

133.4, 143.8, 145.1, 193.7 (C=O). Anal. Calcd for C₁₀H₁₂OS: C, 66.63; H, 6.71; O, 8.88. Found: C, 66.83; H, 6.92; O, 9.18.

1-(2-Furyl)-2,3-dimethylbut-3-en-1-one (3d):¹⁸ colorless oil, bp 70 °C (5.0 mmHg, Kugelrohr). IR (neat): 1673.0 cm⁻¹ ($\nu_{C=O}$). ¹H NMR (CDCl₃, 400 MHz): δ 1.31 (d, 3H, CH₃, J = 6.84 Hz), 1.76 (s, 3H, CH₃), 3.94 (q, CH, J = 6.84 Hz), 4.90 (s, 1H, CH₂=), 4.92 (s, 1H, CH₂=), 6.52 (dd, 1H, furyl, J = 3.42, 1.47 Hz), 7.23 (d, 1H, furyl, J = 3.42 Hz), 7.58 (d, 1H, furyl, J = 1.47 Hz). ¹³C NMR (CDCl₃, 100 MHz): δ 15.3, 20.5, 49.2, 112.1, 113.3, 117.5, 144.6, 146.2, 152.3, 189.8 (C=O). Anal. Calcd for C₁₀H₁₂O₂: C, 73.15; H, 7.37. Found: C, 72.85; H, 7.64.

(E)-1-(2-Thienyl)-2-methylpent-3-en-1-one (3e): pale yellow oil, bp 80 °C (5.0 mmHg, Kugelrohr). IR (neat): 1659.8 cm⁻¹ ($\nu_{C=O}$). ¹H NMR (CDCl₃, 400 MHz): δ 1.31 (d, 3H, CH₃, J = 6.84 Hz), 1.67 (d, 3H, CH₃, J = 5.37 Hz), 3.92 (dq, 1H, CH, J = 7.32, 6.84 Hz), 5.59 (dd, 1H, CH=, J = 16.85, 7.32 Hz), 5.65 (dq, 1H, CH=, J = 16.85, 5.37 Hz), 7.11–7.13 (m, 1H, thienyl), 7.61–7.62 (m, 1H, thienyl), 7.73–7.76 (m, 1H, thienyl). ¹³C NMR (CDCl₃, 100 MHz): δ 13.1, 17.5, 46.4, 125.5, 128.0, 131.8, 132.0, 133.5, 143.7, 194.5 (C=O). Exact mass: calcd for C₁₀H₁₂OS, 180.0609; found, 180.0608. Anal. Calcd

for C₁₀H₁₂OS: C, 66.63; H, 6.71. Found (for a 56:44 mixture of **3e** and **3e'**): C, 66.69; H, 6.75.

(Z)-1-(2-Thienyl)-2-methylpent-3-en-1-one (3e'): pale yellow oil, bp 80 °C (5.0 mmHg, Kugelrohr). IR (neat): 1659.8 cm⁻¹ ($\nu_{C=O}$). ¹H NMR (CDCl₃, 400 MHz): δ 1.29 (d, 3H, CH₃, J = 6.84 Hz), 1.77 (d, 3H, CH₃, J = 5.37 Hz), 4.14 (dq, 1H, CH, J = 8.79, 6.84 Hz), 5.56 (dd, 1H, CH=, J = 9.28, 8.79 Hz), 5.57 (dq, 1H, CH=, J = 9.28, 5.37), 7.11–7.13 (m, 1H, thienyl), 7.61–7.62 (m, 1H, thienyl), 7.73–7.76 (m, 1H, thienyl). ¹³C NMR (CDCl₃, 100 MHz): δ 11.7, 17.9, 41.6, 117.6, 127.5, 130.5, 130.8, 133.5, 143.6, 194.7 (C=O).

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Additions and Corrections

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Judith C. Gallucci, Olivier Gobley, Florence Zaegel, Philippe Meunier, Bernard Gautheron, Holger Lange, Rolf Gleiter, Natasha Kozmina, and Leo A. Paquette*: Solid-State Structural Analysis of the "Naked" Isodicyclopentadienide Anion.

Page 112. The space groups for the Na(isodiCp)(15-crown-5) and the [K-cryptand(2.2.2)]⁺(isodiCp)⁻ structures were incorrectly reported in refs 13 and 14 as *P1*. Both space groups are *P1̄*.

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