Comparison of measured and calculated Stark broadening parameters for neutral-helium lines

M. S. Dimitrijevic

Institute of Physics, P.O. Box 57, 11001 Beograd, Yugoslavia

S. Sahal-Bréchot Observatoire de Paris, 92190 Meudon, France (Received 27 February 1984)

Stark broadening parameters for neutral-helium lines obtained using a semiclassical perturbation formalism are compared with critically selected experimental data. Our data are also compared with the results of Benett and Griem as well as with the semiclassical convergent calculations of Bassalo, Cattani, and Walder.

I. INTRODUCTION

Recently, we have calculated¹ electron and proton impact linewidths and line shifts of 56 neutral-He lines in the ultraviolet, visible, and infrared regions of the spectrum using a semiclassical perturbation formalism.^{2,3} The version of semiclassical approach used in our calculations differs in several respects from the other large-scale calculations $4-7$ of neutral-He lines. Since He is a simple atomic system and since many experimental Stark broadening data are available for He lines, $8,9$ the data we obtained provide an opportunity for testing various approximations included in the semiclassical perturbation formalism. Our data as well as the results of Benett and Griem⁴ and semiclassical convergent calculations of Bassalo et $al.^7$ are compared here with the critically selected^{8,9} experimental $data^{10-22}$ in order to test various sets of assumptions involved in different approaches.

II. THEORY

We calculated electron-impact broadening parameters of isolated neutral He lines using the semiclassical perturbation formalism.^{2,3} The relations for (full) half width (w) and shift (d) in the framework of the abovementioned approach are

$$
w = N \int_0^\infty v f(v) dv \left[\sum_{j \neq i} \sigma_{ij}(v) + \sum_{j' \neq f} \sigma_{fj'}(v) + \sigma_{el} \right], (1)
$$

$$
d = N \int_0^\infty v f(v) dv \int_{R_3}^{R_d} 2\pi \rho \sin(2\phi_p) d\rho . \tag{2}
$$

The inelastic cross section $\sigma_{ij}(v)$ can be expressed by an integration over the impact parameter of the transition probabilities P_{ij} as

$$
\sum_{j\neq i} \sigma_{ij}(v) = \frac{1}{2}\pi R_1^2 + \int_{R_1}^{R_d} 2\pi \rho \, d\rho \sum_{j\neq i} P_{ij}(\rho, v) \,. \tag{3}
$$

Using the adiabatic approximation, the elastic cross section is

$$
\sigma_{\rm el} = 2\pi R_2^2 + \int_{R_2}^{R_d} 8\pi \rho \, d\rho \sin^2\delta \;, \tag{4}
$$

$$
\phi = (\phi_p + \phi_q) ,
$$

\n
$$
\phi_p = \sum_{j \neq i} \phi_{ij} - \sum_{j' \neq f} \phi_{fj'} .
$$
\n(5)

Here, i and f denote, respectively, the initial and final levels and j and j' the corresponding perturbing levels. The phase shifts ϕ_p and ϕ_q , due respectively to the polarization ($-r^{-4}$) and to the quadrupole potential ($-r^{-3}$), are given in Ref. 2 (Sec. 3, Chap. 2).

All of the cutoffs (R_1, R_2, R_3, R_d) are described in Ref. 3 (Sec. 1, Chap. 3). We recall here that we have "symmetrized" the inelastic cross section by replacing v^2 by its modified value due to collisions, i.e., $v^2 - (\hbar \omega_{ii'}/m)$, where $\hbar \omega_{ji'}$ is the threshold energy.

In terms of the S matrix for a collision of a single electron with the atom, one has

$$
\sum_{i,j'\neq j} (P_{jj'}+P_{jj})=2\{1-S_iS_f^{-1}\}_{\text{av}},\tag{6}
$$

where the curly brackets $\{\}_\text{av}$ indicate an angular average.
Griem *et al*.⁴ as well as Sahal-Bréchot^{2,3} adopted a perturbative method based on the series expansion of the Smatrix elements. Griem et al .⁴ neglected the elastic term P_{ij} and took into account elastic collisions via the strong collision term. The low cutoff is determined by the condition

$$
\{1-S_iS_f^{-1}\}_{av}\approx(\frac{3}{4})^{3/2}
$$

or

$$
\sum_{j,j'\neq j} (P_{jj'}+P_{jj}) \approx 1.3 ,
$$

while in Refs. 2 and 3 the lower cutoff for inelastic collisions (R_1) is determined from the condition goinsions (**R**)
 $\sum_{j' \neq j} P_{jj'} = \frac{1}{2}$ and the corresponding one for elastic collisions (R_2) from the condition $P_{jj} = 2$.

The "convergent approach" is based on the analogy between the Dyson-series and Taylor-series expansion of e^x . In this case the divergence for small impact parameters

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does not exist and lower cutoffs are not needed. This approach, developed by Vainshtein and Sobel'man²³ in a two-level approximation, was applied with some modifications by Dyne and O'Mara²⁴ and, recently, in the many-level approximation by Cattani, Yamamoto, Bassalo, and Walder^{7,25,26} to calculate Stark broadening parameters of neutral-helium lines. In Refs. ¹—³ ions had been treated in the impact approximation. To take the ion contribution into account, here we use the approach developed by Griem et al.⁴ and Griem,²⁷ due to its simple applicability to different kinds of perturbing ions. In terms of the parameters A and $R,$ ^{4,27} total width and shift are given by

$$
W_{\text{tot}} = [1 + 1.75A(1 - 0.75R)]w,
$$

 $d_{\text{tot}} = d \pm 2.0A (1 - 0.75R)w$.

The sign, in the shift equation is equal to that of the lowvelocity limit of d.

III. COMPARISON OF MEASURED AND CALCULATED VALUES

Our results' as well as the calculations of Benett and Griem⁶ and Bassalo *et al.*,⁷ are compared with the critically selected^{8,9} experimental data.¹⁰⁻²² The criteria for selection of experimental papers were as follows:^{8,9} (a) an independent and accurate determination of the plasma electron density N_e , (b) a reasonably accurate determination of the electron temperature, (c) a discussion of other interfering broadening mechanisms and appropriate ex-

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COMPARISON OF MEASURED AND CALCULATED STARK . . .

TABLE IV. Average accuracy of different theoretical methods compared to Stark width and shift experimental data for helium lines. The results in parentheses are obtained by excluding the $2p^3D-3d^3D$ line which exhibits a strong unexplained difference between d_m and the calculated shift (especially for d_{DSB} and d_{BG}).

	All experiments included	Experiments with C and D accuracy excluded
$(W_m/W_{\text{DSB}})_{\text{av}}$	1.17 ± 0.04	1.17 ± 0.02
$(W_m/W_{\text{RCW}})_{\text{av}}$	1.07 ± 0.04	1.07 ± 0.04
$(W_m/W_{BG})_{av}$	0.92 ± 0.04	0.93 ± 0.02
$(d_m/d_{\rm DSR})_{av}$	1.20 ± 0.13	$1.13 + 0.03$
	(1.07 ± 0.04)	
$(d_m/d_{\rm BCW})_{\rm av}$	1.23 ± 0.08	1.34 ± 0.09
	(1.27 ± 0.07)	
$(d_m/d_{\text{BG}})_{\text{av}}$	1.14 ± 0.07	$1.14 + 0.03$
	(1.07 ± 0.04)	

perimental problems. Total uncertainties (in electron density and Stark width or shift measurements) are subdivided into four ranges and coded by letters. $8,9$ The letters represent the following uncertainties: A , within 15%; B ,

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within 30%; C, within 50%; D, larger than 50%. Further differentiations are made by singling out slightly better pieces of data among similar ones by assigning a plus sign $(+)$ to indicate the first choice.^{8,9} Key data from selected experiments are given in Table I. Comparisons between various calculations and experimental data are given in Tables II and III for widths and shifts, respectively.

The averaging is first performed within a multiplet and then the obtained values are averaged over the number of multiplets. The average values of the ratios of measured to calculated linewidths and line shifts are presented in Table IU.

The calculations of Bassalo et al .⁷ and of Benett and Griem $⁶$ give, on the average, the better agreement with</sup> linewidth measurements than the present calculations, while the results of Benett and Griem $⁶$ and the present</sup> calculations are in better agreement with experimental shifts. Since the assumed uncertainty of the semiclassical approach is within $20\%^{27}$ and the typical uncertainties of experimental data are within 30%, one can conclude that all three variants of the semiclassical approach are successful in the calculation of the Stark broadening parameters.

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