

PHYSICAL REVIEW LETTERS

VOLUME 28

8 MAY 1972

NUMBER 19

Hydrogen-Antihydrogen Interactions*

B. R. Junker and J. N. Bardsley

Physics Department, University of Pittsburgh, Pittsburgh, Pennsylvania 15213

(Received 2 June 1971; revised manuscript received 31 March 1972)

The interaction potential between hydrogen and antihydrogen is obtained by a variational calculation. A potential maximum is found, but its peak is below the dissociation limit. A classical calculation shows that the cross section for total annihilation decreases monotonically from $31a_0^2$ at 0.01 eV to $4a_0^2$ at 20 eV.

Interest in the atomic interactions involving antimatter has been aroused by the analysis of bubble-chamber reactions and by recent cosmological speculation concerning the possible presence of equal amounts of koinomatter and antimatter in the universe. Various authors¹⁻⁴ have pointed out that the rate of proton-antiproton annihilation in a system containing atoms or antiatoms may be considerably larger than in a system with just baryons and leptons. We will show that the study of the interactions between atoms and antiatoms presents some exacting and fascinating problems for the quantum chemist.

Consider a collision between neutral hydrogen and antihydrogen atoms. As the distance R between the proton and antiproton is reduced from "infinity," the binding energy of the electron and positron is reduced. As $R \rightarrow 0$, the interaction of each lepton with the two baryons is that of a dipole with decreasing moment. Since a dipole of arbitrarily small moment cannot bind an electron or positron, the system is unstable at small separations and both leptons will be emitted. There are critical separations R_{ps} and R_F below which it is energetically more favorable for the leptons to leave the baryons, either as positronium ($R < R_{ps}$) or as a free pair ($R < R_F$). Because of the binding energy of positronium, R_{ps} is greater than R_F .

For slow collisions a simple adiabatic model

can be constructed. If, during the collision, the baryon separation becomes less than R_{ps} , the leptons will be emitted as positronium. Since the binding energy of the lepton pair is now 6.8 eV, in the contrast to the 27.2 eV initially, the nuclear kinetic energy must be reduced by at least 20.4 eV. Hence if the initial kinetic energy is less than 20.4 eV, a bound state of protium will be formed. If the leptons escape while R is close to R_{ps} , they will leave with little kinetic energy and the protium will be formed in one of the highest states consistent with energy conservation. This means a principal quantum number of 24 or 23. The levels of different l and m will be populated with roughly equal probability. After formation, the protium will lose energy by radiation or further collisions until annihilation occurs from a level with small n and l values. In this model R_F plays no role since by the time the nuclear separation reaches R_F , both leptons will have left. In the absence of further collisions the positronium will also suffer annihilation shortly after its formation.

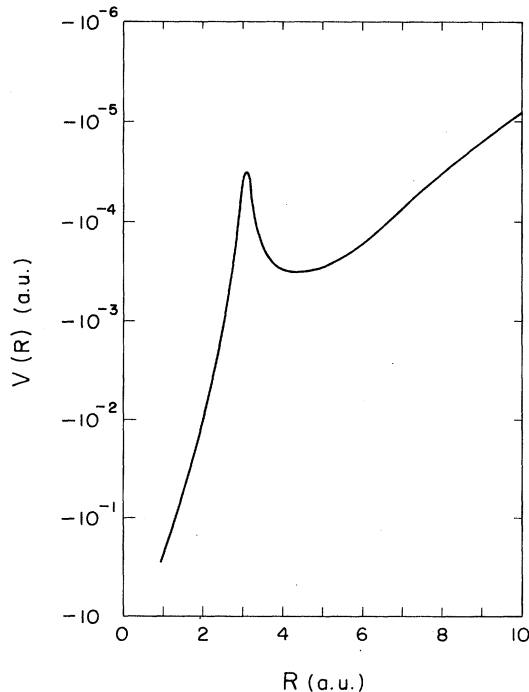
The annihilation cross section then depends only on the probability that R will become smaller than R_{ps} at some time during the collision. Morgan and Hughes¹ estimated this probability with the assumption that the H- \bar{H} interaction is completely attractive. However, Puget² has shown that there may be a potential barrier since the

TABLE I. Energy of the H- \bar{H} system with the nuclei at a fixed separation R .

R (a.u.)	H- \bar{H} interaction energy (a.u.)	$e^- - e^+$ binding energy (a.u.)
0.975	-0.277 201	0.001 560
1.0	-0.257 076	0.007 076
1.5	-0.055 216	0.138 549
2.0	-0.010 219	0.260 219
2.5	-0.001 244	0.351 244
3.0	-0.000 040	0.416 707
3.05	-0.000 033	0.422 165
3.10	-0.000 031	0.427 451
3.15	-0.000 039	0.432 579
3.5	-0.000 165	0.464 451
4.0	-0.000 301	0.500 301
5.0	-0.000 280	0.550 280
6.0	-0.000 154	0.583 487
10.0	-0.000 008	0.750 008

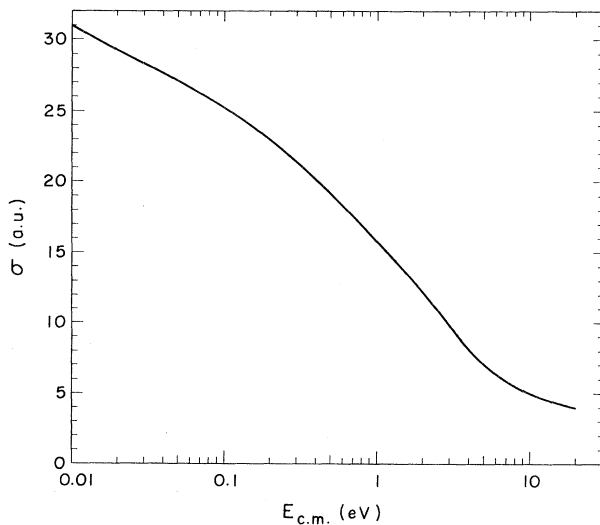
overlap of the lepton charge clouds leads to a repulsive interaction, in contrast to the attractive interaction produced for H-H. Using simple wave functions, Puget obtained a barrier of height 0.8 eV near $1.8a_0$. Such a barrier would clearly reduce the annihilation cross section significantly.

We have therefore calculated the H- \bar{H} interaction by the variational method using a trial wave function with up to 75 configurations. Each configuration is made up from a product of orbitals,

FIG. 1. Interaction energy between H and \bar{H} .TABLE II. The H- \bar{H} interaction energy at $R=3.1$ a.u.

No. of terms in the wave function	Interaction energy (a.u.)
25	+0.000 700
47	+0.000 097
60	+0.000 034
75	-0.000 031

which are either Slater orbitals based on one of the nuclei or standard elliptical orbitals⁵ based on both nuclei. The nonlinear orbital parameters, one for each Slater orbital and two for each elliptical orbital, were optimized to give the lowest energy with a wave function of 21 or 25 terms. The results of the 75-term calculation are shown in Table I and Fig. 1. At large R the potentials shows the attractive Van der Waals interaction, and at small R the $p-\bar{p}$ Coulomb interaction produces a strong attraction. Between these regions there is a peak at 3.089 a.u., as determined by a cubic spline curve fitted through the points in Table I. The height of the peak is 3.1×10^{-5} a.u. (0.85 meV) below the dissociation limit. The binding energy of the lepton pair reduces to almost zero at 0.975 a.u., suggesting that $R_{ps} \approx 0.96$ a.u. The rate of convergence of the calculation can be judged from Table II. It should be noted that in the 60-term calculation the energy at 3.1 a.u. was slightly higher than the dissociation limit. Thus, it was necessary to go to the 75-term function in order to demonstrate that the

FIG. 2. Cross section for the formation, with subsequent annihilation, of protium and positronium in H- \bar{H} collisions.

potential maximum does not represent a barrier to zero-energy collisions with zero angular momentum.

The cross section for the formation of positronium and protium pairs, calculated using the adiabatic model together with classical trajectories, is shown in Fig. 2. The cross section decreases monotonically from $31a_0^2$ at 0.01 eV to $4a_0^2$ at 20 eV. These results should be accurate over the higher energies in this range, but quantum effects on the nuclear motion may influence the results at low energies.

The maximum in the $H-\bar{H}$ potential is probably a feature common to all interaction potentials between neutral atoms and antiatoms. This suggests the possibility that for some pairs the potential at the maximum might be greater than the dissociation energy, thus providing a potential barrier which could keep matter and antimatter apart. This could perhaps give some support to the recent speculation⁶ that ball lightning is caused by the congregation of large amounts of antimatter in the atmosphere. Preliminary studies on $He-\bar{H}$, with 7-term wave functions, reveal a small barrier which will probably disappear in

more accurate calculations. Our results for systems in which one or other atom is charged show no potential barriers.

We are very grateful to the National Aeronautics and Space Administration Goddard Space Center for providing the computational facilities for this calculation, to Dr. D. L. Morgan for introducing us to the problem, and to Dr. D. L. Morgan and Dr. J. Bredekamp for running our program.

*Supported by the National Aeronautics and Space Administration under Contract No. NGL-39-011-035.

¹D. L. Morgan and V. Hughes, *Phys. Rev. D* **2**, 1389 (1970).

²J. L. Puget, CERN Report No. TH1201, 1970 (unpublished), and *Nature (London)*, *Phys. Sci.* **230**, 173 (1971).

³G. Steigman, thesis, New York University, 1968 (unpublished), and *Nature (London)* **224**, 477 (1969).

⁴See also A. S. Wightman, *Phys. Rev.* **77**, 521 (1950).

⁵J. C. Browne and F. A. Matsen, *Phys. Rev.* **135**, A1227 (1964).

⁶D. E. T. F. Ashby and C. Whitehead, *Nature (London)*, *Phys. Sci.* **230**, 180 (1971).

Metastable Autoionizing States of Highly Excited Heavy Ions*

I. A. Sellin, D. J. Pegg, and P. M. Griffin

University of Tennessee, Knoxville, Tennessee 37916, and Oak Ridge National Laboratory, Oak Ridge, Tennessee 37830

and

Winthrop W. Smith

University of Connecticut, Storrs, Connecticut 06268

(Received 20 March 1972)

We report the respective electronic energies [-5.658 ± 0.024 and -6.369 ± 0.033 keV] and electron emission lifetimes [0.91 ± 0.04 and 0.66 ± 0.04 nsec] of the longest-lived autoionizing quartet states of lithiumlike Cl^{14+} and Ar^{15+} ions, firmly establishing the energy and lifetime scaling with Z . We discuss observable relativistic corrections to non-relativistic variational energy calculations. Production of metastable states belonging to lower charge states is discussed, and it is concluded that many excited heavy ions should exhibit them.

Sufficient numbers of metastable Cl^{14+} and Ar^{15+} ions have been created by passage of chlorine and argon beams through $15\text{-}\mu\text{g}/\text{cm}^2$ carbon foils to render possible the measurement of the total electronic energies and autoionization lifetimes of the lowest, longest-lived, nonradiative quartet states [$(1s2s2p)^4P_{5/2}^o$] of these three-electron ions. 41-MeV chlorine beams from the Oak

Ridge tandem accelerator and 87- and 60-MeV argon beams from the Oak Ridge isochronous cyclotron were used for these experiments. In each case a few tenths of 1% of the emergent three-electron beam fraction has been found to exist in the lowest quartet state having highest angular momentum ($J = \frac{5}{2}$); it is then probable that $\geq 1\%$ of the three-electron beam fraction is