

NEW RESULTS ON ELECTRON BROADENING OF SOME UV LINES OF NII, CII/IV AND SiII/III/IV

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Abstract. The electron damping constants of some UV lines of astrophysical interest are calculated within the impact and semiclassical approximations and are compared with experimental results where these are available (N II, UV(1) and C II, UV(1)).

The good agreement between experimental results and the semi-classical calculations of Sahal-Brechot (1969a, b) and Chapelle and Sahal-Brechot (1970) has led us to perform other calculations on UV lines of astrophysical interest. First we give a short account of our method. The widths are calculated within the impact approximation (Baranger, 1958). Under this condition the profile is Lorentzian and the width is given by the quantum formula (77c) of Baranger (1958). The validity criterion is

$$\gamma\tau \ll 1$$

where γ is the width of the line (damping constant) and τ a typical collision time. We have shown (Sahal-Brechot, 1969a, b; Chapelle and Sahal-Brechot, 1970) that the impact approximation is valid for both electron and proton collisions in stellar atmospheres. For this reason we have also calculated the damping constants for proton and He^+ collisions.

The scattering amplitudes and cross-sections which interfere in the calculation of the width are evaluated within the semi-classical approximation (Alder *et al.*, 1956); first, the perturber is assumed to be a classical particle moving along a classical path unperturbed by the collision; for radiating ions this path is a hyperbola. This is the main improvement compared with the Baranger-Griem calculations (Griem *et al.*, 1962; Griem, 1964); the use of a hyperbolic path instead of a straight one increases the resulting widths by a factor between 3 and 5. Secondly, the second order perturbation theory for the interaction electrostatic potential is used; dipole inelastic interactions, dipole (polarization) and quadrupole (20% of the total width) elastic interactions are taken into account; the quadrupole inelastic interactions are negligible; the inelastic interactions are important only for electron collisions. Symmetrized cross-sections (Alder *et al.*, 1956; Seaton, 1962) are used; the impact parameter cut-off is chosen in a way which copes with the conservation of the flux of particles, i.e. the *S*-matrix is unitary (Seaton, 1962).

The semi-classical approximation is obviously valid when most of the collisions are weak, i.e. the close (or strong) collisions are not important. The validity criterion is the following;

$$\Delta E/kT \ll 1$$

where ΔE is a typical energy gap between the interacting levels of the radiating atom and kT is the temperature. Nevertheless the comparisons between semi-classical calculations, experiments and the more recent elaborate quantum calculations (Bely and Griem, (1970) for MgII 3s-3p, Barnes and Peach (1970) for CaII 4s-4p) show that the semi-classical calculations are accurate to within 30%, despite the importance of the strong collisions (75% in the case of $Mg^+ 3s-3p$ where $\Delta E/kT > 4$). This agreement probably comes from the care taken with the physical nature of the model (use of symmetrized cross-sections and a unitary S-matrix).

From the results presented in this paper it can be seen that our calculations for the UV resonance lines CII (1) and NII (1) (Figures 1 and 2) fit the recent experiments of Fortna (1969) and yet the typical energy gap for these two lines is of the order of 10 eV! The close collisions are predominant and the typical impact parameter ρ_{eff} is very small (see Table I). Table I shows also that the impact approximation is valid for both electrons and protons: γ is the calculated (or experimental) value of the damping constant for a density, $N_e = 10^{15} \text{ cm}^{-3}$, and $T = 2 \times 10^4 \text{ K}$; the typical collision

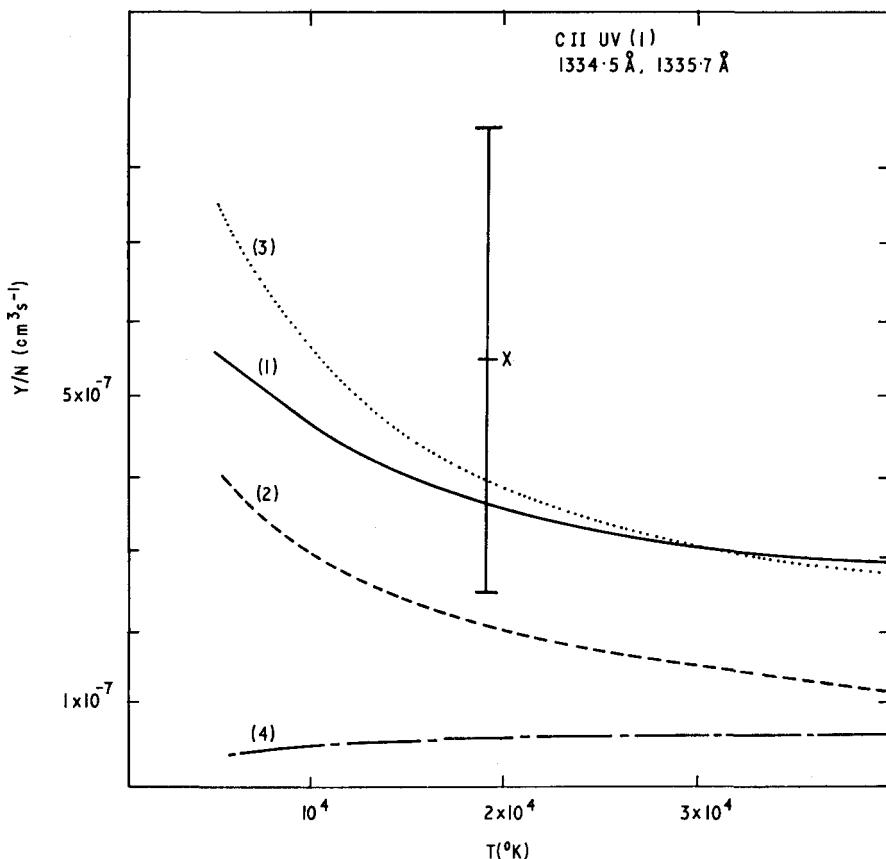


Fig. 1.

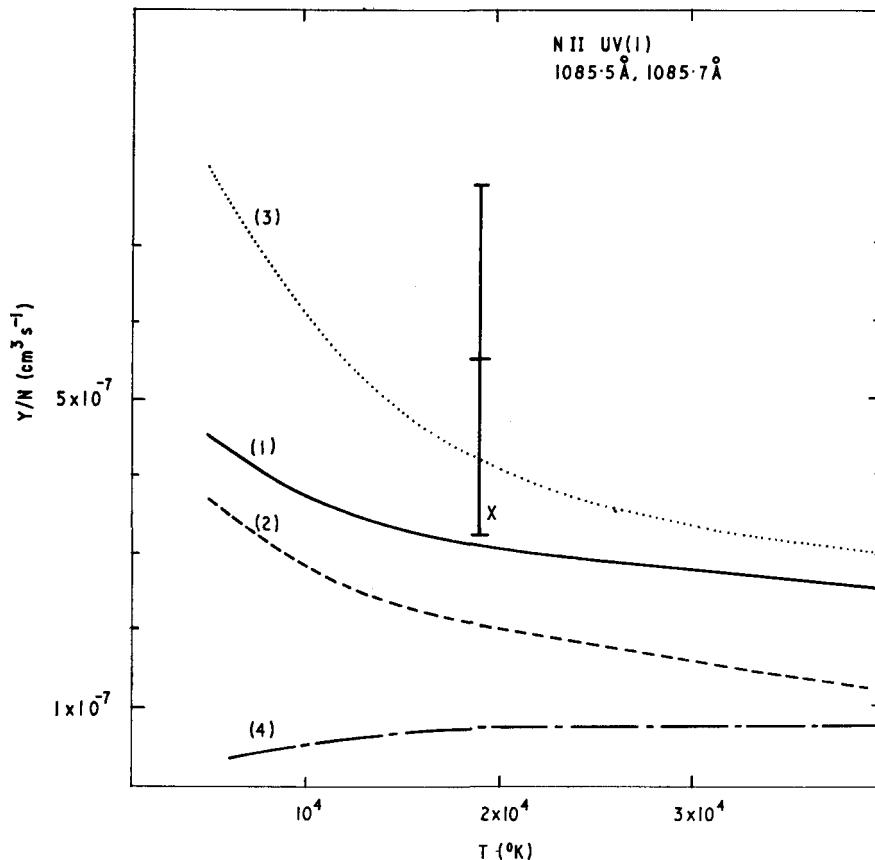


Fig. 2.

Fig. 1-2. Electron Damping constant of CII, UV (1) and NII, UV (1) (for a density of 1 cm^{-3}) versus the electronic temperature. Curve (1) this work. (2) this work, elastic collisions only. (3) Formula (102) of Brechot and Van Regemorter (1964) (No (1) of this paper). (4) Lindholm (1942). \times semi-empirical (Griem, 1968). Classical damping constant: experiment of Fortna (1969)

$$\begin{cases} 1.2 \times 10^9 \text{ for CII (1)} \\ 1.9 \times 10^9 \text{ for NII (1)} \end{cases}$$

TABLE I

Q_{eff}/a_0	$\Delta E \text{ eV}$		close coll. contrib.	l		$\gamma\tau$	
	electrons	protons		electrons	protons	electrons	protons
CII UV (1) 6	10	9	80%	1 to 2	125	3×10^{-7}	2×10^{-5}
NII UV (1) 4	6	11	80%	1	75	2×10^{-7}	1×10^{-5}

Q_{eff} typical impact parameter for $T = 2 \times 10^4 \text{ K}$

a_0 Bohr radius

ΔE typical energy difference between the interacting levels

l typical angular momentum of the perturber

γ damping constant for $N_e = 10^{15} \text{ cm}^{-3}$ and $T = 2 \times 10^4 \text{ K}$

τ typical collision time for $T = 2 \times 10^4 \text{ K}$

TABLE II
Damping constant in $\text{cm}^3 \text{sec}^{-1}$ (the damping constant is proportional to the density N) for a density of 1 cm^{-3} as a function of the temperature

Multiplet	$T(10^4 \text{ K})$	$\gamma_{e/N_e} \times 10^6$	Elastic collis. contrib. %	Typical $kT/\Delta E$	$\gamma_{eI/N} \times 10^6$	$\gamma_{L/N_e} \times 10^6$	$\gamma_{\text{B-VR}/N_e} \times 10^6$	$\gamma_{\text{H}^+/N_{\text{H}}^+} \times 10^7$	$\gamma_{\text{He}^+/N_{\text{He}}^+} \times 10^7$
$2s\ 2p^3\ 2D - 2s^2\ 2p^2\ 3P$ (1084 Å)	0.5	0.5	80	0.04	1.89	10^{-2}	0.8	10^{-3}	10^{-2}
	1	0.3	70	0.08		0.1	0.6	10^{-2}	10^{-2}
	2	0.3	65	0.16		0.1	0.4	10^{-2}	10^{-2}
	4	0.2	55	0.32		0.1	0.3	10^{-1}	10^{-1}
N II									
$2s\ 2p^3\ 2D - 2s^2\ 2p^3\ 2P^0$ (1335.7 Å)	0.5	0.6	75	0.04	1.25	10^{-2}	0.8	10^{-2}	10^{-2}
	1	0.5	65	0.08		10^{-2}	0.6	10^{-2}	10^{-1}
	2	0.3	55	0.16		0.1	0.4	10^{-1}	10^{-1}
	4	0.3	50	0.32		0.1	0.3	10^{-1}	10^{-1}
C III									
$2p^2 P - 2s^2 S$ (1549 Å)	0.5	1.3	85	0.05	0.9	1.0	2.0	10^{-3}	10^{-3}
	1	0.8	80	0.10		1.0	1.4	10^{-3}	10^{-2}
	2	0.5	80	0.20		1.0	1.0	10^{-2}	10^{-2}
	4	0.4	75	0.40		1.0	1.0	10^{-2}	10^{-1}
C IV									
$4f^2 F - 3d^2 D$ (1168.9 Å)	0.5	5.2	70	0.1	1.6	0.3	5.6	0.1	0.1
	1	4.3	60	0.2		0.3	4.0	0.1	1
	2	3.3	55	0.4		0.4	2.8	1	1
	4	2.7	50	0.8		0.4	2.0	1	1
Si III									
$3s^2\ 3d^2 D - 3s^2\ 3p\ 2p^0$ (1262.9 Å)	0.5	4.4	100	0.07	1.39	2.2	7.9	3.0	3.0
	1	3.3	100	0.14		2.5	5.6	4.0	4.0
	2	2.7	100	0.28		2.8	3.9	6.0	6.0
	4	2.4	85	0.56		3.1	2.8	8.0	7.0

Table II (continued)

Multiplet	$T(10^4 \text{K})$	$\gamma_{e/N_e} \times 10^6$	Elastic collis. $kT/\Delta E$	Typical contrib. %	$\gamma_{el/N} \times 10^5$	$\gamma_{ULN_e} \times 10^6$	$\gamma_{B-VR/N_e} \times 10^6$	$\gamma_{H^+/N^{He^+}} \times 10^6$
$3s\ 3p^2\ ^2S-3s^2\ 3p\ ^2P^0$ (1306.8 Å)	0.5	1.1	100	0.07	1.30	0.1	1.4	10^{-2}
	1	0.8	100	0.14		0.1	0.9	10^{-2}
	2	0.5	100	0.28		0.1	0.7	10^{-1}
	4	0.5	80	0.56		0.2	0.5	10^{-1}
$3p^3\ ^2P^0-3s\ 3p^2\ ^2S$ (1485.2 Å)	0.5	4.4	100	0.07	1.01	0.2	0.7	0.9
	1	3.3	100	0.14		0.3	1.5	1.0
	2	2.5	100	0.28		0.3	1.0	1.1
	4	2.1	100	0.56		0.3	0.7	1.2
$3s^2\ 4s\ 3^2S-3s^2\ 3p\ ^2P^0$ (1530.0 Å)	0.5	3.3	100	0.2	0.95	1.0	4.8	10^{-2}
	1	2.3	95	0.4		1.1	3.4	10^{-1}
	2	2.0	75	0.8		1.2	2.4	10^{-1}
	4	1.8	50	1.6		1.4	1.7	10^{-1}
$3s\ 3p^2\ ^2D-3s^2\ 3p\ ^2P^0$ (1813 Å)	0.5	1	100	0.05		~ 0.1	1.0	10^{-2}
	1	0.7	100	0.1	0.67	~ 0.1	0.7	10^{-2}
	2	0.5	100	0.2		~ 0.1	0.5	10^{-2}
	4	0.3	100	0.4		~ 0.1	0.3	10^{-1}
$6f\ ^2F-3d\ ^2D$ (2501.4 Å)	0.5	71.5	32	15	0.35	68.4	61.8	17.6
	1	85.4	20	30		76.8	43.7	21.3
	2	93.6	15	60		86.2	30.9	26.3
	4	95.3	10	120		96.8	21.9	25.9
$5f\ ^2F-3d^2\ D$ (2905 Å)	0.5	28.8	40	10	0.26	23.7	32.7	5.6
	1	33.4	25	20		26.6	23.1	6.9
	2	35.9	20	40		29.9	16.4	8.7
	4	36.4	15	80		35.5	11.6	9.3
Sum								
$3s\ 4s\ ^3S-3s\ 3p\ ^3P^0$ (995 Å)	0.5	2.5	100	0.15	2.24	0.6	6.4	10^{-3}
	1	1.7	100	0.30		0.7	4.5	10^{-2}
	2	1.3	90	0.60		0.8	3.2	~ 0.1
	4	1.8	50	1.20		0.9	2.3	0.1

Table II (continued)

Multiplet	$T(10^4 \text{ K})$	$\gamma_{e/N_e} \times 10^6$	Elastic collis. contrib. %	Typical $kT/\Delta E$	$\gamma_{el/N} \times 10^5$	$\gamma_{B-VR/N_e} \times 10^6$	$\gamma_{H^+/N_e H^+} \times 10^7$	$\gamma_{He^+/N_e He^+} \times 10^7$
$3s\ 3d^3D - 3s\ 3p\ ^3P^0$ (1111 Å)	0.5	1.7	100	0.05	1.8	0.1	2.4	10^{-1}
	1	1.2	100	0.1		0.1	1.7	10^{-1}
	2	0.9	90	0.2		0.1	1.2	10^{-1}
	4	0.7	90	0.4		0.2	0.8	~ 1
$3s\ 3p\ ^1P^0 - 3s^2\ ^1S$ (1206.5 Å)	0.5	1.5	100	0.04	1.53	~ 0.1	1.7	10^{-1}
	1	1.1	90	0.08		~ 0.1	1.2	10^{-1}
	2	1.1	60	0.16		~ 0.1	0.9	10^{-1}
	4	0.8	60	0.32		~ 0.1	0.6	10^{-1}
$3p^2\ ^3P - 3s\ 3p\ ^3P^0$ (1296 Å)	0.5	3.0	100	0.04	1.32	0.3	4.0	10^{-1}
	1	2.0	95	0.08		0.3	2.9	10^{-1}
	2	1.8	75	0.16		0.4	2.0	10^{-1}
	4	1.3	75	0.32		0.4	1.4	~ 1
$3d^1\ ^1D - 3p\ ^1P^0$ (1206.5 Å)	0.5	3.3	70	0.04	1.53	0.4	4.9	10^{-1}
	1	2.6	60	0.08		0.4	3.5	10^{-1}
	2	2.2	55	0.16		0.5	2.5	~ 1
	4	1.8	45	0.32		0.6	1.7	~ 1
$3p^2\ ^1D - 3p\ ^1P^0$ (2541.8 Å)	0.5	1.8	80	0.04	0.34	0.01	1.9	10^{-2}
	1	1.4	70	0.08		0.01	1.3	10^{-1}
	2	1.1	65	0.16		0.1	0.9	10^{-1}
	4	0.8	60	0.32		0.1	0.7	10^{-1}
SiIV								
$3p\ ^2P - 3s\ ^2S$ (1397 Å)	0.5	2.0	85	0.05	1.1	10^{-2}	2.0	10^{-2}
	1	1.0	80	0.10		10^{-2}	1.0	10^{-2}
	2	1.0	70	0.20		10^{-2}	1.0	10^{-1}
	4	1.0	65	0.40		10^{-1}	1.0	10^{-1}
$4f^2F - 3d^2D$ (1066.6 Å)	0.5	6.0	60	0.1	1.9	10^{-1}	6.0	10^{-1}
	1	5.0	50	0.2		10^{-1}	4.0	10^{-1}
	2	4.0	40	0.4		10^{-1}	3.0	1
	4	3.0	35	0.8		10^{-1}	2.0	1

Table II (continued)

Multiplet	$T(10^4 \text{ K})$	$\gamma_{e/N_e} \times 10^5$	Elastic collis. contrib. %	$kT/\Delta E$	$\gamma_{el/N} \times 10^5$	$\gamma_{L/N_e} \times 10^5$	$\gamma_{B-VR/N_e} \times 10^5$	$\gamma_{H^+/N_{H^+}} \times 10^5$	$\gamma_{He^+/N_{He^+}} \times 10^5$
$3d^2D-3p^2P$ (1125 Å)	0.5	0.2	75	0.04	1.7	10^{-3}	0.2	10^{-4}	10^{-4}
	1	0.2	60	0.08		10^{-3}	0.1	10^{-4}	10^{-3}
	2	0.1	60	0.16		10^{-3}	0.1	10^{-3}	10^{-3}
	4	0.1	55	0.32		10^{-2}	0.1	10^{-3}	10^{-3}
$6f^2F-4d^2D$ (1533.2 Å)	0.5	4.5	30	0.9	0.9	6.0	13.8	0.88	0.86
	1	4.9	20	1.8		6.7	9.7	1.4	1.2
	2	4.4	15	3.6		7.6	6.9	1.7	1.4
	4	3.4	15	7.2		8.5	4.9	2.2	1.9
$6g^2G-4f^2F$ (1672.6 Å)	0.5	5.4	20	50	0.79	12.6	21.5	2.3	2.1
	1	4.3	20	100		14.1	15.2	3.0	2.6
	2	3.2	20	200		15.9	10.7	3.9	3.3
	4	2.3	20	400		17.8	7.6	4.8	4.2
$6p^2P-4d^2D$ (1796.8 Å)	0.5	2.2	75	0.9	0.69	0.3	2.1	0.1	0.1
	1	2.0	60	1.8		0.3	1.5	0.1	0.1
	2	1.7	50	3.6		0.3	1.0	0.2	0.2
	4	1.5	40	7.2		0.4	0.7	0.2	0.2

 T electron temperature γ_{e/N_e} electron damping constant (our semi-classical results) $\gamma_{cl/N}$ classical damping constant for $N = 10^{14} \text{ cm}^{-3}$ γ_{L/N_e} Lindholm (1941) formula γ_{B-VR/N_e} Formula (102) of Brechot and Van Regemorter (1964) (Formula (1) of this paper) $\gamma_{H^+/N_{H^+}}$ proton damping constant Cour semi-classical results $\gamma_{He^+/N_{He^+}}$ damping constant by He^+ collisions (our semi-classical results).

time, τ is of the order of ϱ_{eff}/v ; the velocity of the perturber v is of the order of $\sqrt{(8kT)/(\pi m)}$ where m is the reduced mass. The typical impact parameter ϱ_{eff} is evaluated *a posteriori* (Chapelle and Sahal-Brechot, 1970) in the following manner: first, the cross-section σ is evaluated and then the predominant impact parameter ϱ_{eff} is deduced (one can write $\sigma = \pi \varrho_{\text{eff}}^2$ for instance).

From Figures 1 and 2 it can be seen that the semi-empirical formula of Griem (1968) and the simple formula of Brechot and Van Regemorter (1964), analogous to the Lindholm (1943) adiabatic formula but with a hyperbolic classical path,

$$\gamma = N_e 3.86 \times 10^{-5} Z^{4/5} |\alpha_i - \alpha_f|^{2/5} T^{-1/2} \quad (1)$$

also give reliable results for these two lines. N_e is the electron density in cm^{-3} , Z is the charge of the radiating ion, α_i and α_f are the polarisabilities, in atomic units, of the radiating ion in the stage i or f , the initial and final levels of the transition involved. Nevertheless, these two formulae must be used with caution if a result with accuracy better than a factor of two is required. The Lindholm (1941) formula gives a rather poor result and likewise the classical value of the damping constant gives results which are greatly overestimated at stellar densities ($N \approx 10^{14} \text{ cm}^{-3}$).

In Table II the damping constants of several Si II, III, IV, C II, IV and N II (I) lines are collected. The typical energy difference ΔE gives the order of the uncertainty of the results. If $kT/\Delta E > 10$, the uncertainty is less than 20%, if $kT/\Delta E < \frac{1}{2}$, it is of the order of 30% or 40%. The classical value γ_{cl}/N (evaluated for $N_e = 10^{14} \text{ cm}^{-3}$), the Lindholm value (1941) and the Brechot and Van Regemorter (1964) value (formula 1 of this paper) are given for comparison.

When $kT/\Delta E \ll 1$ the simple B-VR formula is rather useful; indeed, the elastic contribution is predominant and the adiabatic assumption is valid. Under this condition the very simple formula can give quite reliable results. Nevertheless, the semi-classical calculations are also very simple to perform, the Fortran IV program of the Observatoire de Meudon gives the result after a few seconds (IBM 360-65 computer), the input data are only the oscillator strengths and the energy differences of the interacting levels of the transitions involved.

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DISCUSSION

G. K. Oertel: Elastic collisions contribute strongly to the electron broadening of ionic lines. The impact theory with hyperbolic classical path is therefore applicable to only a small portion of the total damping constant in most cases — The good agreement with other calculations and with experiment is therefore very surprising.