Campestarenes: new building blocks with 5-fold symmetry

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Fig 4 MALDI-TOF MS of 1i



Fig 5 MALDI-TOF MS of **1**j





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Fig 8 MALDI-TOF MS of **1m**



Fig 9 MALDI-TOF MS of **1n**







Fig 12 $^{\rm 13}C$ NMR spectrum of ${\bf 6g\text{-}Et}$ in CDCl_3



Fig 14 $^{\rm 13}C$ NMR spectrum of ${\bf 6g-tBu}$ in ${\rm CDCI_3}$





Fig 18 $^{\rm 13}\text{C}$ NMR spectrum of 2g-Et in CDCl_3















Fig 27 $^{\rm 11}\text{B}$ NMR spectrum of 10n in CD_3OD





Fig 29 ¹H NMR spectrum of 10o in CD₃OD



Fig 30 ¹¹B NMR spectrum of **100** in CD₃OD





















Fig 40 COSY spectrum of **1f** in DMSO- d_6 . Red denotes correlation between CH₂ and CH₃ on the ethyl group.



Fig 42 ¹H NMR spectrum of **1h** in DMSO- d_6



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Fig 46 ¹H NMR spectra in DMSO- d_6 of a) **1i**, b) de-methylated **1i** using 5 equivalents of BBr₃ to **1i** and c) using 15 equivalents of BBr₃ to **1i**. x denotes methanol.



Fig 47 UV-Vis spectrum of **1f** (8.5 x 10^{-5} mol.L⁻¹ in DMSO)



Fig 48 UV-Vis spectrum of 1g (10 x 10⁻⁵ mol.L⁻¹ in DMSO)



Fig 49 UV-Vis spectrum of **1h** (9.7 x 10^{-5} mol.L⁻¹ in DMSO)



Fig 50 UV-Vis spectrum of **1i** (13 x 10^{-5} mol.L⁻¹ in DMSO)



Fig 51 UV-Vis spectrum of 1j (5.1 x 10⁻⁵ mol.L⁻¹ in DMSO)



Fig 52 UV-Vis spectrum of **1k** (approx. $5.1 \times 10^{-5} \text{ mol.L}^{-1}$ in DMSO)



Fig 53 UV-Vis spectrum of **1I** (approx. 9.1×10^{-5} mol.L⁻¹ in DMSO)



Fig 54 UV-Vis spectrum of 1m (approx. 6.7 x 10⁻⁵ mol.L⁻¹ in DMSO)



Fig 55 UV-Vis spectrum of $\mathbf{1n}$ (6.8 x 10⁻⁵ mol.L⁻¹ in DMSO)



Fig 56 UV-Vis spectrum of **1o** (3.6 x 10^{-5} mol.L⁻¹ in DMSO)







Fig 58 IR spectrum of **1g**







Fig 60 IR spectrum of 1i







Fig 62 IR spectrum of **1k**











Fig 66 IR spectrum of **10**