

## Construction of potential curves for diatomic molecular states by the IPA method

Asen Pashov, Włodzimierz Jastrzębski, Paweł Kowalczyk *Computer Physics Communications*, **128(3)**, 622-634 (2000)

DOI: [10.1016/S0010-4655\(00\)00010-2](https://doi.org/10.1016/S0010-4655(00)00010-2)

Software for calculation the transitions energies [ $\text{cm}^{-1}$ ] between state A and state B with example files:

{attachments}