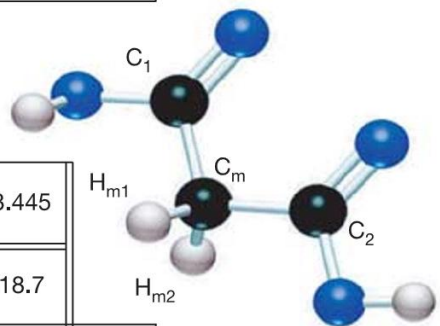
(a)

	C_1	C_2	C_m	H_{m1}
C_1	5.893			
C_2	0.227	1.057		
C_m	0.935	1.070	−3.445	
H_{m1}	−1.5	1.4	−18.7	
H_{m2}	2.0	1.0	−0.9	−22.0



(b)