ADDITIONS AND CORRECTIONS

2009, Volume 113A

Solvejg Jørgensen* and Allan Gross: Theoretical Investigation of the Reaction between Carbonyl Oxides and Ammonia

Page 10284. Figures 2−5 should be the following figures.

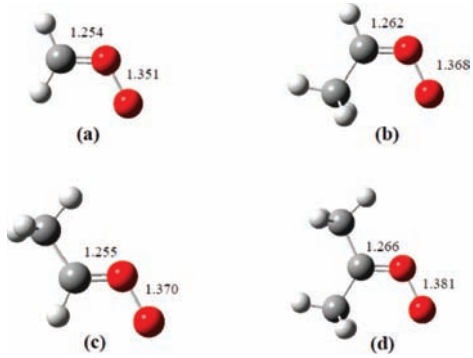

Figure 2. Structures of the carbonyl oxides studied in this work optimized at B3LYP/6-311++G(2d,2p) (white = hydrogen, gray = carbon, and red = oxygen atoms): (a) H_2COO, (b) *syn*-$(CH_3)HCOO$, (c) *anti*-$(CH_3)HCOO$, and (d) $(CH_3)_2COO$.

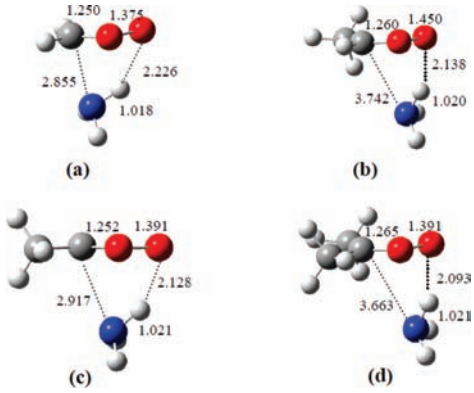

Figure 3. Structures of the prereactive complex studied in this work optimized at B3LYP/6-311++G(2d,2p) (blue = nitrogen atoms): (a) $H_2COO\cdots H-NH_2$, (b) *syn*-$(CH_3)HCOO\cdots H-NH_2$, (c) *anti*-$(CH_3)_2COO\cdots H-NH_2$, and (d) $(CH_3)_2COO\cdots H-NH_2$. Distances are given in angstroms.

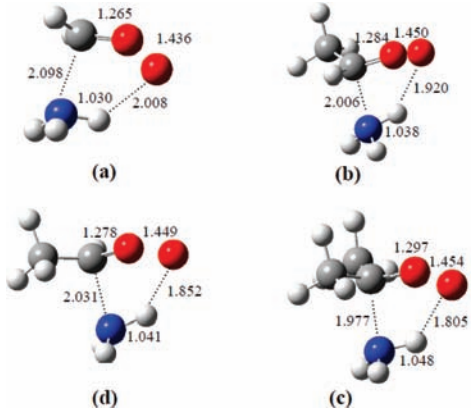

Figure 4. Structures of the transition-state complex studied in this work optimized at B3LYP/6-311++G(2d,2p): (a) $[H_2COO\cdots H\cdots NH_2]^{\#}$, (b) [*syn*-$(CH_3)HCOO\cdots H\cdots NH_2]^{\#}$, (c) [*anti*-$(CH_3)HCOO\cdots H\cdots NH_2]^{\#}$, and (d) $[(CH_3)_2COO\cdots H\cdots NH_2]^{\#}$. Distances are given in angstroms.

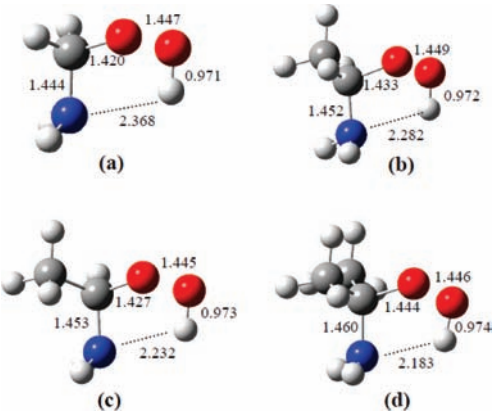

Figure 5. Structures of the products studied in this work optimized at B3LYP/6-311++G(2d,2p): (a) $H_2C(OOH)(NH_2)$, (b) *syn*-$(CH_3)HC(OOH)(NH_2)$, (b) *anti*-$(CH_3)HC(OOH)(NH_2)$, and (d) $(CH_3)_2C(OOH)(NH_2)$. Distances are given in angstroms.

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