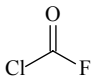
18 CClFO Carbonyl chloride fluoride **C_s**

IR, *ab initio* calculations
(CCSD(T)/TZ2Pf)



r_e	Å [a])	θ_e	deg [a])
C=O	1.173(1)	O=C–F	124.0(1)
C–F	1.323(1)	O=C–Cl	126.4(1)
C–Cl	1.721(1)		

[a]) Uncertainties were unidentified, possibly estimated standard errors.

Demaison, J., Perrin, A., Bürger, H.: J. Mol. Spectrosc. **221** (2003) 47.

II/25B(**3**, 32)