

| Reactant | Reaction class | Product | E_b kcal mol ⁻¹ | E_{react} kcal mol ⁻¹ | k (300 K) s ⁻¹ |
|---------------------|------------------------------|---|---------------------------------|--|-------------------------------------|
| HMVKAO ₂ | -OH 1,5-H-shift | CH ₃ -C(=O)-CH(O [•])-CH ₂ OOH | 21.6 | 20.4 | 5.0 × 10 ⁻⁴ |
| | α-OH 1,4-H-shift | CH ₃ -C(=O)-C [•] (OH)-CH ₂ OOH | 24.7 | -6.2 | 3.3 × 10 ⁻³ |
| | -CH ₃ 1,6-H-shift | C [•] H ₂ -C(=O)-CH(OH)-CH ₂ OOH | 23.1 | 10.2 | 5.9 × 10 ⁻⁴ |
| HMVKBO ₂ | -OH 1,5-H-shift | CH ₃ -C(=O)-CH(OOH)-CH ₂ O [•] | 22.5 | 20.6 | 8.8 × 10 ⁻⁵ ^a |
| | α-OH 1,4-H-shift | CH ₃ -C(=O)-CH(OOH)-C [•] HOH | 25.1 | 6.5 | 3.2 × 10 ⁻⁵ |
| | -CH ₃ 1,6-H-shift | C [•] H ₂ -C(=O)-CH(OOH)-CH ₂ OH | 27.4 | 10.0 | 3.8 × 10 ⁻⁵ |
| | HO ₂ elimination | CH ₃ -C(=O)-CH=CHOH + HO ₂ | 30.0 | -1.5 | 6.1 × 10 ⁻¹⁰ |