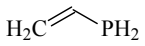
291 **C_2H_5P** **Vinylphosphine** **C_s (*syn*)**
MW Ethenylphosphine



An *ab initio* calculation at the MP2/6-311G (2d,p) level reproduces the experimental rotational constants satisfactorily.

Dréan, P., Colmont, J.-M., Lesarri, A., López, J.C.: J. Mol. Spectrosc. **176** (1996) 180.
II/25B(**3**, 843)