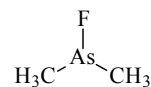


294 **C₂H₆AsF**ED, *ab initio*
calculations**Fluorodimethylarsine**

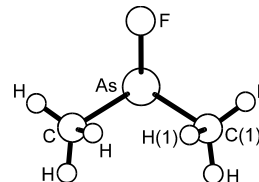
Dimethylarsinous fluoride

C_s

| r_g | Å ^{a)} |
|-------|-----------------|
| As–C | 1.955(1) |
| As–F | 1.758(1) |

| r_a^0 | Å ^{a)} |
|---------|-----------------|
| As–C | 1.951(1) |
| As–F | 1.754(1) |
| C–H | 1.061(3) |

| θ_a^0 | deg ^{a)} |
|--------------|-------------------|
| C–As–F | 95.3(5) |
| C–As–C | 96.9(8) |
| As–C–H | 106.7(7) |
| τ^b | 57(11) |



The harmonic force constants were calculated at the MP2/6-31G** level of theory.
The nozzle was at about 293 K.

^{a)} Estimated standard errors.

^{b)} Torsional angle F–As–C(1)–H(1), $\tau = 0^\circ$ for *syn* position; positive value for the anticlockwise rotation.

Downs, A.J., Greene, T.M., McGrady, G.S., Townson, N., Brain, P.T., Pulham, C.R., Rankin, D.W.H., Robertson, H.E., Smart, B.A., Alberts, I.L.: Inorg. Chem. **35** (1996) 6952.