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First-principles calculations of pressure broadening for the case of N_2 -perturbed 118 GHz fine-structure line in $O_2(X^3\Sigma_g^2)$

M. Gancewski¹, H. Jóźwiak¹, E. Quintas-Sánchez², R. Dawes², F. Thibault³, P. Wcisło¹

1. Institute of Physics, Faculty of Physics, Astronomy and Informatics, Nicolaus Copernicus University in Toruń, ul. Grudziądzka 5, 87-100 Toruń, Poland

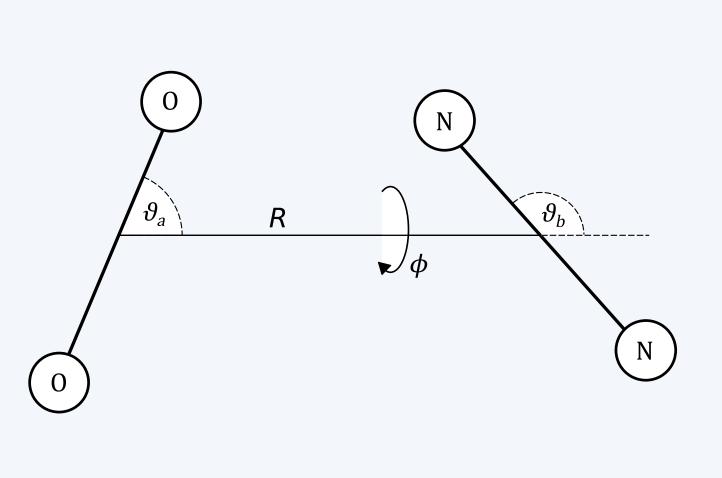
2. Department of Chemistry, Missouri University of Science and Technology, Rolla, Missouri 65409-0010, USA 3. Univ. Rennes, CNRS, IPR (Institut de Physique de Rennes)-UMR 6251, Rennes F-35000, France

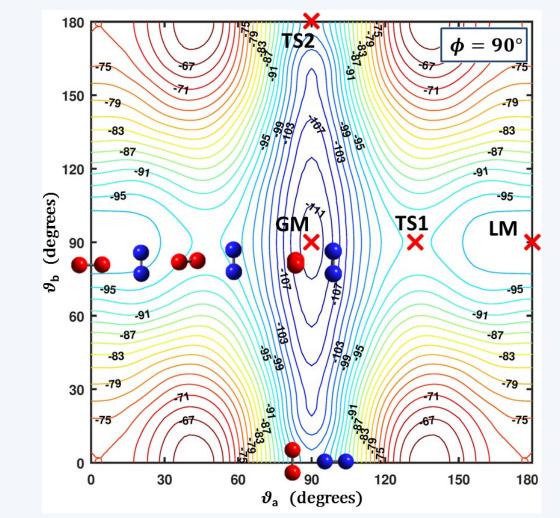
What is this about?

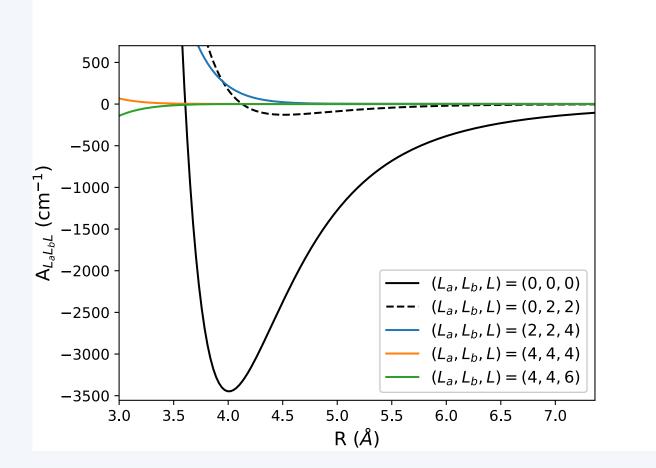
- The general problem: *How is the absorption* (emission) spectrum of O_2 modified when it is immersed in the large bath of N_2 molecules?
- This model question is important in general considerations of atmospheric physics (~78% of N₂ & ~21% of O₂ in our atmosphere...)
- We want to attack this problem *ab initio* style, using methods of theoretical physics and quantum chemistry
- Many spectral lines & bands of O_2 will be affected by its interactions with N_2 in a complicated manner
- A torough study of these is very challenging (preparation of the potential, solving the Schrödinger & kinetic equations...)
- We thus begin this program by adressing a simpler problem: <u>How much a single isolated line of O₂ is broadened by collisions with N₂ at different gas temperatures?</u>
- Here, we consider both (rigid) molecules to be in their ground states & we focus on the N_2 -induced perturbations of the isolated *fine-structure* line (centered at ~118 GHz) of $O_2(^3\Sigma)$

Our methodology

- In $O_2(X^3\Sigma)$ each rotational level splits into three *fine-structure* levels due to spin-spin interaction; we study the broadening of the $j=1 \rightarrow j=0$ *fine-structure* transition, occurring <u>within the first</u> rotational level
- By using angular momentum recoupling techniques [1, 2, 3] we separate the O_2 spin from the problem, and solve the Schrödinger equation for the collision of two spin-free particles on the 4-dimensional potential energy surface (PES)



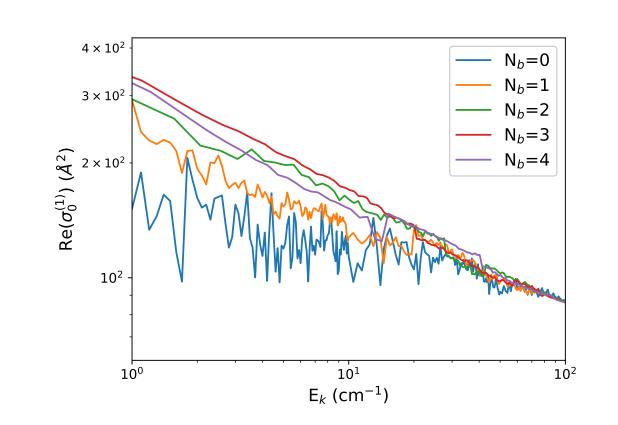


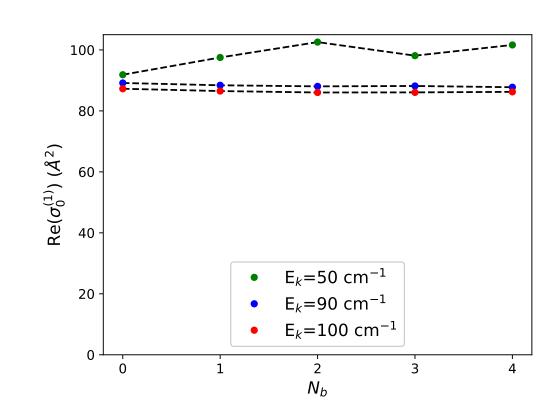


- Our new 4-dimensional O_2 - N_2 PES is computed using the (AE)UCCSD(T)-F12b/CBS methodology with aid of the "AUTOSURF" code [1, 4]; the Schrödinger equation is solved numerically using our "BIGOS" quantum scattering code [5]
- We transform the obtained <u>spin-free scattering matrices</u> back into the representation which <u>include the quantum spin number</u> we are thus in a position to <u>compute the line-broadening for a purely *fine-structure* transition in the single <u>rotational level</u></u>
- The "BIGOS" code uses the transformed S-matrices to calculate the pressure broadening cross-section for the N₂ perturbation of the $j=1 \rightarrow j=0$ fine-structure transition in O₂
- The energy-averaged cross sections yield the pressure broadening parameter for this line (*i.e.*, the half-width at half maximum, HWHM) as a function of the gas temperature [1]

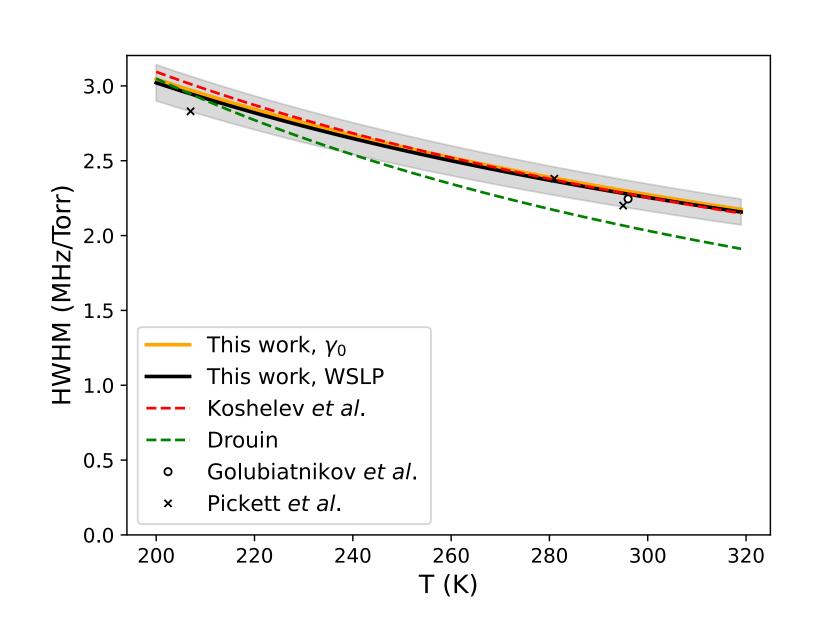
The results

- ullet We calculated the broadening cross sections in the kinetic energy range from 1 cm $^{-1}$ to \sim 100 cm $^{-1}$
- Here, cross sections are given for first 5 rotational levels of N_2 (labeled with N_b) as a function of kinetic energy; at high enough E_ν they become independent of the rotational state of the thermal bath (see the figure on the right)





- We extrapolate the cross sections given in the figure into the higher kinetic energies (up to ~4 000 cm⁻¹); we then assume that for N_2 rotational states with N_b >4, the cross sections are the same as for the N_b =4; thus in the evaluation of the statistical partition sums over bath variables we are able to include the N_2 levels up to N_b =31
- By suitable energy-averaging of the cross sections we obtain both the <u>speed-independent</u> & <u>speed-dependent</u> N_2 -broadening coefficients for the O_2 118 GHz line. The temperature dependence of the speed-dependent HWHM is obtained by averaging the speed-dependent Lorentz profiles over the Maxwellian distribution



$\gamma(T) = \gamma(T_0) \left(\frac{T_0}{T}\right)^{n_{\gamma}},$			
Experiment	T (K)	γ (MHz/Torr)	n_{γ}
Pickett et al. ¹⁹	207	2.83(6)	
	281	2.38(5)	
	295	2.2(1)	
Golubiatnikov <i>et al.</i> ²⁰	297	2.245(20)	
Drouin ²¹	296	2.06(6)	1.00(5)
Makarov <i>et al.</i> ¹⁵	300	2.255(14)	0.79(5)
Koshelev et al. 16	296	2.278(2)	0.781(6)
This work, WSLP	296	2.28(9)	0.7231(1)
This work, γ_0	296	2.30(9)	0.7245(6)

• Our *ab initio* calculations agree with the experimental data (see [1] for the relevant papers) within the estimated 4% uncertainty of numerical calculations (shaded area). This paves the way for more elaborate future analyses of the general line-shape problem for atmospheric O₂ perturbations

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