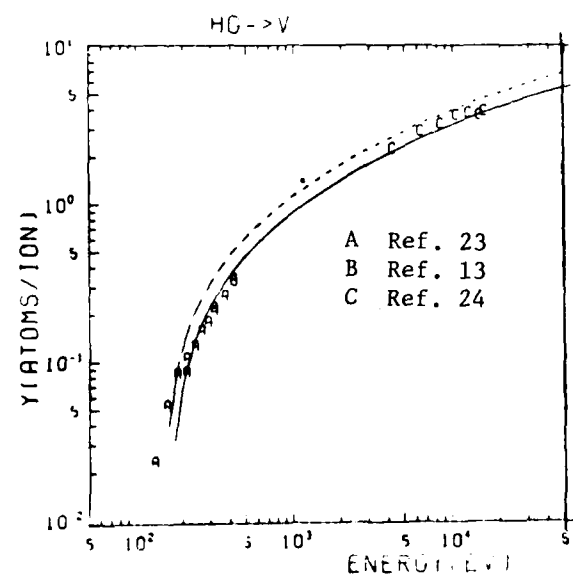
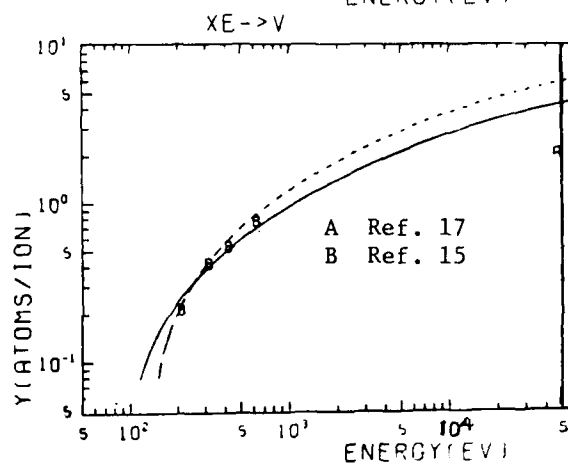
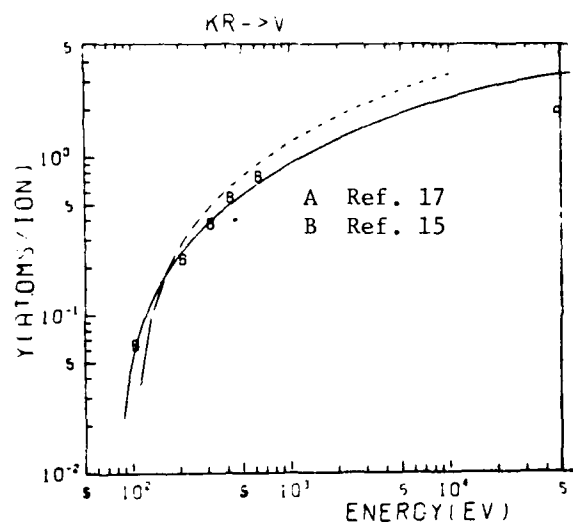
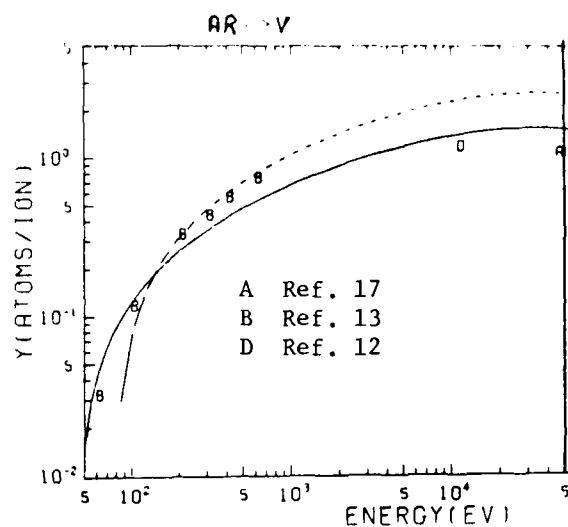
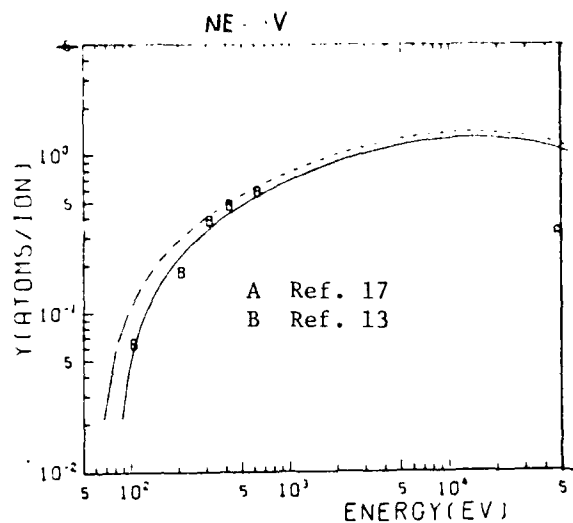
Graphical Data H-1.16

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ , Xe^+ and Al^+ on Al. The solid line is a best fit of Eq. 1 (section H-1.32) to the data. From Ref. 12.



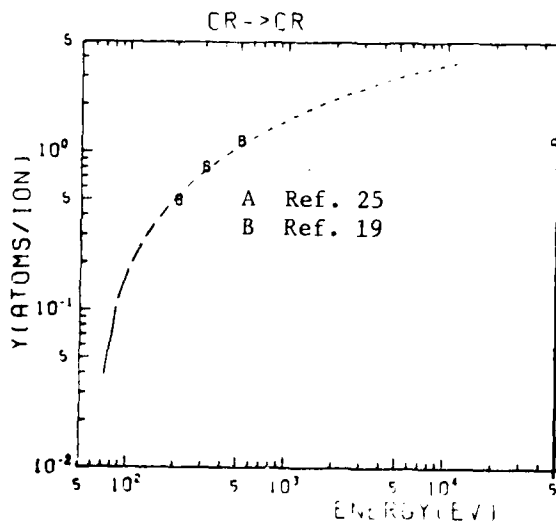
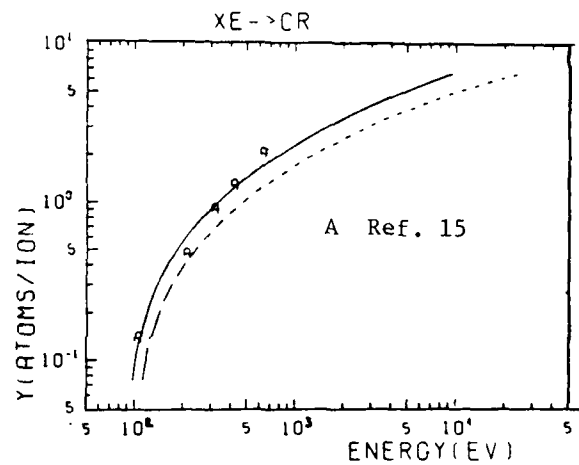
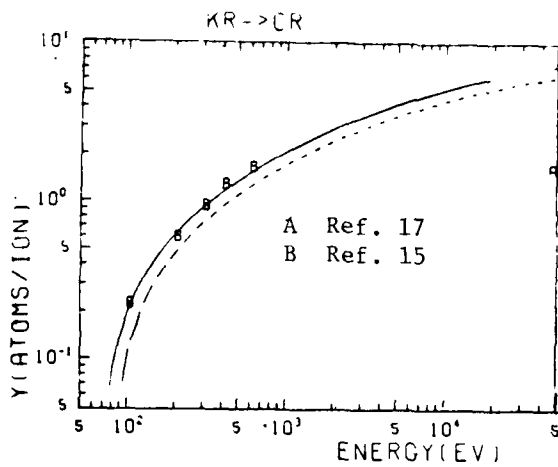
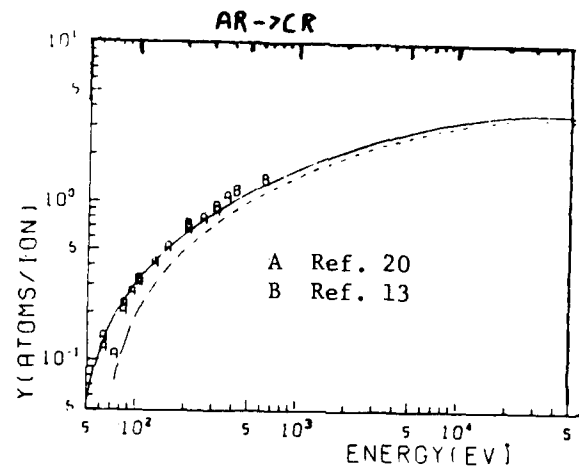
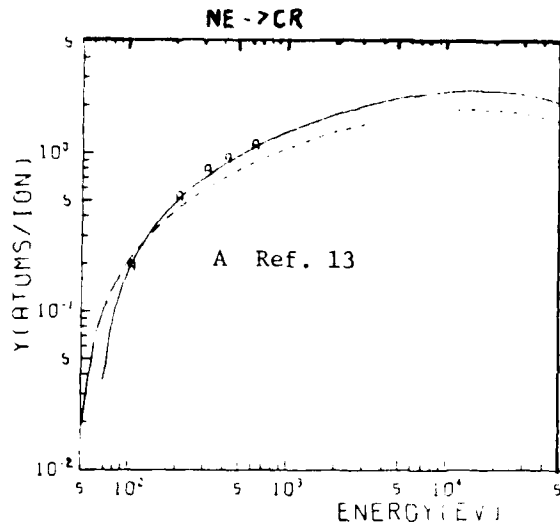
Graphical Data H-1.17

Energy dependence of the sputtering yield for Ar^+ , Kr^+ and Xe^+ impact on Ti. The solid line is a best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12.



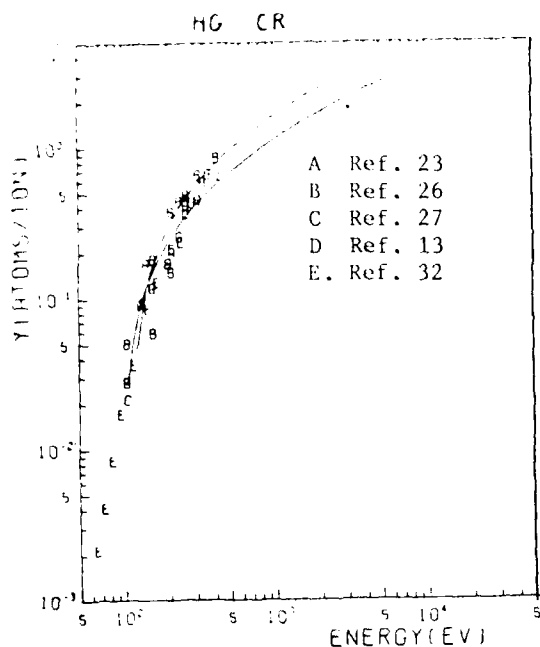
Graphical Data H-1.18

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ , Xe^+ and Hg^+ impact on V. The solid line is a best fit of Eq. 1 (section H-1.32) to the data. From Ref. 12.



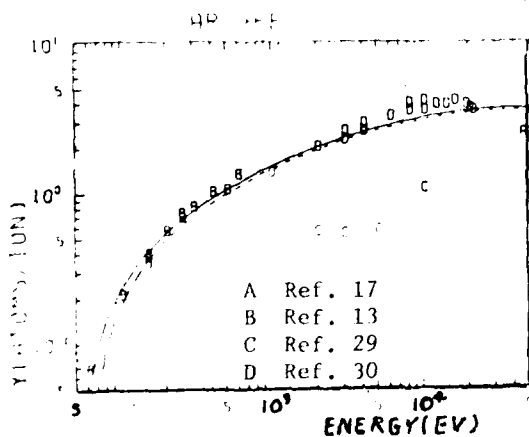
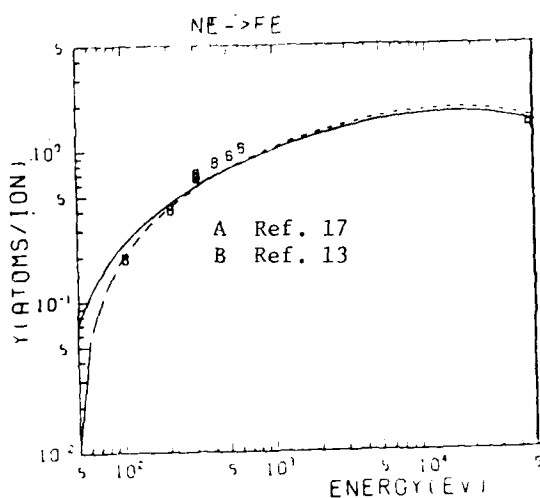
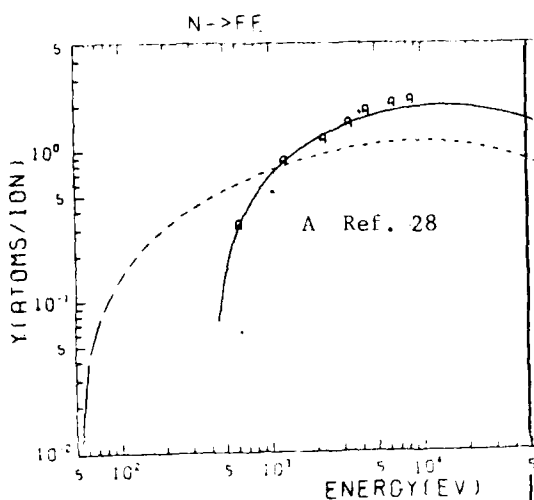
Graphical Date H-1.19.

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ and Xe^+ impact on Cr. The solid line is a best fit of Eq. 1 (section H-1.32) to the data. From Ref. 12



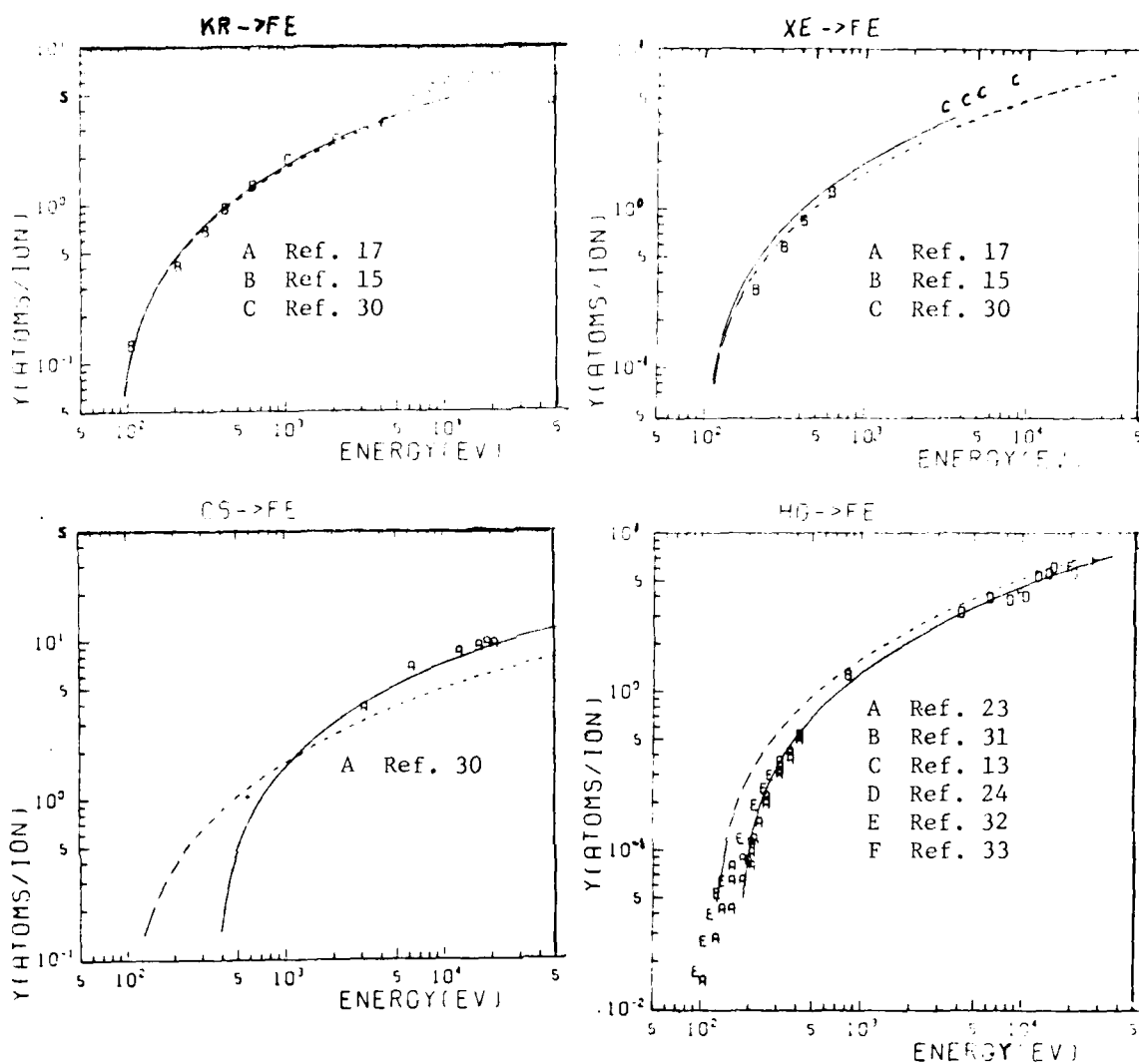
Graphical Data H-1.20

Energy dependence of the sputtering yield for Hg^+ impact on Cr. The solid line is a best fit of Eq. 1 (Section H-1. to the data. From Ref. 12.



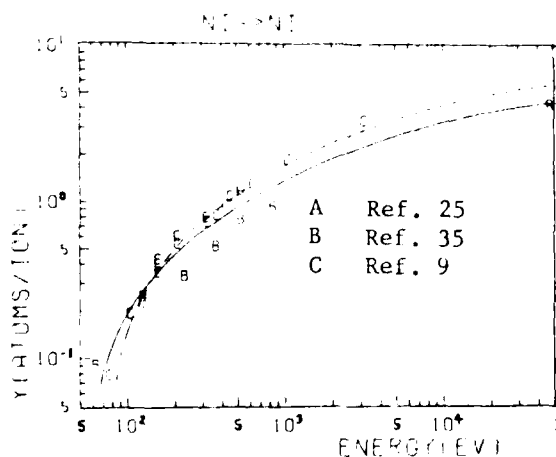
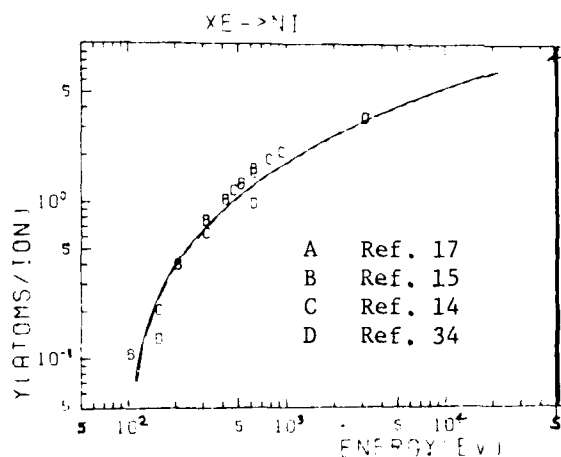
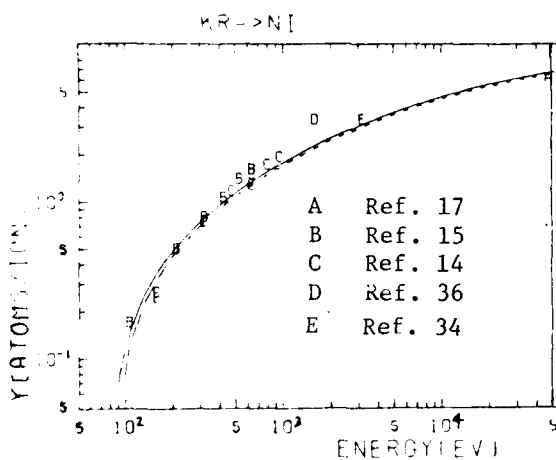
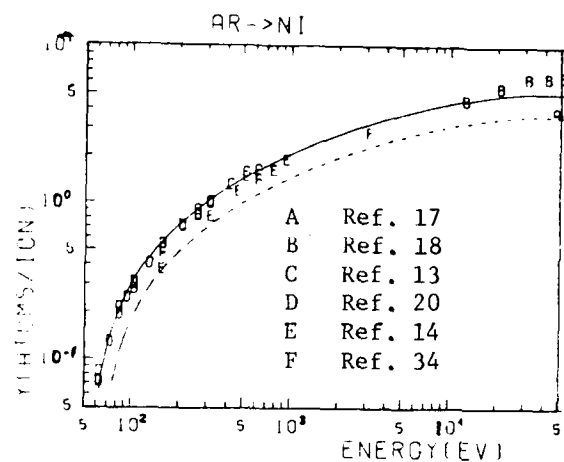
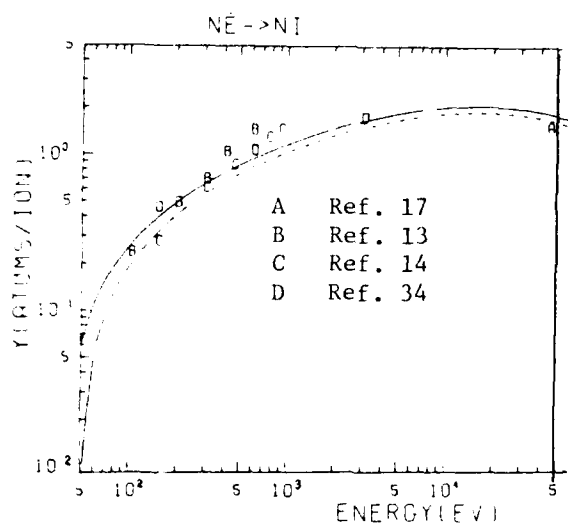
Graphical Data H-1.21

Energy dependence of the sputtering yield for N^+ , Ne^+ and Ar^+ impact on Fe. The solid line is a best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12



Graphical Data H-1.22

Energy dependence of the sputtering yield for Kr⁺, Xe⁺, Cs⁺ and Hg⁺ impact on Fe. The solid line is a best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12.



Graphical Data H-1.23

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ , Xe^+ and Ni^+ impact on Ni. The solid line in the best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12.

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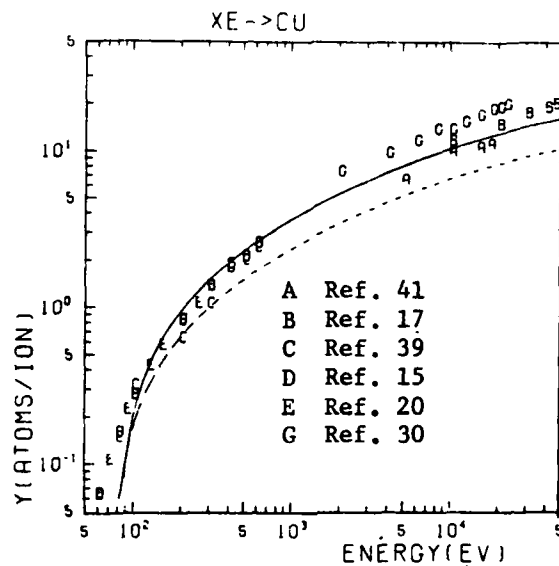
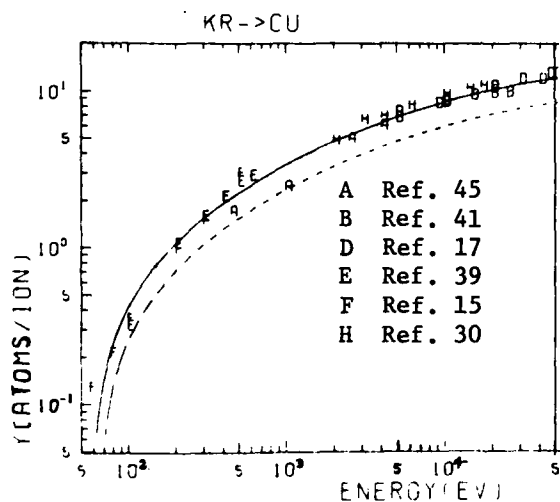
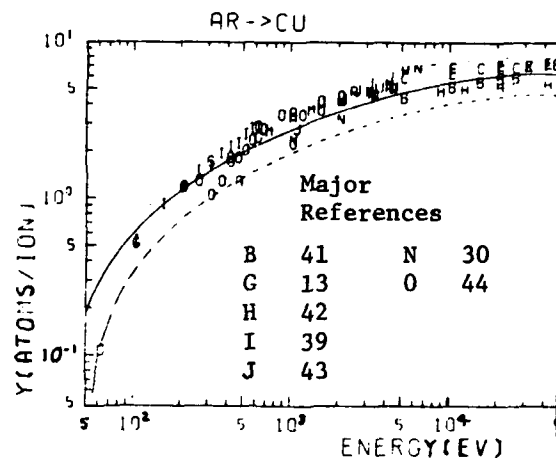
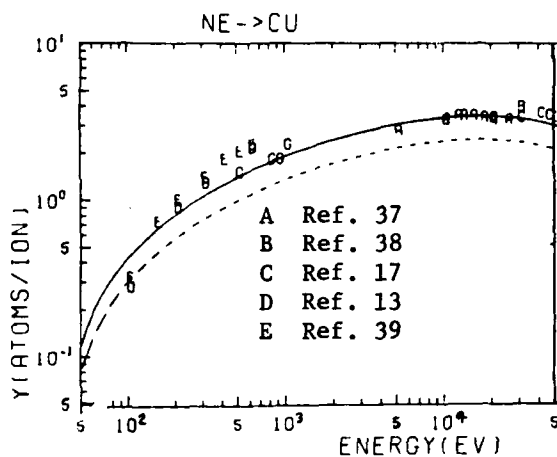
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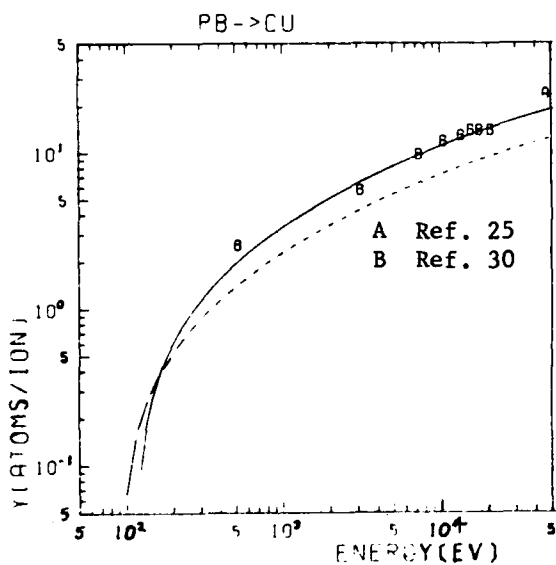
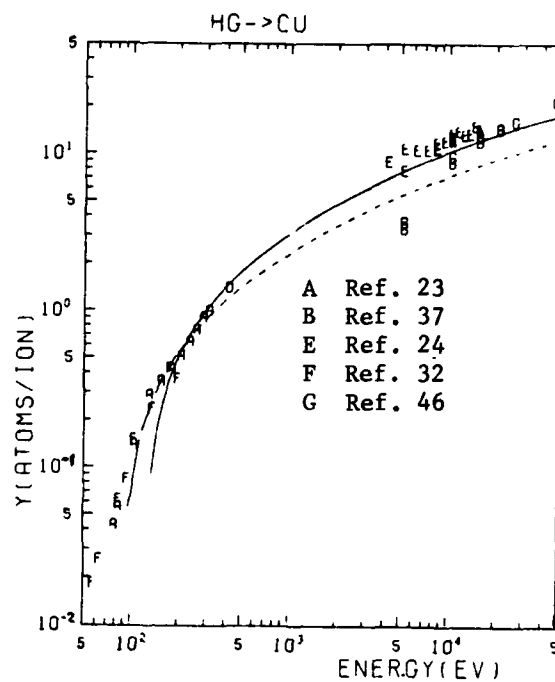
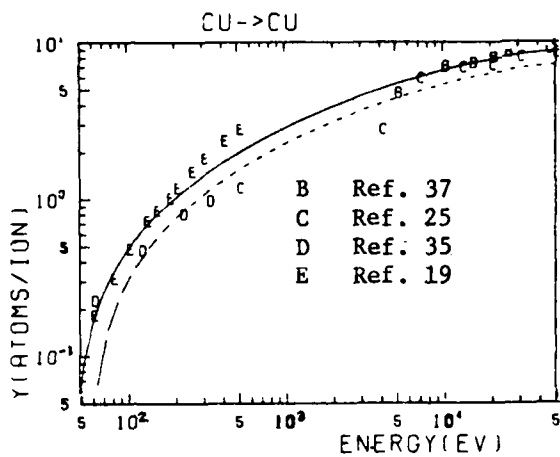
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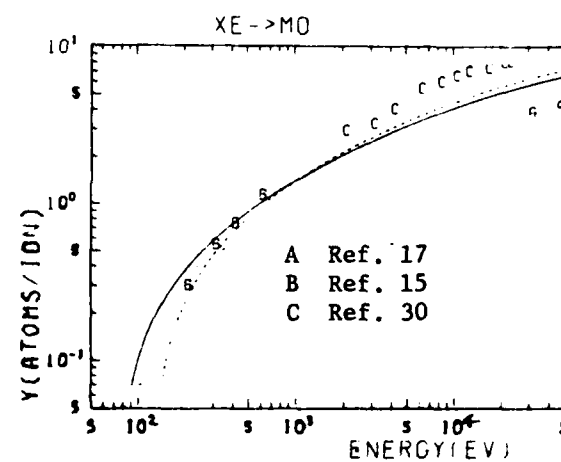
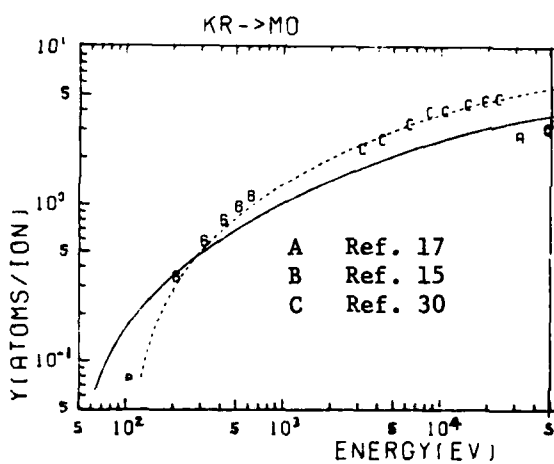
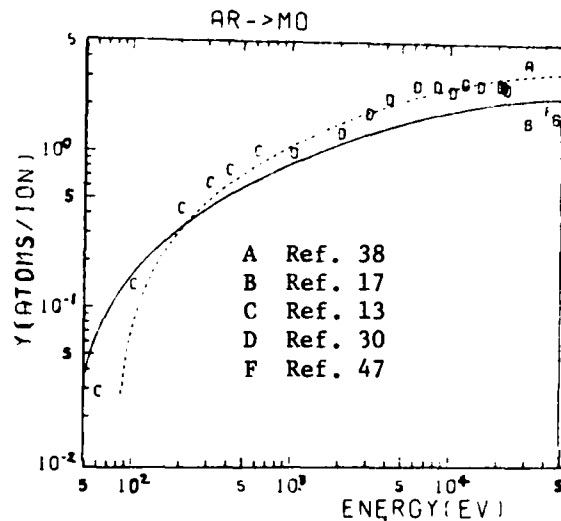
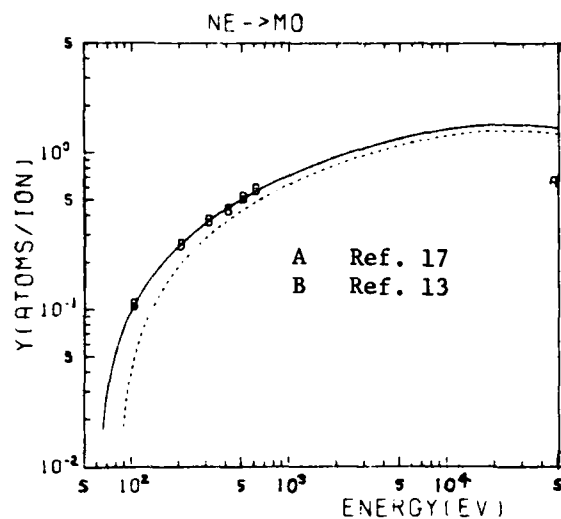
Graphical Data H-1.24

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ and Xe^+ impact on Cu. The solid line is the best fit of Eq. 1 (section H-1.32) to the data. From Ref. 12.



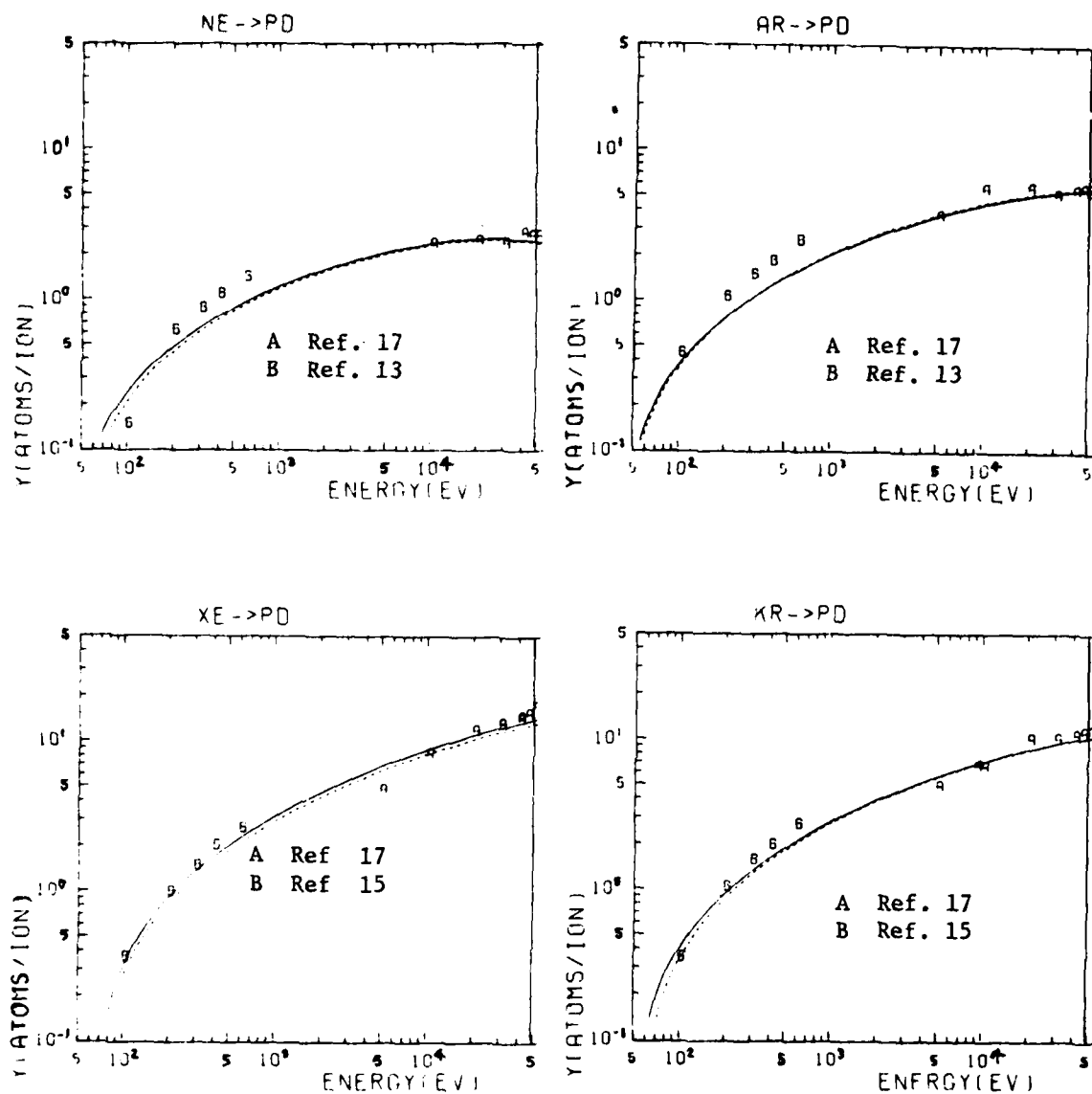
Graphical Data H-1.25

Energy dependence of the sputtering yield for Cu^+ , Hg^+ and Pb^+ impact on Cu. The solid line is the best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12.



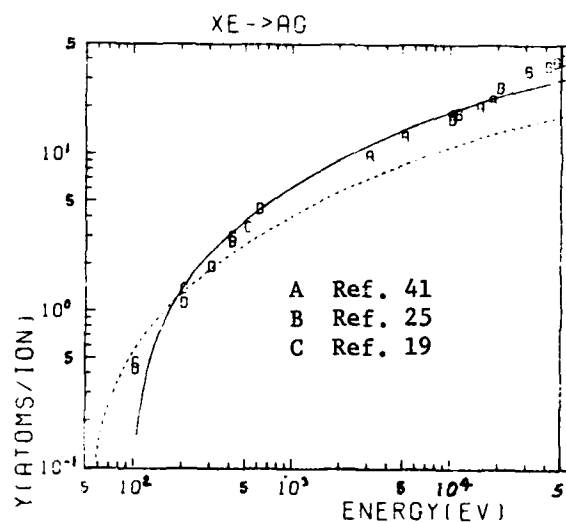
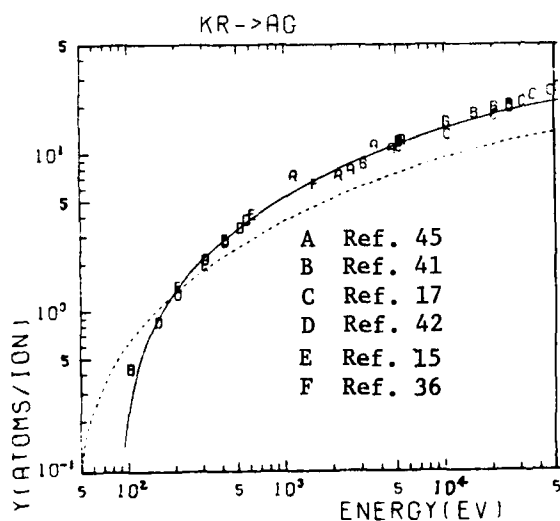
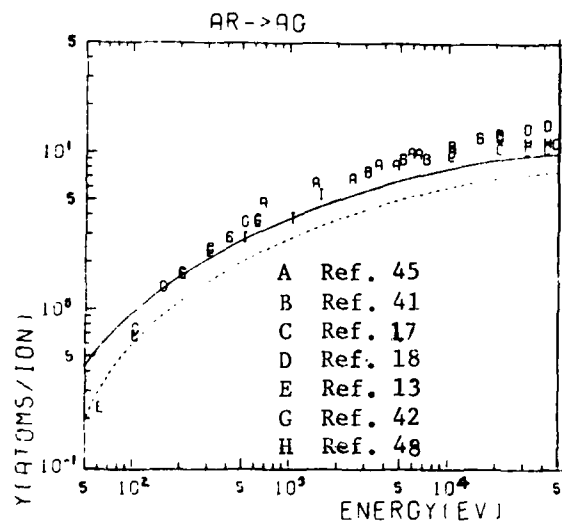
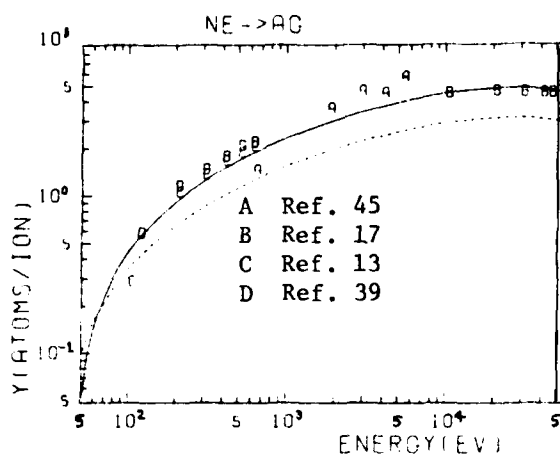
Graphical Data H-1.26

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ and Xe^+ impact on Mo. The solid line is the best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12.



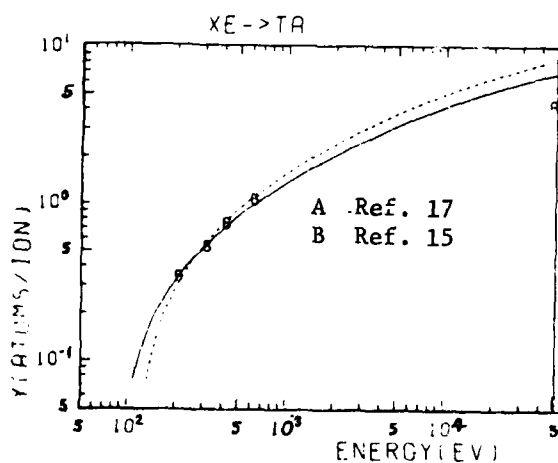
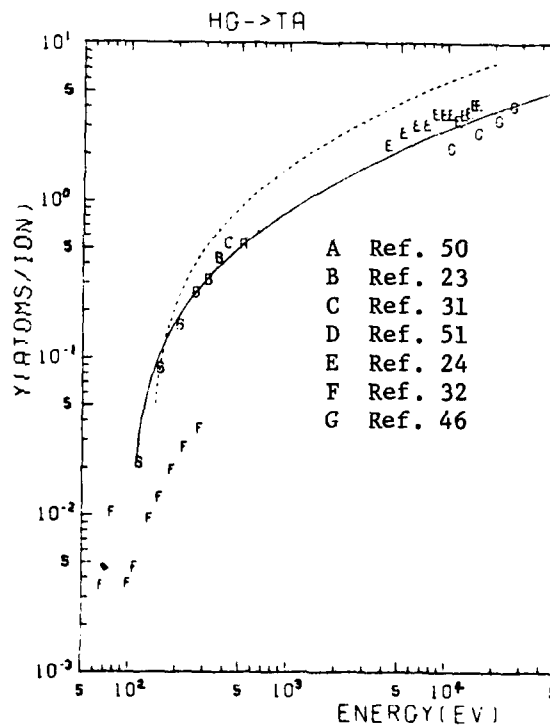
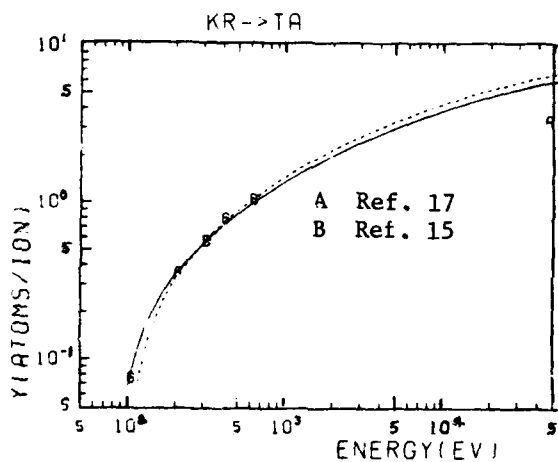
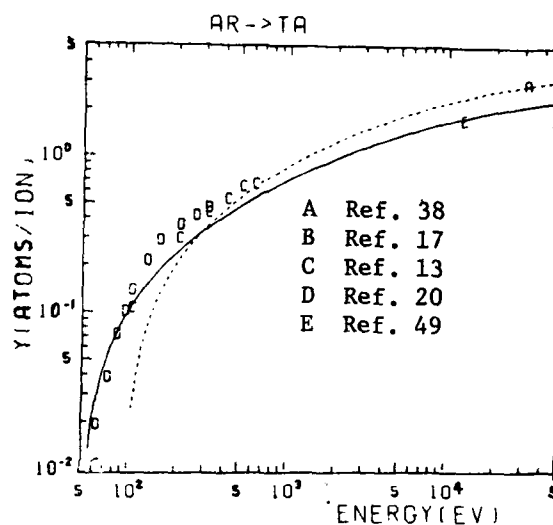
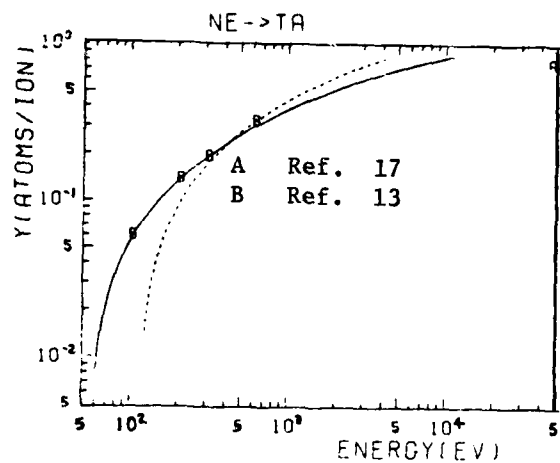
Graphical Data H-1.27

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ and Xe^+ impact on Pd. The solid line is the best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12.



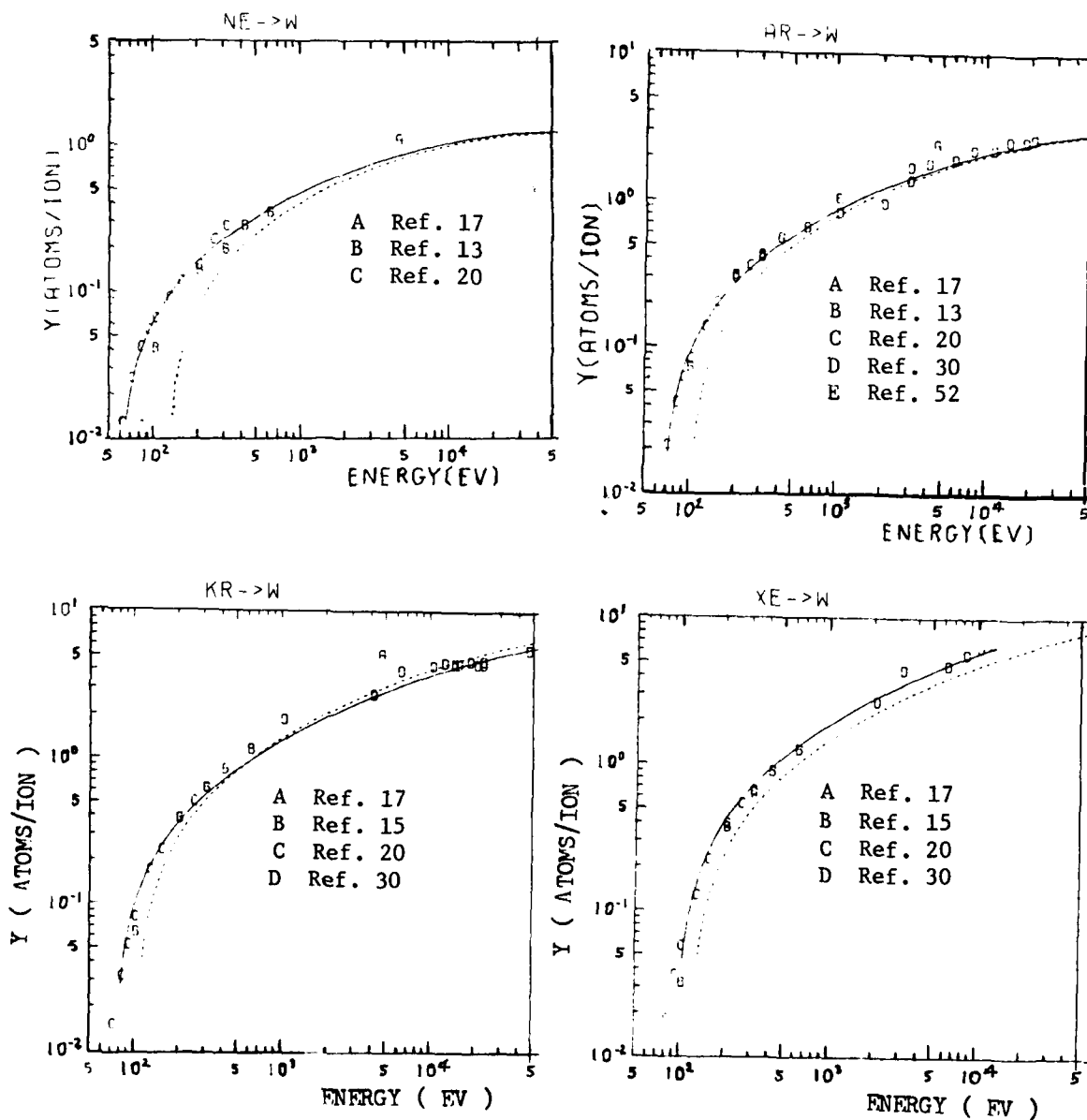
Graphical Data H-1.28

Energy dependance of the sputtering yield for Ne^+ , Ar^+ , Kr^+ and Xe^+ impact on Ag. The solid line is the best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12.



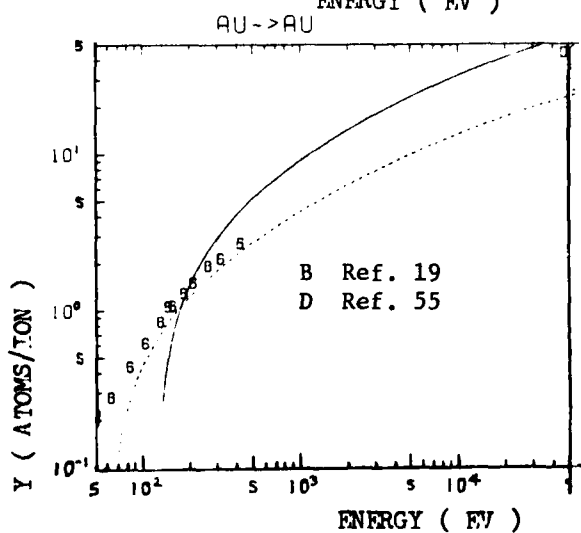
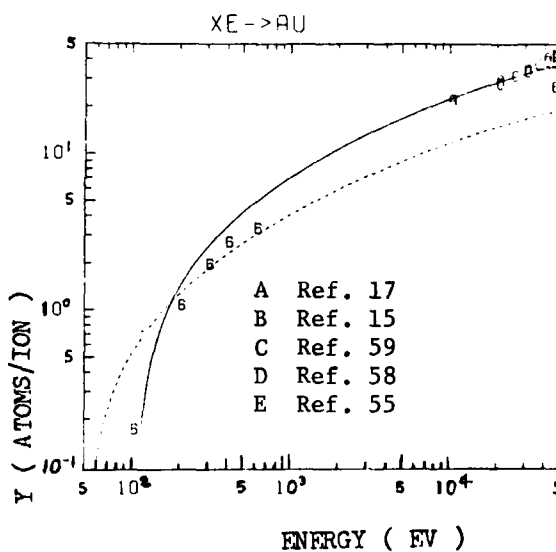
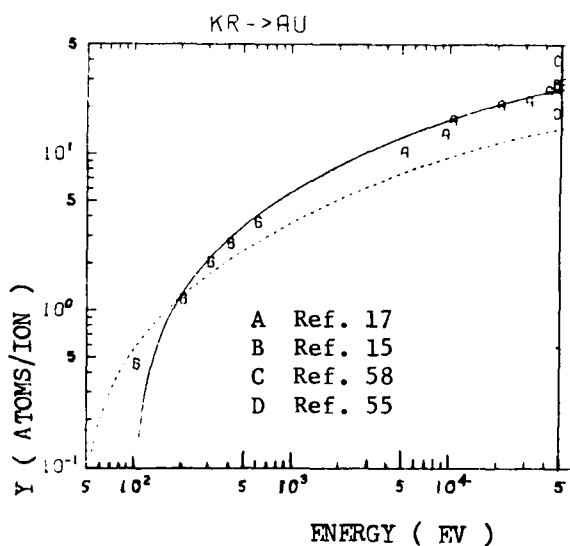
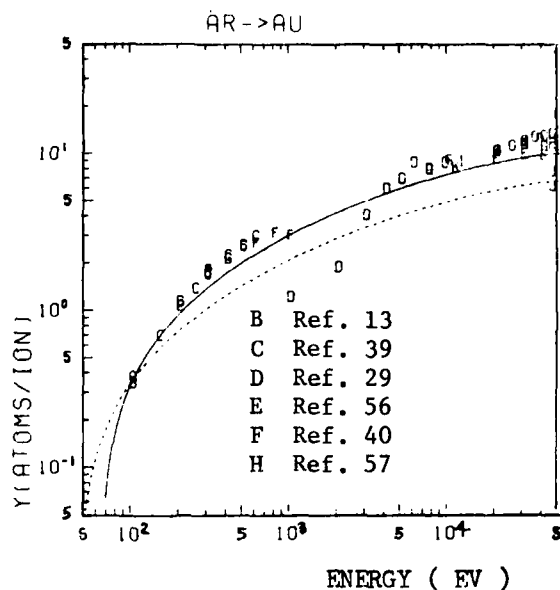
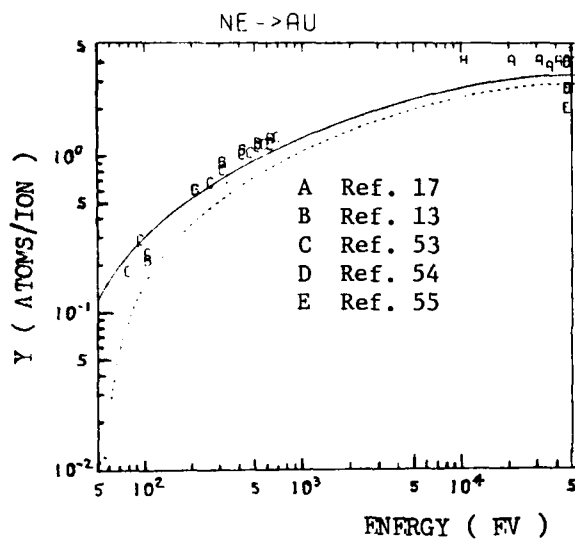
Graphical Data H-1.29

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ , Xe^+ and Hg^+ impact on Ta. The solid line is the best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12.



Graphical Data H-1.30

Energy dependence of the sputtering yield for Ne⁺, Ar⁺, Kr⁺ and Xe⁺ impact on W. The solid line is the best fit of Eq. 1. (Section H-1.32) to the data. From Ref. 12.



Graphical Data H-1.31

Energy dependence of the sputtering yield for Ne^+ , Ar^+ , Kr^+ , Xe^+ and Au^+ impact on Au. The solid line is the best fit of Eq. 1 (Section H-1.32) to the data. From Ref. 12

Tabular Data H-1.32

Semi-Empirical Formulation for Sputtering Yields

due to Heavy Particle Impact

Symbols

- E = Projectile energy (eV)
 E_{th} = Threshold energy for sputtering (found by fitting to the data) (eV).
 P = Dimensionless factor found by fitting to the data.
 S_n = Reduced elastic stopping cross section, dimensionless.
 (Expressed in analytical form below).
 ϵ = Reduced energy, dimensionless.
 C_e = A factor related to nuclear charge and mass of projectile and target;
 given in tabular form (eV^{-1})

Yield Expression.

$$Y = P S_n \left\{ 1 - \left(\frac{E_{th}}{E} \right)^{1/2} \right\} \quad \text{atoms/ion} \quad (1)$$

where

$$S_n = \frac{3.441 \sqrt{\epsilon} \log(\epsilon + 2.718)}{1 + 6.355 \sqrt{\epsilon} + \epsilon(-1.708 + 6.882 \sqrt{\epsilon})} \quad (2)$$

$$\epsilon = C_e E \quad (3)$$

The factors P and E_{th} are found by fitting the equations to the available data; the factor C_e is computed. All three factors are tabulated on a subsequent page.

Limitations.

Formulation has been tested for a variety of projectile target combinations and exhibits no obvious limitations. For light ion impact (H^+ , D^+ , He^+) at low energies (10 keV and below) the formulation of section H-1.13 fits the data better. The formulation appears to be inadequate close to threshold (see note 3 below).

Accuracy.

Represents empirical data to better than $\pm 25\%$.

Source.

N. Matsunami, Y. Yamanura, Y. Itikawa, N. Itoh, K. Kazumata, S. Miyagawa, K. Morita, and N. Shimizu. "Energy Dependence of Sputtering Yields of Monatomic Solids". Institute of Plasma Physics, Nagoya University, Nagoya, Japan. Report No. IPPJ-AM-14. To be published in Radiation Effects Letters.

Notes.

- (1) The fitted curve represented by Eq. 1 is shown in the Graphical Data H-1.15 through H-1.31 by the solid line.

- (2) The formulation for Y is similar to that by Sigmund [Phys. Rev. 184, 383 (1969) and 187, 768 (1969)] Sigmund gives analytical forms of the constant P but does not include the threshold energy factor (essentially Sigmund has $E_{th} = 0$).
- (3) The values for threshold energy obtained from the fit of Eq. 1 to the data differ from those given in H-1.13 for the cases where comparison is possible; they also disagree badly with theoretical values. The implication is that this equation is inadequate close to threshold.

Tabular Data H-1.32 (continued)

Empirical parameters for the sputtering equation of Matsunami et al.

Target	Projectile	C_e	P	E_{th}	Target	Projectile	C_e	P	E_{th}
Be	He	1.389E-3	0.9835	89.52	Cr	He	1.999E-4	0.7059	74.16
	Ne	9.382E-5	2.682	98.27		Ne	2.712E-5	6.471	62.70
	Ar	2.715E-5	3.620	73.10		Ar	1.093E-5	9.284	41.79
	Kr	5.988E-6	5.346	89.39		Kr	3.299E-6	19.10	66.76
	Xe	2.360E-6	7.536	138.8		Xe	1.493E-6	30.33	34.17
C	D	2.239E-3	0.1092	31.88	Hg	Hg	6.727E-7	21.28	96.39
	He	9.196E-4	0.4505	167.4	Fe	He	1.814E-4	0.5800	67.71
	Ar	2.184E-5	6.768	348.9		N	4.053E-5	5.994	407.7
	Kr	5.010E-6	1.936	93.73		Ne	2.509E-5	4.504	35.85
	Xe	2.007E-6	1.485	18.07		Ar	1.025E-5	9.133	52.56
Al	Hg	8.189E-7	12.44	286.5		Kr	3.134E-6	17.23	81.77
	He	4.085E-4	0.9009	151.4	Co	Xe	1.440E-6	27.00	96.42
	Ne	4.489E-5	4.565	93.17		Cs	1.396E-6	37.56	348.9
	Al	2.895E-5	4.748	39.44		Hg	6.515E-7	29.28	164.7
	Ar	1.592E-5	6.936	71.28		Pb	6.162E-7	38.45	25.37
Si	Kr	4.177E-6	11.85	81.07	Ni	He	1.734E-4	0.5597	89.25
	Xe	1.775E-6	15.34	86.28		Ne	2.430E-5	6.902	81.92
	Hg	7.556E-7	23.74	133.6		Ar	1.002E-5	11.64	65.54
	He	3.740E-4	0.4875	85.02		Co	5.260E-6	5.886	41.87
	Ne	4.183E-5	2.835	90.95		Kr	3.098E-6	15.47	86.61
Ar	Ar	1.497E-5	3.987	64.20	D	Xe	1.433E-6	22.70	82.71
	Kr	3.964E-6	6.912	92.71		Hg	6.515E-7	34.86	146.8
	Xe	1.692E-6	11.59	159.3	He	Ni	3.573E-4	0.05155	82.13
	Pb	6.810E-7	15.52	2.435		D	3.514E-4	0.1168	58.71
Ti	He	2.221E-4	0.2605	56.38		He	1.654E-4	0.5689	54.52
	Ne	2.944E-5	2.725	61.79		O	3.139E-5	1.670	82.86
	Ar	1.168E-5	3.933	57.15		Ne	2.322E-5	5.311	39.09
	Kr	3.450E-6	6.948	86.89		Ar	9.577E-6	12.69	53.94
	Xe	1.556E-6	10.37	118.6	Ni	Ni	4.832E-6	10.43	52.65
V	Hg	6.933E-7	13.84	106.2		Kr	2.964E-6	17.53	75.63
	He	2.108E-4	0.2536	90.93		Xe	1.371E-6	24.35	93.92
	Ne	2.840E-5	3.415	79.57		Hg	6.239E-7	36.43	115.4
	Ar	1.139E-5	3.899	44.73					
	Kr	3.409E-6	8.466	79.38					
	Xe	1.548E-6	12.38	83.06					
	Hg	6.936E-7	19.18	160.2					

Tabular Data H-1.32 (continued)

Empirical parameters for the sputtering equation of Matsunami et al.

Target	Projectile	C_e	P	E_{th}	
Cu	D	3.366E-4	0.2694	405.3	
	He	1.589E-4	0.3420	49.10	
	N	3.629E-5	4.756	22.06	
	Ne	2.269E-5	9.678	37.07	
	Ar	9.477E-6	16.01	33.93	
	Cu	4.452E-6	22.27	44.46	
	Kr	2.980E-6	31.52	56.50	
	Xe	1.391E-6	49.07	79.87	
	Cs	1.350E-6	68.60	186.1	
	Hg	6.375E-7	63.21	122.1	
	Pb	6.035E-7	68.25	111.3	
	Ge	He	1.411E-4	0.3502	79.96
		Ne	2.073E-5	5.178	67.53
Ar		8.850E-6	8.296	29.12	
Kr		2.862E-6	14.82	72.07	
Xe		1.358E-6	20.54	93.73	
Hg		6.310E-7	24.89	110.8	
Zr	He	1.069E-4	0.1647	163.3	
	Li	6.788E-5	0.6451	101.8	
	Ne	1.648E-5	3.008	70.80	
	Ar	7.294E-6	6.061	53.12	
	Kr	2.477E-6	6.196	60.16	
	Xe	1.211E-6	12.75	99.15	
	Hg	5.777E-7	12.60	91.80	
Nb	D	2.163E-4	0.02584	52.37	
	He	1.036E-4	0.2386	191.0	
	Li	6.585E-5	0.8341	124.0	
	Ne	1.603E-5	2.896	63.97	
	Ar	7.118E-6	7.456	75.26	
	Kr	2.427E-6	10.75	90.38	
	Nb	1.984E-6	10.39	367.5	
	Xe	1.190E-6	25.50	207.5	
	Cs	1.156E-6	22.39	4.010	
	Hg	5.690E-7	12.41	105.9	

Target	Projectile	C_e	P	E_{th}
Ru	H	2.119E-4	8.914E-3	327
	D	2.098E-4	0.01602	131.5
	He	1.006E-4	0.1457	99.28
	Ne	1.565E-5	3.848	59.02
	Ar	6.980E-6	5.405	39.32
	Kr	2.396E-6	9.982	44.13
	Xe	1.179E-6	19.59	73.07
	Cs	1.146E-6	31.21	253.7
	Hg	5.659E-7	19.81	135.0
	Pb	5.374E-7	31.26	135.1
	Ne	1.490E-5	4.774	72.74
	Ar	6.696E-6	10.55	58.70
	Kr	2.323E-6	21.83	86.21
	Xe	1.152E-6	28.50	91.43
Rh	He	9.221E-5	0.2683	90.57
	Ne	1.454E-5	6.251	75.52
	Ar	6.549E-6	13.88	68.49
	Kr	2.281E-6	24.30	76.39
	Xe	1.133E-6	31.30	80.13
	Hg	5.485E-7	41.58	96.42
	He	8.973E-5	0.6540	75.61
Pd	Ne	1.422E-5	6.440	41.67
	Ar	6.435E-6	13.96	41.77
	Kr	2.256E-6	28.52	49.03
	Xe	1.126E-6	43.72	63.04
	Hg	5.468E-7	52.60	78.25
	He	1.831E-4	0.1271	268.3
	D	1.815E-4	0.2666	44.60
Ag	Ne	1.387E-5	13.09	45.34
	Ar	6.292E-6	24.67	23.39
	Kr	2.212E-6	60.33	84.88
	Ag	1.443E-6	89.25	89.07
	Xe	1.106E-6	94.51	95.93
	Hg	5.384E-7	140.	123.6
	He	1.831E-4	0.1271	268.3

Tabular Data H-1.32 (continued)

Empirical parameters for the sputtering equation of Matsunami et al.

Target Projectile	C_0	P	E_{th}	Target Projectile	C_0	P	E_{th}
Sn Ar	5.982E-6	10.44	51.82	Ir He	4.664E-5	0.07918	18.1
Sn	1.249E-6	26.60	56.43	Ne	8.016E-6	4.323	64.22
Hf Ne	8.665E-6	3.125	63.57	Ar	3.889E-6	12.39	57.87
Ar	4.172E-6	7.804	39.20	Kr	1.518E-6	22.01	74.44
Kr	1.608E-6	16.92	68.16	Xe	8.169E-7	42.31	92.65
Xe	8.578E-7	30.57	118.3	Hg	4.268E-7	46.20	93.30
Hg	4.440E-7	22.56	109.0	Pt He	4.588E-5	0.2193	179.3
Ta Ne	8.527E-6	2.457	54.64	Ne	7.900E-6	5.177	56.34
Ar	4.112E-6	5.801	50.15	Ar	3.837E-6	14.84	54.99
Kr	1.589E-6	17.35	81.60	Kr	1.501E-6	32.93	83.10
Xe	8.488E-7	22.93	86.49	Xe	8.093E-7	50.28	92.13
Hg	4.402E-7	19.56	99.34	Hg	4.236E-7	67.79	112.3
W He	4.908E-5	0.05713	165.3	Au H	9.300E-5	0.05583	372.8
Ne	8.395E-6	3.031	57.92	D	9.253E-5	0.1261	152.0
Ar	4.054E-6	7.493	65.92	He	4.513E-5	0.4742	68.88
Kr	1.571E-6	16.32	73.80	N	1.169E-5	4.327	30.89
Xe	8.408E-7	32.15	95.61	Ne	7.780E-6	7.831	28.05
Cs	8.191E-7	30.21	51.47	Ar	3.783E-6	26.05	62.12
Hg	4.369E-7	16.31	92.18	Kr	1.482E-6	76.97	97.73
Re Ne	8.264E-6	3.864	77.98	Xe	8.003E-7	126.0	105.2
Ar	3.996E-6	11.18	71.67	Au	4.259E-7	221.0	120.7
Kr	1.552E-6	24.96	85.87	Hg	4.194E-7	134.6	36.26
Xe	8.319E-7	33.11	96.64	Th Ne	6.688E-6	3.436	82.55
Hg	4.330E-7	32.46	105.6	Ar	3.300E-6	7.758	53.79
Os He	4.744E-5	0.1555	187.0	Kr	1.325E-6	20.21	78.22
Ne	8.144E-6	3.839	60.74	Xe	7.292E-7	24.34	83.14
Ar	3.946E-6	12.11	84.66	U Ne	6.517E-6	5.003	70.40
Kr	1.537E-6	25.67	91.50	Ar	3.223E-6	11.79	63.84
Xe	8.261E-7	35.74	95.12	Kr	1.299E-6	27.14	75.83
				Xe	7.169E-7	27.44	122.3
				Hg	3.844E-7	35.55	60.23

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H-3. SECONDARY ELECTRON EMISSION BY ION IMPACT

CONTENTS

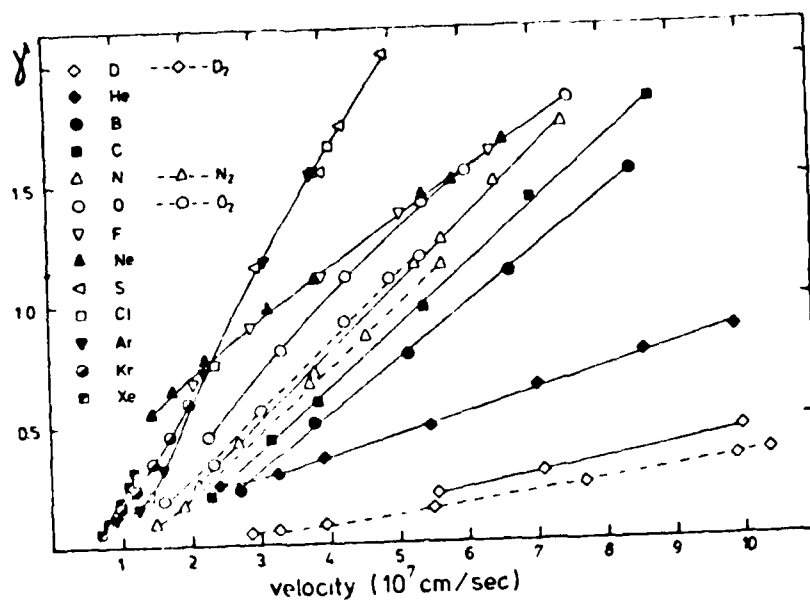
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H-3.1 Secondary Electron Emission Coefficient for Various Ions on Al.....	
H-3.2 Secondary Electron Emission Coefficients for Ne^+ , Ar^+ , Kr^+ and Xe^+ impact on Mo and Cu.....	
H-3.3 Secondary Electron Emission Coefficients for H^+ , H_2^+ , H_3^+ , He^+ , Ne^+ and Ar^+ on Au and W.....	
H-3.4 Secondary Electron Emission Coefficients for Ar^+ Impact on Mo, Ta, Zr, Al, Cu and Ni.....	

INTRODUCTION

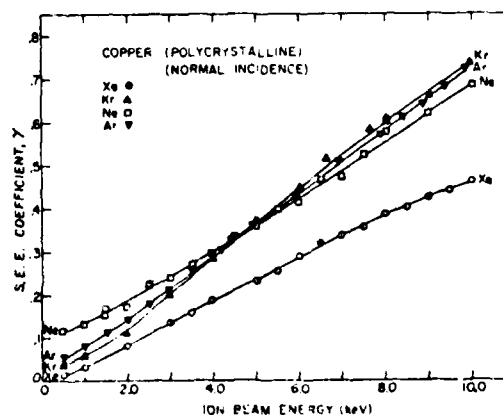
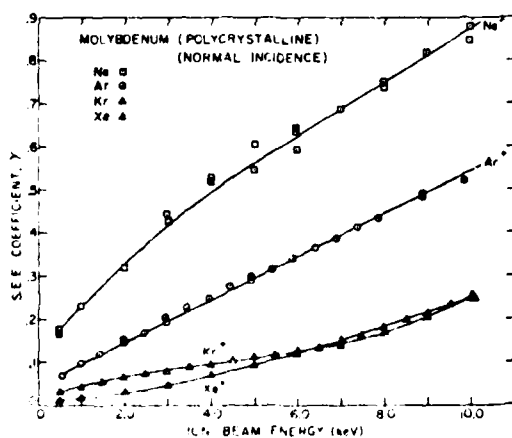
This section updates Section H-3 of Technical Report H-78-1 "Compilation of Data Relevant to Rare Gas and Rare Gas - Monohalide Excimer Lasers. Volume II" (U.S. Army Missile Research and Development Command, Redstone Arsenal, Alabama, December 1977). For low energy impact of light projectiles the earlier compilation cited above remains quite adequate as there has been no significant additional data. What we show here is some additional data for projectile energies of 500eV and greater. Data given are for normal incidence on polycrystalline targets. Some excellent data on single crystal targets is given in reference 4 but not reproduced here.

References.

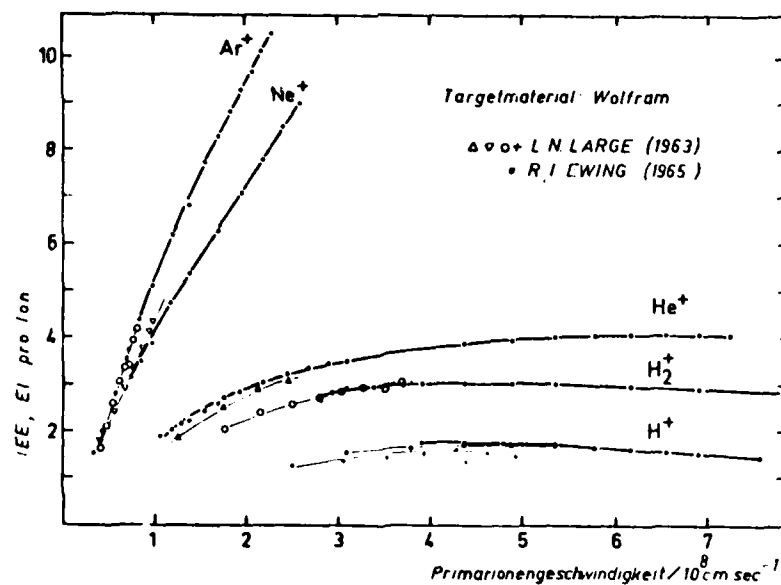
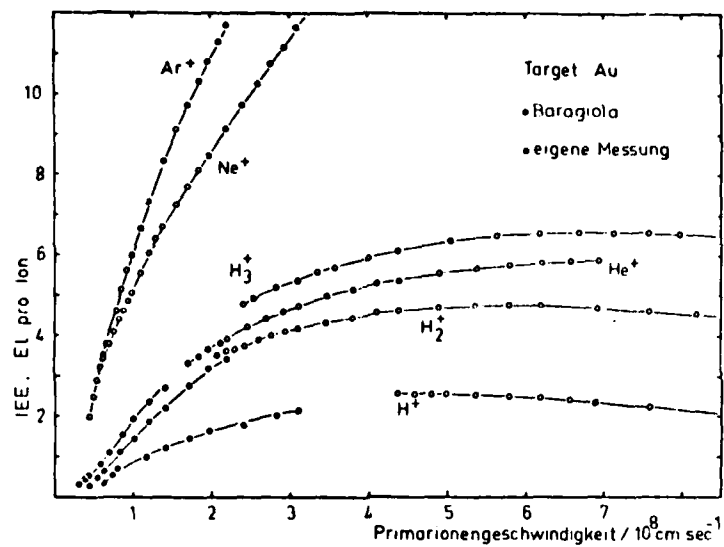
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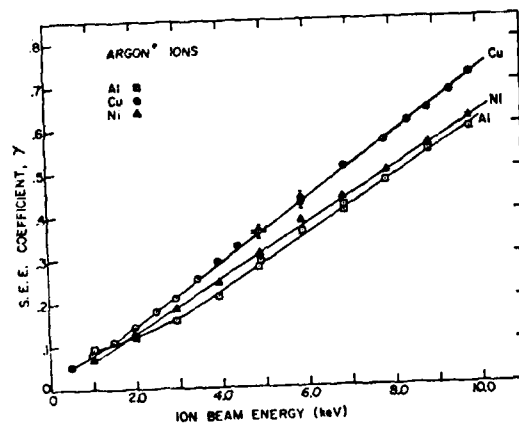
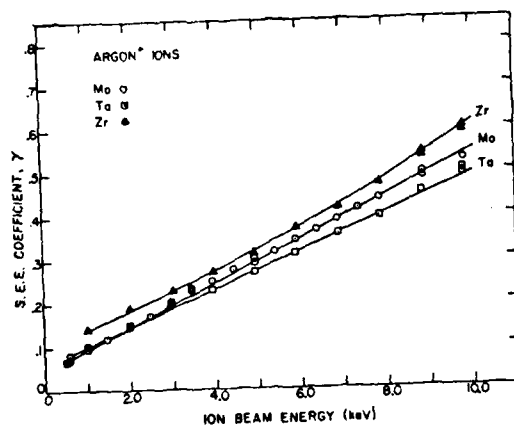
Graphical Data H-3.1. Secondary electron emission coefficient for various ions on Al as a function of projectile velocity. From Ref. 1.



Graphical Data H-3.2. Secondary electron emission coefficients for Ne^+ , Ar^+ , Kr^+ and Xe^+ impact on Mo and Cu. From Ref. 2



Graphical Data H-3.3 Secondary electron emission coefficient for H^+ , H_2^+ , H_3^+ , He^+ , Ne^+ and Ar^+ on Au and W as a function of impact velocity. From Ref. 3 Includes also data from Ref. 5



Graphical Data H-3.4. Secondary electron emission coefficients for Ar⁺ impact on Mo, Ta, Zr, Al, Cu and Ni. From Ref. 2.

H-5. ION REFLECTION FROM SURFACES

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Graphical Data H-5.2. Reflection Coefficients of heavy ions	3203

INTRODUCTION

This section presents data on coefficients for reflection of ions from surfaces as the ratio of particles reflected (ions plus atoms, integrated over all exit angles and energies) to particles incident. This supplements coverage of such processes given previously as section H-5 of Technical Report H-78-1, "Compilation of Data Relevant to Rare Gas-Rare Gas and Rare Gas - Monohalide Excimer Lasers Volume II" (U.S. Army, Missile Research and Development Command, Redstone Arsenal, Alabama, December 1977). For light ion impact (H,D,He) there is excellent coverage due in large measure to the need for such data in the Fusion Energy Program. We provide an extensive presentation of data and indicate a reliable method for interpolating values for cases that have not yet been considered (H-5.1). For heavier species we provide what little data are in the literature (H-5.2). For heavy projectiles at low energies one might make order to magnitude estimates by using some average of the data given by the scaling procedure of H-5.1.

Tabular Data H-5.1. Reflection of Light Ions (H,D,He).

A substantial quantity of experimental data is available which agrees well with theoretical predictions. There is no evidence that the reflection coefficient for an incoming ion (e.g. H^+) is any different from that for the corresponding atom (e.g. H).

It is most convenient to plot the reflection coefficient R_N (particles reflected, integrated over all exit angles and energies : by particles incident) as a function of the reduced energy ϵ defined as follows:

$$\epsilon = \frac{M_2}{M_1 + M_2} \frac{a}{Z_1 Z_2 e^2} E$$

Lindhard has chosen for 'a' the Thomas-Fermi screening length a_{TF}

$$a_{TF} = 0.4685 (Z_1^{2/3} + Z_2^{2/3})^{-1/2}$$

Z_1 and Z_2 are the charges and M_1 and M_2 are the masses of the ion and the target atom. e is the electron charge.

With $e^2 = 14.39 \text{ eV } \text{\AA}$ one gets

$$\epsilon = 32.55 \frac{M_2}{M_1 + M_2} \frac{1}{Z_1 Z_2 (Z_1^{2/3} + Z_2^{2/3})^{1/2}} E(\text{keV}) = \epsilon_L E(\text{keV}) \quad (1)$$

In the following figure we show theoretical model calculations of R_N plotted as a function of reduced energy ϵ , for a variety of projectile-target combinations. We choose to present theory rather than experiment because it covers a greater energy range particularly down to low energies. Experiment is in good agreement with theory as is shown in the primary references given below.

To obtain R_N for a projectile-target combination given on the graph the factor ϵ_L should be identified from the table, the reduced energy of interest should be computed and the relevant R_N read from the graph. For a projectile-target combination not on the graph the factor ϵ_L should be calculated from the formulae given above, the reduced energy of interest ϵ evaluated and data points read from the graph for a target of nuclear charge closest to that of interest. For compounds one may estimate R_N by evaluating the reflection coefficients for each constituent atomic species and take a sum weighted according to the atomic composition of the target compound or alloy.

This scaling procedure should provide adequate estimates of R_N to within a factor of two, except for low Z targets (e.g. C) where the error will be larger. The data obtained by this procedure will be appropriate to normal incidence on a polycrystalline target.

- Reference. (i) The graph is taken from: W. Eckstein and H. Verbeek "Data on Light Ion Reflection", Report IPP 9/32 (Max Planck Institut für Plasma Physik, Garching, West Germany, August 1979. This report contains extensive theoretical and experimental data gathered from various published and unpublished sources.
- (ii) The theoretical simulation is by the MARLOWE code described most fully in : M. T. Robinson and I. M. Torrens, Phys. Rev. B9, 5008 (1974). See also S. Oen, and M. T. Robinson, Nucl. Instr. and Methods 132, 647 (1976).

Target atom			ϵ_L				
Element	Z_2	M_2	H	D	T	^3He	^4He
C	6	12.0	2.414	2.242	2.092	0.9814	0.9200
Al	13	26.98	0.9449	0.9123	0.8819	0.4223	0.4087
Si	14	28.09	0.8604	0.8318	0.8050	0.3862	0.3742
Ti	22	47.90	0.4871	0.4774	0.4680	0.2266	0.2222
Fe	26	55.85	0.3934	0.3866	0.3800	0.1845	0.1814
Ni	28	58.69	0.3575	0.3516	0.3459	0.1682	0.1655
Cu	29	63.54	0.3420	0.3368	0.3317	0.1614	0.1590
Nb	41	92.91	0.2188	0.2165	0.2142	0.1047	0.1037
Mo	42	95.95	0.2121	0.2099	0.2078	0.1016	0.1006
Ag	47	107.87	0.1832	0.1816	0.1799	0.08814	0.08735
Ta	73	180.95	0.1032	0.1026	0.1021	0.05024	0.04997
W	74	183.92	0.1014	0.1008	0.1003	0.04937	0.04911
Au	79	197.0	0.09305	0.09258	0.09212	0.04538	0.04515

Table showing computed ϵ_L factors for H,D, ^3He and ^4He impact on a variety of targets.

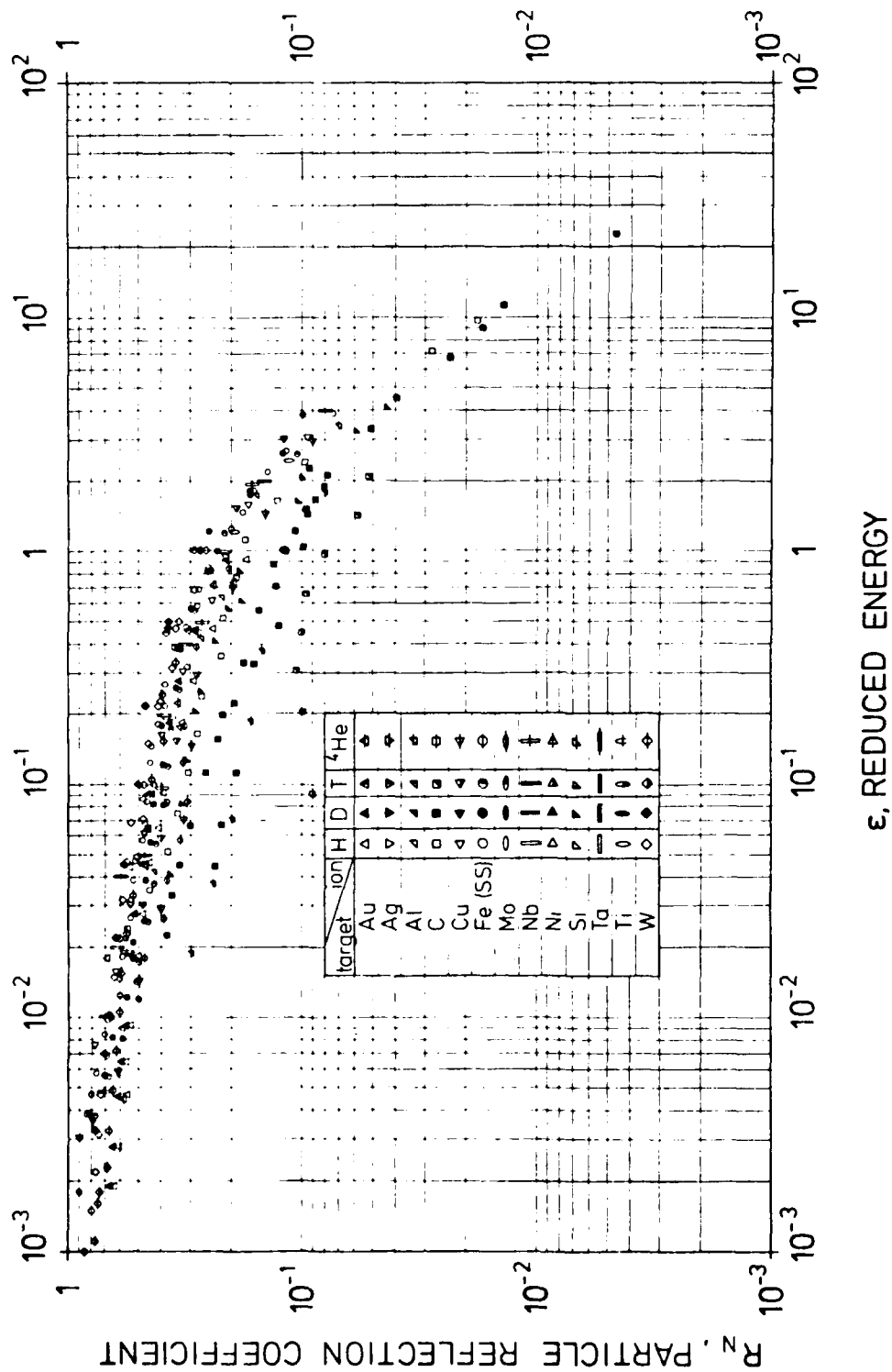
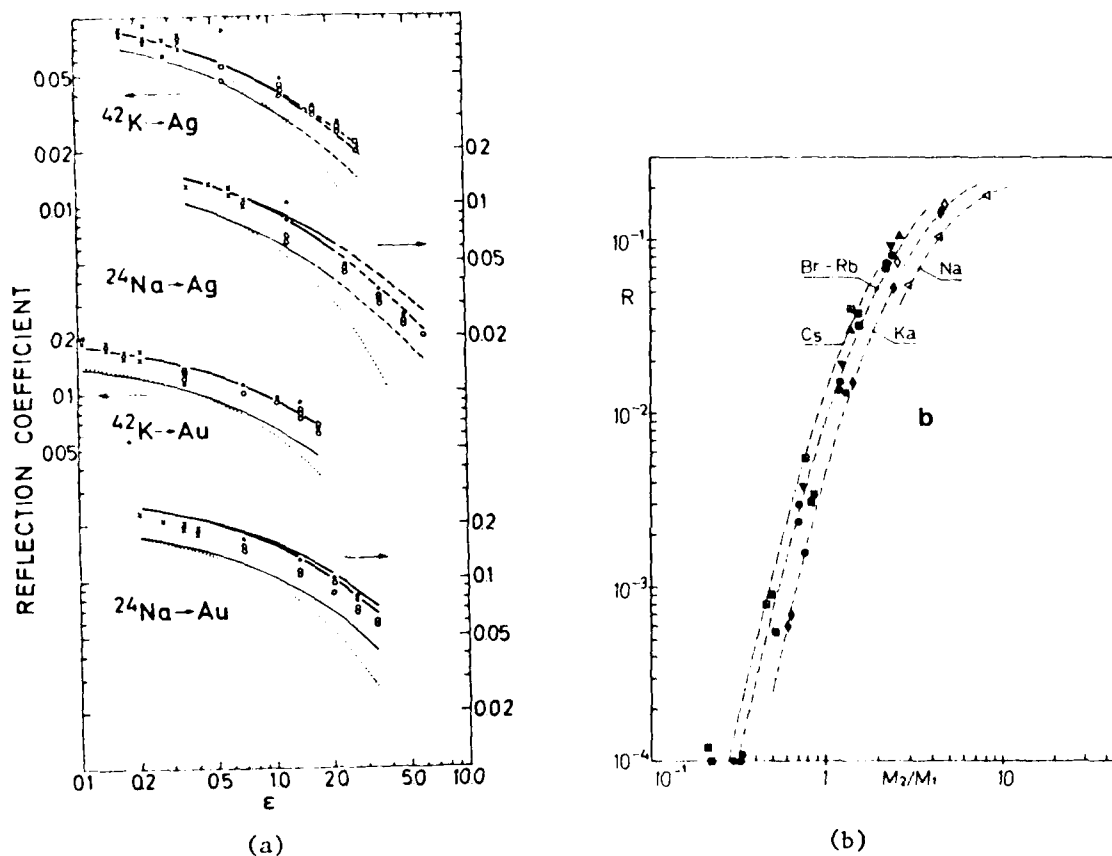


Fig. H-5.2: Calculated particle reflection coefficients for several ion-target combinations versus the reduced energy ϵ . Reproduced from reference (i).



Graphical Data H-5.2. Reflection coefficients for heavy particle scattering from surfaces.

- K and Na impact on Ag and Au shown as a function of reduced energy (defined by Eq. 1 of preamble to Graphical Data H-5.1). J. Bottiger et. al., Radiation Effects, 11, 133 (1971).
- Impact of 60 keV Na, K (written Ka), Br, Rb and Cs on Al, Cu, Ag and Au shown for each projectile as a function of target to projectile mass ratio (M_2/M_1). J. Bottinger et. al., Nucl. Instr. Meth. 170, 499 (1980).

I. SECONDARY ELECTRON SPECTRA
CONTENTS

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I - 1. Energy Spectra of Secondary Electrons from Electron Impact Ionization	3205
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(data presented in this chapter either extend or supersede the
data given previously in Chapter I of Volume V, pages 2241-2374.)

I - 1. ENERGY SPECTRA OF SECONDARY ELECTRONS
FROM ELECTRON IMPACT IONIZATION

Page

I-1.1

Secondary Electron Energy Spectra for
Electron Impact Ionization of CO₂ 3206

Tabular Data I-2.1
Single Differential Cross Section (Secondary
Electron Spectra) for $e^- + CO_2$ Collisions
(Units of $10^{-18} \text{ cm}^2/\text{eV}$)

Secondary Electron Energy (eV)	Primary Electron Energy (eV)			
	50	100	200	400
1	16.08	22.38	15.66	12.66
2	25.47	27.62	19.29	12.50
3	23.37	26.62	20.12	14.44
4	16.57	19.94	17.52	11.43
5	14.60	17.53	14.44	10.51
6	13.51	14.90	12.90	9.32
8	12.79	12.85	10.64	7.70
10	12.50	10.12	8.76	6.30
12	11.26	8.55	7.39	5.47
15	9.61	8.18	5.77	4.52
20		6.18	4.46	3.26
25		4.18	3.22	2.38
30		3.14	2.57	1.76
35		2.87	1.77	1.32
40		2.65	1.32	1.04
50			0.81	0.687
65			0.54	0.401
80			0.49	0.277
100				0.188
120				0.137
140				0.131
160				0.129
180				0.123

Accuracy: The quoted uncertainty in these data is $\pm 17\%$

Reference: These data were taken from T.W. Shyn and W.E. Sharn, Phys. Rev. A 20, 2332 (1979).

I - 2. ENERGY SPECTRA OF SECONDARY ELECTRONS
FROM PROTON IMPACT IONIZATION

Page

I-2.1	
Secondary Electron Energy Spectra for	
Proton Impact Ionization of Ar	

Tabular Data I-2.1

Single Differential Cross Section (Secondary
Electron Spectra) for H^+ + Ar Collisions

(Units of cm^2/eV)

ENERGY (eV)	5 keV	ENERGY (eV)	10 keV	20 keV	ENERGY (eV)	50 keV
1.	6.44-1A	1.	7.14-1A	2.44-17	1.	3.24-17
2.	8.72-1A	2.	1.10-17	2.30-17	2.	2.91-17
3.	8.44-1A	3.	1.39-17	2.25-17	3.	2.55-17
4.	7.27-1A	4.	1.26-17	2.08-17	4.	2.22-17
5.	6.25-1A	5.	1.16-17	1.89-17	5.	2.09-17
6.	5.57-1A	6.	1.04-17	1.70-17	6.	1.80-17
7.	5.82-1A	7.	9.92-1A	1.60-17	7.	1.68-17
8.	4.43-1A	8.	8.40-1A	1.40-17	8.	1.53-17
9.	3.85-1A	9.	7.68-1A	1.38-17	9.	1.44-17
10.	4.46-1A	10.	7.36-1A	1.27-17	10.	1.32-17
11.	4.25-1A	12.	6.13-1A	1.15-17	12.	1.19-17
12.	3.24-1A	14.	5.29-1A	9.26-1A	14.	9.98-1A
13.	3.14-1A	16.	4.81-1A	8.09-1A	16.	9.07-1A
14.	3.26-1A	18.	2.69-1A	6.12-1A	18.	7.56-1A
15.	4.47-1A	20.	2.06-1A	5.13-1A	20.	6.80-1A
16.	4.38-1A	22.	1.64-1A	4.43-1A	22.	5.32-1A
17.	1.75-1A	24.	1.31-1A	3.75-1A	30.	4.24-1A
18.	8.14-1A	26.	1.03-1A	3.18-1A	35.	3.36-1A
19.	6.59-1A	28.	8.35-1A	2.77-1A	40.	2.72-1A
20.	5.46-1A	30.	6.78-1A	2.42-1A	50.	1.79-1A
22.	3.94-1A	32.	5.41-1A	2.08-1A	75.	6.58-1A
24.	2.85-1A	34.	4.33-1A	1.80-1A	100.	2.35-1A
26.	2.10-1A	36.	3.48-1A	1.58-1A	125.	8.25-2A
28.	1.59-1A	38.	2.80-1A	1.38-1A	150.	3.08-2A
30.	1.08-1A	40.	2.23-1A	1.18-1A	175.	1.52-2A
32.	8.19-2A	45.	1.32-1A	8.22-1A	200.	1.05-2A
34.	6.36-2A	50.	8.10-2A	5.83-1A	225.	3.38-2A
36.	3.78-2A	55.	5.06-2A	4.06-1A		

Note: "Energy (eV)" refers to secondary electron energy and the keV energies, atop the data columns, refer to the incident H^+ energies.

Accuracy: The estimated uncertainty in these data is 20% except near threshold (<10 eV) where it may be higher.

Reference: These data were taken from M. E. Rudd, L. H. Toburen, and N. Stolterfoht, Atomic Data and Nuclear Data Tables 23, 405 (1979).

Tabular Data I-2.1 (continued)
 Single Differential Cross Section (Secondary
 Electron Spectra) for $H^+ + Ar$ Collisions
 (Units of cm^2/eV)

ELECTRON ENERGY (eV)	Proton Energy (keV)							
	5	7	10	15	20	30	50	70
1.5	2.31-17	2.48-17	2.38-17	2.36-17	2.87-17	2.77-17	2.52-17	2.67-17
2.0	2.27-17	2.19-17	2.40-17	2.49-17	3.03-17	2.86-17	2.67-17	2.72-17
3.0	1.62-17	1.90-17	2.14-17	2.38-17	2.86-17	2.46-17	2.46-17	2.51-17
5.0	9.95-18	1.27-17	1.51-17	1.88-17	2.33-17	2.24-17	2.07-17	2.07-17
7.5	6.19-18	8.94-18	1.17-17	1.54-17	1.48-17	2.01-17	1.85-17	1.84-17
10.0	1.95-18	6.19-18	8.51-18	1.17-17	1.65-17	1.75-17	1.65-17	1.58-17
15.0	3.00-18	4.20-18	5.35-18	7.43-18	9.71-18	1.26-17	1.21-17	1.14-17
20.0	7.47-19	1.33-18	2.43-18	3.74-18	5.44-18	7.60-18	8.82-18	8.34-18
30.0	1.50-19	3.36-19	7.64-19	1.56-18	2.36-18	3.75-18	5.18-18	5.28-18
50.0	5.05-20	2.69-20	8.93-20	2.66-19	5.07-19	1.14-18	1.87-18	2.25-18
75.0	1.21-20	6.69-21	1.18-20	3.57-20	8.40-20	2.87-19	6.61-19	9.03-19
100.0	2.01-21	2.01-21	4.32-21	8.75-21	1.61-20	7.05-20	2.47-19	4.05-19
130.0	2.42-21	9.57-22	1.33-21	2.39-21	2.94-21	1.32-20	7.28-20	1.62-19
160.0	1.55-21	9.34-23	4.60-22	1.03-21	9.93-22	3.34-21	2.07-20	6.56-20
200.0	1.55-22	6.75-22	2.01-22	4.11-22	5.56-22	2.11-21	7.31-21	2.45-20
250.0		4.51-23	1.89-22	1.38-22	3.11-23	3.03-22	4.24-22	3.16-21
300.0			4.07-24			5.99-23	7.46-23	8.16-22

Note: "Energy (eV)" refers to secondary electron energy.

Accuracy: The estimated uncertainty in these data is 20% except near threshold (<10 eV) where it may be higher.

Reference: These data were taken from M. E. Rudd, L. H. Toburen, and N. Stolterfoht, Atomic Data and Nuclear Data Tables 23, 405 (1979).

Tabular Data I-2.1 (continued)

Single Differential Cross Section (Secondary

Electron Spectra) for $H^+ + Ar$ Collisions

(Units of cm^2/eV)

Proton Energy (keV)							
ENERGY (eV)	50	100	150	200	250	300	ENERGY (eV)
1.0	2.36-17	2.16-17	1.74-17	1.40-17	1.27-17	1.10-17	1.0
1.3	2.35-17	2.16-17	1.75-17	1.43-17	1.30-17	1.12-17	1.3
1.6	2.34-17	2.15-17	1.77-17	1.46-17	1.33-17	1.14-17	1.6
2.0	2.34-17	2.15-17	1.80-17	1.50-17	1.35-17	1.17-17	2.0
2.5	2.29-17	2.13-17	1.83-17	1.54-17	1.38-17	1.20-17	2.5
3.0	2.27-17	2.12-17	1.87-17	1.59-17	1.42-17	1.23-17	3.0
4.0	2.19-17	2.06-17	1.87-17	1.62-17	1.44-17	1.27-17	4.0
5.0	2.06-17	1.92-17	1.77-17	1.56-17	1.39-17	1.24-17	5.0
6.0	1.90-17	1.79-17	1.67-17	1.49-17	1.33-17	1.19-17	6.0
8.0	1.63-17	1.56-17	1.46-17	1.33-17	1.19-17	1.08-17	8.0
10.0	1.52-17	1.46-17	1.36-17	1.23-17	1.10-17	9.69-18	10.0
13.0	1.31-17	1.16-17	1.06-17	9.79-18	8.41-18	7.67-18	13.0
16.0	1.12-17	9.61-18	8.24-18	7.35-18	6.35-18	5.67-18	16.0
20.0	8.70-18	7.25-18	5.99-18	5.13-18	4.37-18	3.80-18	20.0
25.0	6.65-18	5.46-18	4.36-18	3.57-18	2.96-18	2.53-18	25.0
30.0	4.89-18	4.44-18	3.45-18	2.77-18	2.28-18	1.92-18	30.0
40.0	3.09-18	3.27-18	2.46-18	1.94-18	1.55-18	1.29-18	40.0
50.0	1.96-18	2.41-18	1.89-18	1.41-18	1.13-18	9.36-19	50.0
60.0	1.26-18	1.71-18	1.41-18	1.07-18	8.54-19	7.05-19	60.0
80.0	5.57-19	9.33-19	8.50-19	7.12-19	5.60-19	4.64-19	80.0
100.0	2.45-19	5.62-19	5.67-19	4.97-19	3.56-19	3.24-19	100.0
130.0	8.83-20	2.75-19	2.15-19	2.89-19	2.51-19	2.11-19	130.0
160.0	2.00-20	1.41-19	1.96-19	1.94-19	1.72-19	1.54-19	160.0
200.0	7.92-21	7.28-20	1.36-19	1.65-19	1.69-19	1.72-19	200.0
250.0	6.84-22	1.36-20	4.31-20	5.86-20	5.74-20	5.05-20	250.0
300.0	3.23-22	3.60-21	1.86-20	3.24-20	3.65-20	3.36-20	300.0
350.0	1.82-22	1.37-21	7.01-21	1.78-20	2.32-20	2.34-20	350.0
400.0	9.67-23	2.97-22	2.42-21	9.05-21	1.45-20	1.64-20	400.0
450.0	6.39-23	1.52-22	8.55-22	4.16-21	8.75-21	1.09-20	450.0
500.0	3.66-23	8.08-23	3.47-22	1.79-21	4.94-21	7.41-21	500.0
550.0	2.37-23	6.62-23	1.62-22	6.04-22	2.58-21	4.86-21	550.0
600.0	2.09-23	3.51-23	8.84-23	3.46-22	1.32-21	3.00-21	600.0
650.0	1.57-23	1.95-23	5.70-23	1.74-22	6.53-22	1.74-21	650.0
700.0	9.79-24	1.18-23	4.05-23	1.61-22	3.23-22	9.87-22	700.0
750.0	5.55-24	8.30-24	2.77-23	6.57-23	1.78-22	5.46-22	750.0
800.0	3.43-24	6.91-24	1.83-23	4.73-23	1.09-22	2.94-22	800.0
850.0	3.19-24	4.07-24	1.31-23	3.54-23	6.88-23	1.69-22	850.0
900.0	5.44-24	1.23-24	1.08-23	2.65-23	4.34-23	9.93-23	900.0
950.0	8.18-24	8.03-25	7.76-24	2.02-23	3.01-23	6.35-23	950.0
1000.0	7.36-24	2.32-24	5.17-24	1.61-23	2.44-23	4.77-23	1000.0

Note: "Energy (eV)" refers to secondary electron energy.

Accuracy: The estimated uncertainty in these data is 20% except near threshold (<10 eV) where it may be higher.

Reference: These data were taken from M. E. Rudd, L. H. Toburen, and N. Stolterfoht, Atomic Data and Nuclear Data Tables 23, 405 (1979).

Tabular Data I-2.1 (continued)

Single Differential Cross Section (Secondary

Electron Spectra) for H^+ + Ar Collisions

(Units of cm^2/eV)

0.25 Mev		0.3 Mev		0.5 Mev		1.0 Mev	
ENERGY (eV)		ENERGY (eV)		ENERGY (eV)		ENERGY (eV)	
1.	1.18-17	1.	1.36-17	1.	1.07-17	1.	6.81-18
2.	1.47-17	2.	1.43-17	2.	1.12-17	2.	6.76-18
4.	1.52-17	4.	1.42-17	4.	1.08-17	4.	6.10-18
6.	1.29-17	6.	1.26-17	6.	9.23-18	6.	5.25-18
8.	1.21-17	8.	1.12-17	8.	8.21-18	8.	4.82-18
10.	1.17-17	10.	1.00-17	10.	7.34-18	10.	4.31-18
15.	6.45-18	12.	8.62-18	15.	5.01-18	15.	2.63-18
20.	4.11-18	14.	6.96-18	20.	2.81-18	20.	1.40-18
30.	2.11-18	16.	5.73-18	25.	1.74-18	25.	8.13-19
40.	1.41-18	18.	4.69-18	30.	1.22-18	50.	2.84-19
50.	1.09-18	20.	3.89-18	40.	7.96-19	100.	1.04-19
60.	8.17-19	25.	2.53-18	50.	5.86-19	125.	7.44-20
70.	6.41-19	30.	9.59-19	75.	3.15-19	150.	6.17-20
80.	5.44-19	35.	5.26-19	100.	2.03-19	175.	6.92-20
90.	4.43-19	40.	3.30-19	150.	1.10-19	200.	1.45-19
100.	3.16-19	150.	1.73-19	200.	1.73-19	250.	2.15-20
125.	2.59-19	200.	1.79-19	250.	3.76-20	300.	1.92-20
150.	1.93-19	250.	5.55-20	300.	2.56-20	400.	8.90-21
200.	1.61-19	300.	3.66-20	400.	1.38-20	500.	5.75-21
250.	5.16-20	350.	2.57-20	500.	8.12-21	750.	2.40-21
300.	3.17-20	400.	1.83-20	750.	3.00-21	1000.	1.32-21
350.	2.45-20	500.	8.60-21	1000.	8.88-22	1250.	8.23-22
400.	1.59-20	600.	3.18-21	1200.	1.85-22	1500.	5.46-22
450.	9.54-21	700.	1.07-21	1400.	2.83-23	1750.	3.82-22
500.	5.26-21	800.	2.25-22	1600.	8.35-24	2000.	2.08-22
600.	1.24-21	900.	4.11-23	1800.	2.54-24	2500.	2.08-23
700.	2.32-22	1000.	2.42-24	2000.	2.20-26	3000.	1.06-24
800.	3.96-23					3500.	2.68-25
900.	3.87-24						

Note: "Energy (eV)" refers to secondary electron energy and the MeV energies, atop the data columns, refer to the incident H^+ energies.

Accuracy: The estimated uncertainty in these data is 20% except near threshold (<10 eV) where it may be higher.

Reference: These data were taken from M. E. Rudd, L. H. Toburen, and N. Stolterfoht, Atomic Data and Nuclear Data Tables 23, 405 (1979).

Tabular Data I-2.1 (continued)
Single Differential Cross Section (Secondary
Electron Spectra) for $H^+ + Ar$ Collisions
(Units of cm^2/eV)

ENERGY (eV)	1.5 MeV	ENERGY (eV)	2.0 MeV	ENERGY (eV)	3.0 MeV	3.67 MeV	4.2 MeV
1	5.04-18	15	1.28-18	15	1.51-18	8.87-18	1.03-18
2	5.10-18	20	9.61-19	20	8.88-18	5.51-19	4.94-18
4	4.88-18	30	3.92-19	30	2.88-18	2.23-19	1.42-19
6	4.47-18	40	2.67-19	40	2.06-18	1.56-19	1.03-19
8	4.12-18	50	2.09-19	50	1.57-18	1.23-19	7.91-20
10	3.55-18	75	1.15-19	75	8.91-20	7.08-20	4.67-20
15	2.04-18	100	7.37-20	100	5.78-20	4.52-20	2.91-20
20	1.04-18	125	5.30-20	125	4.20-20	3.22-20	2.29-20
25	6.04-19	150	4.55-20	150	3.51-20	2.84-20	2.47-20
30	4.37-19	175	5.44-20	175	4.45-20	3.68-20	2.37-20
40	3.00-19	200	2.16-19	200	9.65-20	9.26-20	4.98-20
50	2.18-19	250	1.52-20	250	1.23-20	9.08-21	7.15-21
75	1.25-19	300	1.06-20	300	8.12-21	6.16-21	5.14-21
100	8.20-20	400	6.07-21	400	4.93-21	3.52-21	2.90-21
125	5.88-20	500	3.94-21	500	3.25-21	2.23-21	1.96-21
200	1.32-19	600	2.56-21	600	2.22-21	1.96-21	1.32-21
250	1.68-20	700	1.86-21	800	1.27-21	7.72-22	7.90-22
500	4.47-21	800	1.43-21	1000	7.53-22	5.04-22	4.94-22
750	1.83-21	900	1.12-21	1250	4.77-22	3.33-22	3.04-22
1000	1.01-21	1000	8.80-22	1500	3.17-22	2.29-22	2.08-22
1250	6.36-22	1250	5.40-22	2000	1.82-22	1.16-22	1.14-22
1500	4.38-22	1500	3.75-22	2500	1.21-22	7.85-23	7.57-23
1750	3.22-22	2000	2.31-22	3000	8.14-23	5.42-23	5.21-23
2000	2.43-22	2500	1.52-22	3500	5.11-23	3.52-23	3.08-23
2500	1.51-22	3000	8.93-23	4000	3.06-23	2.15-23	2.45-23
3000	7.32-23	3500	5.71-23	4500	1.50-23	1.95-23	2.24-23
3500	1.29-23	4000	2.21-23	5000		2.63-24	1.67-23
		4500	4.80-24	5500		2.09-24	9.06-24

Note: "Energy (eV)" refers to secondary electron energy and the MeV energies, atop the data columns, refer to the incident H^+ energies.

Accuracy: The estimated uncertainty in these data is 20% except near threshold (<10 eV) where it may be higher.

Reference: These data were taken from M. E. Rudd, L. H. Toburen, and N. Stolterfoht, Atomic Data and Nuclear Data Tables 23, 405 (1979).

Tabular Data I-2.1 (concluded)
 Single Differential Cross Section (Secondary
 Electron Spectra) for $H^+ + Ar$ Collisions
 (Units of cm^2/eV)

Proton Energy (MeV)					
ENERGY (eV)	0.3	0.4	0.5	4.2	5.0
1.17	2.44-17			5.36-18	6.18-18
1.36	2.45-17			5.46-18	5.69-18
1.58	2.43-17			5.27-18	5.22-18
1.85	2.21-17			4.67-18	4.88-18
2.15	2.06-17			4.48-18	4.63-18
2.51	1.90-17			4.27-18	4.56-18
2.93	1.81-17			3.89-18	4.36-18
3.41	1.70-17			3.70-18	4.61-18
3.98	1.61-17	1.40-17	1.09-17	3.62-18	4.20-18
4.64	1.56-17	1.30-17	1.09-17	3.69-18	4.13-18
5.41	1.41-17	1.20-17	1.11-17	3.69-18	3.76-18
6.31	1.30-17	1.10-17	1.02-17	3.83-18	3.61-18
7.36	1.19-17	1.03-17	9.24-18	3.47-18	2.96-18
8.58	1.09-17	1.01-17	8.56-18	2.99-18	2.64-18
10.0	9.62-18	8.71-18	7.01-18	2.60-18	1.90-18
11.7	8.16-18	7.79-18	6.91-18	2.16-18	1.81-18
13.6	6.35-18	6.08-18	5.47-18	1.58-18	1.39-18
15.8	5.02-18	4.84-18	4.42-18	1.20-18	9.78-19
18.5	3.85-18	3.52-18	3.22-18	8.48-19	7.02-19
21.5	2.93-18	2.62-18	2.28-18	5.60-19	4.88-19
25.1	2.24-18	1.95-18	1.67-18	3.46-19	2.89-19
29.3	1.80-18	1.51-18	1.28-18	2.49-19	2.21-19
34.1	1.42-18	1.20-18	9.91-19	1.86-19	1.57-19
39.8	1.14-18	9.71-19	8.00-19	1.46-19	1.40-19
46.4	9.08-19	7.77-19	6.47-19	1.29-19	1.03-19
54.1	7.12-19	6.06-19	4.94-19	9.26-20	8.18-20
63.1	5.62-19	4.74-19	3.88-19	7.58-20	6.59-20
73.6	4.38-19	3.72-19	3.04-19	6.07-20	5.24-20
85.8	3.43-19	2.87-19	2.38-19	4.63-20	4.71-20
100.	2.66-19	2.29-19	1.78-19	3.68-20	3.46-20
117.	2.11-19	1.79-19	1.45-19	2.97-20	2.36-20
136.	1.70-19	1.41-19	1.14-19	2.40-20	2.00-20
158.	1.44-19	1.15-19	9.28-20	2.25-20	2.05-20
185.	1.14-19	1.12-19	9.27-20	2.90-20	2.86-20
215.	7.01-20	6.25-20	4.65-20	1.31-20	8.13-21
251.	4.82-20	4.38-20	3.34-20	6.67-21	5.98-21
293.	3.46-20	3.08-20	2.34-20	4.92-21	3.86-21
341.	2.44-20	2.22-20	1.67-20	3.41-21	2.78-21
398.	1.65-20	1.58-20	1.20-20	2.67-21	2.09-21
464.	1.03-20	1.11-20	8.41-21	1.75-21	1.48-21
541.	5.62-21	8.28-21	6.31-21	1.30-21	1.12-21
631.	2.44-21	5.92-21	4.58-21	1.05-21	8.38-22
736.	7.13-22	3.37-21	3.12-21	7.80-22	5.39-22
858.	1.83-22	1.36-21	1.82-21	5.84-22	3.61-22
1000.	6.32-23	3.29-22	8.22-22	4.41-22	2.94-22
1166.	3.10-23	4.87-23	2.29-22	3.44-22	2.40-22
1359.		1.20-23	5.55-23	2.53-22	1.90-22
1585.				1.63-22	1.23-22
1848.				1.11-22	7.55-23
2154.				7.28-23	5.53-23
2512.				5.70-23	5.10-23
2929.				3.96-23	3.16-23
3415.				2.42-23	
3981.				1.80-23	
4642.				1.02-23	

Note: "Energy (eV)" refers to secondary electron energy.

Accuracy: The estimated uncertainty in these data is 20% except near threshold (<10 eV) where it may be higher.

Reference: These data were taken from M. E. Rudd, L. H. Toburen, and N. Stolterfoht, Atomic Data and Nuclear Data Tables 23, 405 (1979).

I - 3. ENERGY SPECTRA OF SECONDARY ELECTRONS
FROM HEAVY - PARTICLE IMPACT IONIZATION

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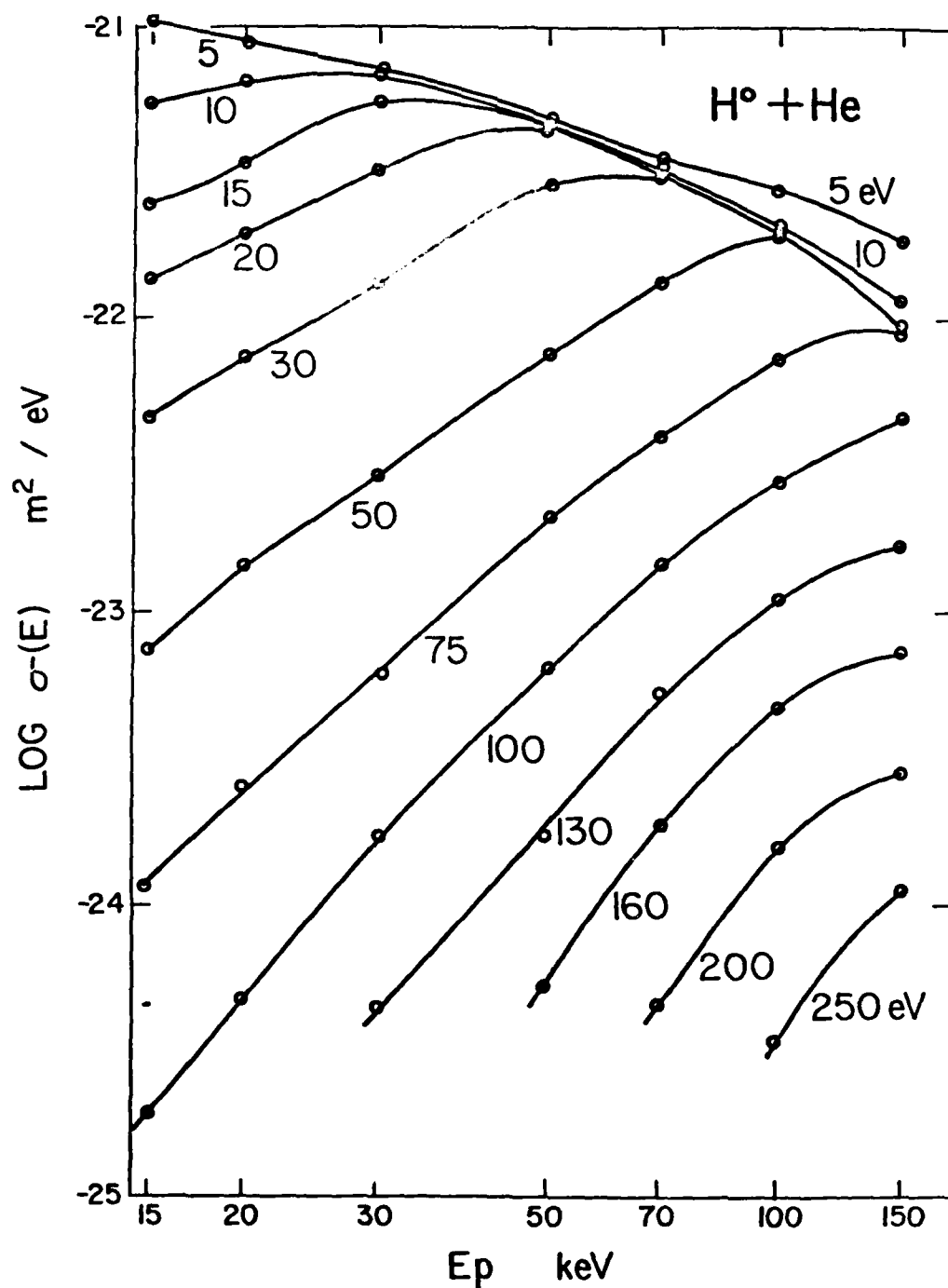
Tabular Data I-3.1
Single Differential Cross Section (Secondary
Electron Spectral) for H + He Collisions
(Units of m^2/eV)

Electron energy	15 keV	20 keV	30 keV	50 keV	70 keV	100 keV	150 keV
1.5	1.19-21	8.15-22	8.00-22	6.97-22	7.27-22	6.24-22	5.07-22
2.0	1.06-21	7.83-22	6.69-22	5.88-22	5.64-22	5.40-22	4.16-22
3.0	1.16-21	8.29-22	6.35-22	5.55-22	4.09-22	3.68-22	3.47-22
5.0	1.06-21	8.63-22	6.87-22	4.76-22	3.56-22	2.72-22	2.11-22
7.5	8.08-22	8.05-22	7.48-22	4.97-22	3.34-22	2.36-22	1.44-22
10.0	5.52-22	6.58-22	6.89-22	4.84-22	3.17-22	2.10-22	1.25-22
15.0	2.45-22	3.44-22	5.53-22	4.75-22	3.23-22	1.97-22	1.01-22
20.0	1.27-22	1.91-22	3.19-22	4.41-22	3.44-22	1.98-22	9.34-23
30.0	4.63-23	7.46-23	1.31-22	2.83-22	3.15-22	2.09-22	9.23-23
50.0	7.46-24	1.42-23	2.93-23	7.63-23	1.31-22	1.90-22	9.65-23
75.0	1.13-24	2.57-24	6.02-24	2.14-23	3.97-23	7.10-23	1.01-22
100.0	1.90-25	4.78-25	1.69-24	6.48-24	1.47-23	2.77-23	4.65-23
130.0		7.62-26	4.31-25	1.62-24	5.27-24	1.10-23	1.72-23
160.0				5.30-25	1.84-24	4.70-24	7.23-24
200.0					4.51-25	1.58-24	2.78-24
250.0					7.77-26	4.45-25	1.17-24
300.0					1.93-26	1.26-25	4.21-25

Accuracy: The quoted overall uncertainty is +17% above 30 keV increasing to +25% at 15 keV with additional uncertainties for the lower electron energies (<10 eV).

Reference: The data were taken from M. E. Rudd, J. S. Risley, J. Fryar, and R. C. Rolfes, Phys. Rev. A 21, 506 (1980).

Graphical Data I-3.2



Single differential cross section (secondary electron spectrum) for $H^{\circ} + He$ collisions. These data were taken from M. E. Rudd, J. S. Risley, J. Fryar, and R. G. Rolfes, Phys. Rev. A 21, 506 (1980).

Comments and References

In addition to the data presented in this section, there is also recent data on the energy and angular distribution of secondary electrons (double differential cross sections) in several cases, but no integration over the angle to obtain single differential cross sections was reported. The systems studied are:

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J. NUCLEAR DATA

(No new entries here. See Vol. V for data.)

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