

Supporting Information

2-Aminophenolate Ligands for Phosphorus(V): A Lithium Salt Featuring the Chiral $[P(OC_6H_4NR)_3]^-$ Anion

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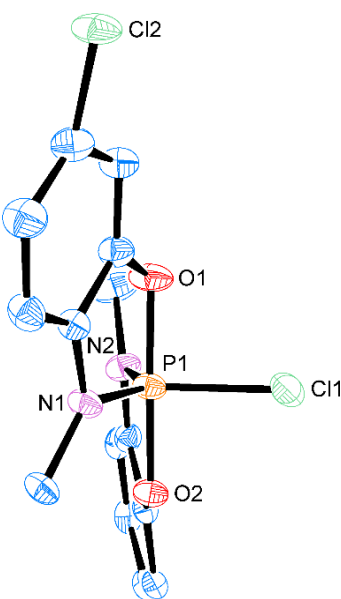


Fig. S1 The minor component (7%) in the molecular structure for intermediate **3a** (thermal ellipsoids are displayed at 50% probability level). Hydrogen atoms are omitted for clarity. Bond lengths and angles are similar to those found in the major structure.

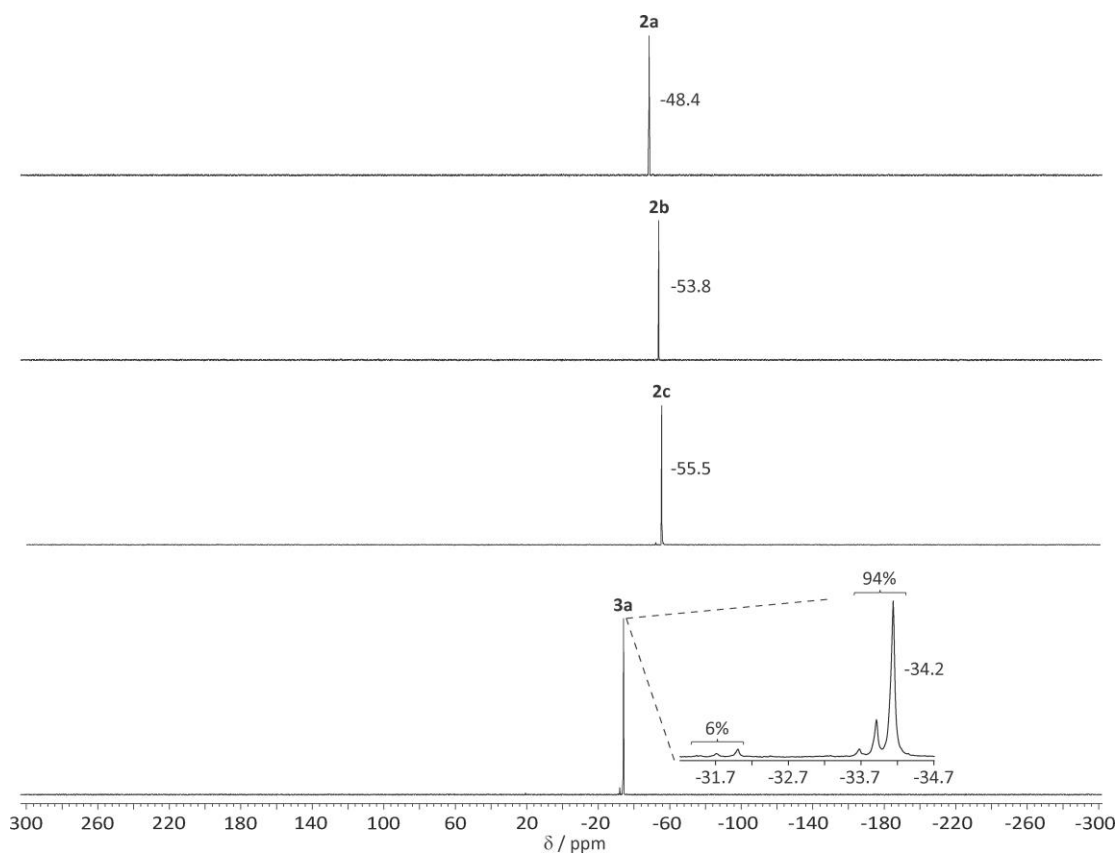


Fig. S2 $^{31}\text{P}\{^1\text{H}\}$ NMR spectra (121 MHz, CDCl_3 , 298 K) of phosphoranes **2a-c** and intermediate **3a**.

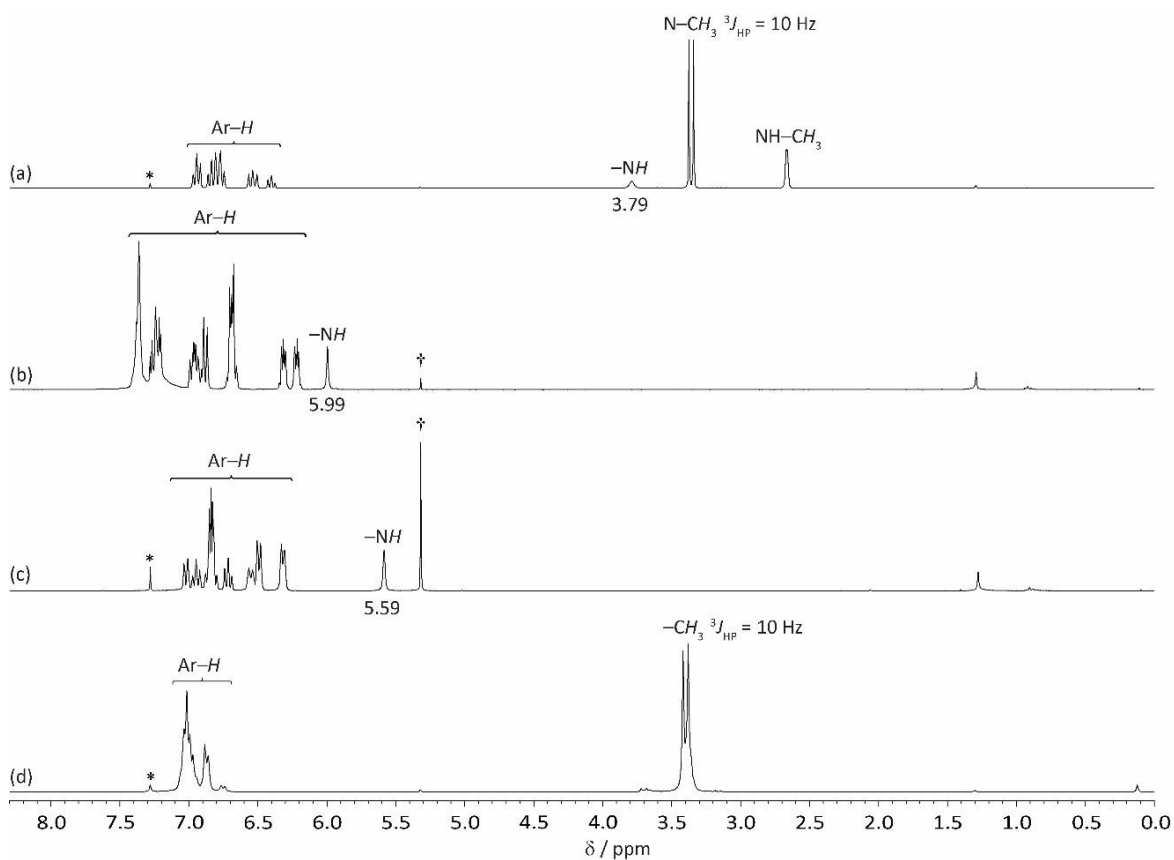


Fig. S3 ^1H NMR spectra (300 MHz, CDCl_3 , 298 K) of phosphoranes: (a) **2a**; (b) **2b**; (c) **2c**; (d) **3a** recrystallized from CH_2Cl_2 (* indicates the residual CHCl_3 and † indicates the residual CH_2Cl_2 solvent).

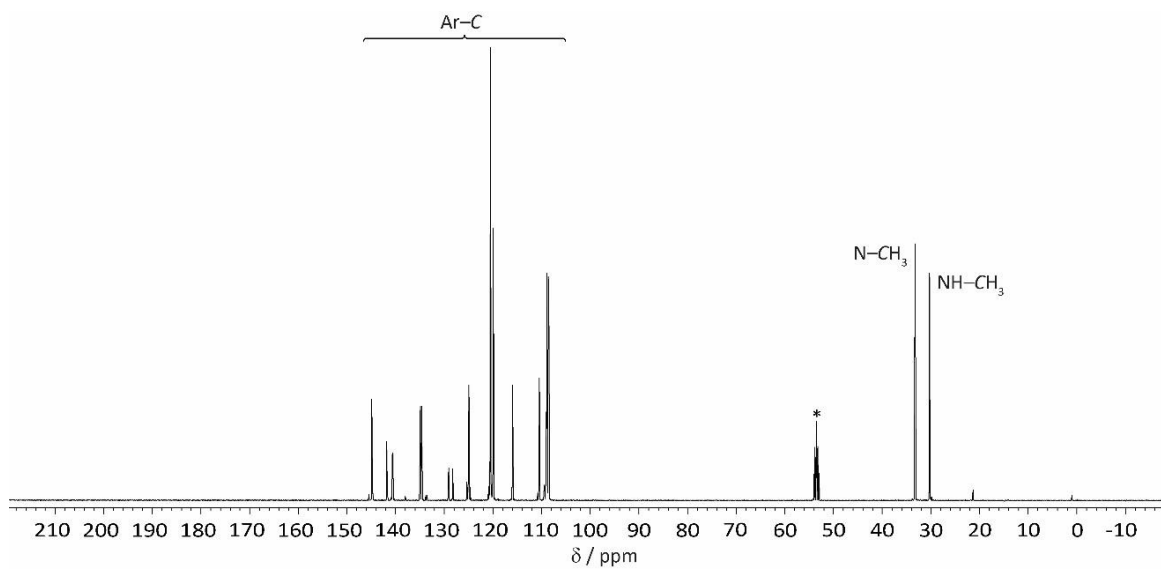


Fig. S4 $^{13}\text{C}\{^1\text{H}\}$ NMR spectrum (100.5 MHz, CD_2Cl_2 , 298 K) of phosphorane **2a** (* indicates the solvent).

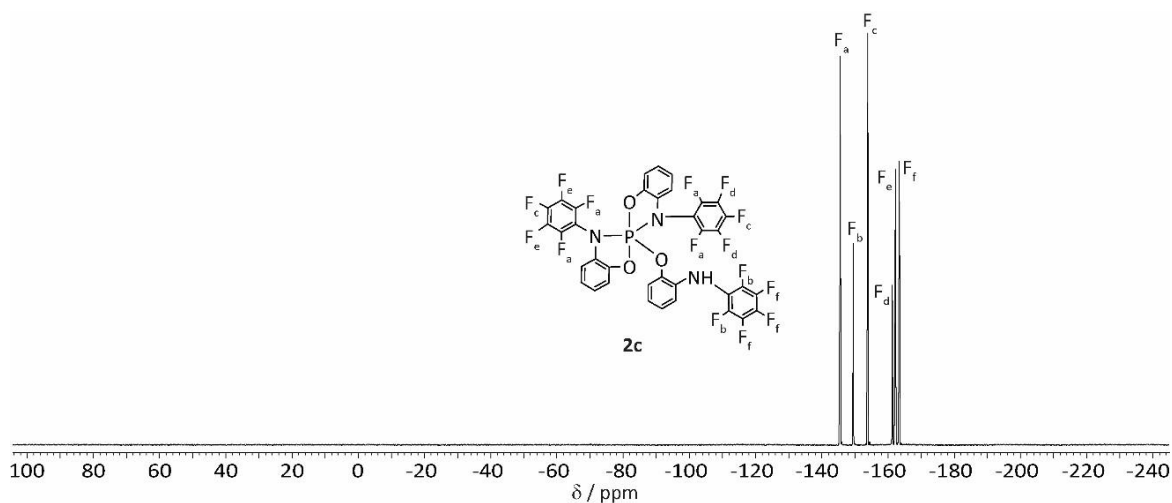


Fig. S5 $^{19}\text{F}\{^1\text{H}\}$ NMR spectrum (282 MHz, CDCl_3 , 298 K) of phosphorane **2c**.

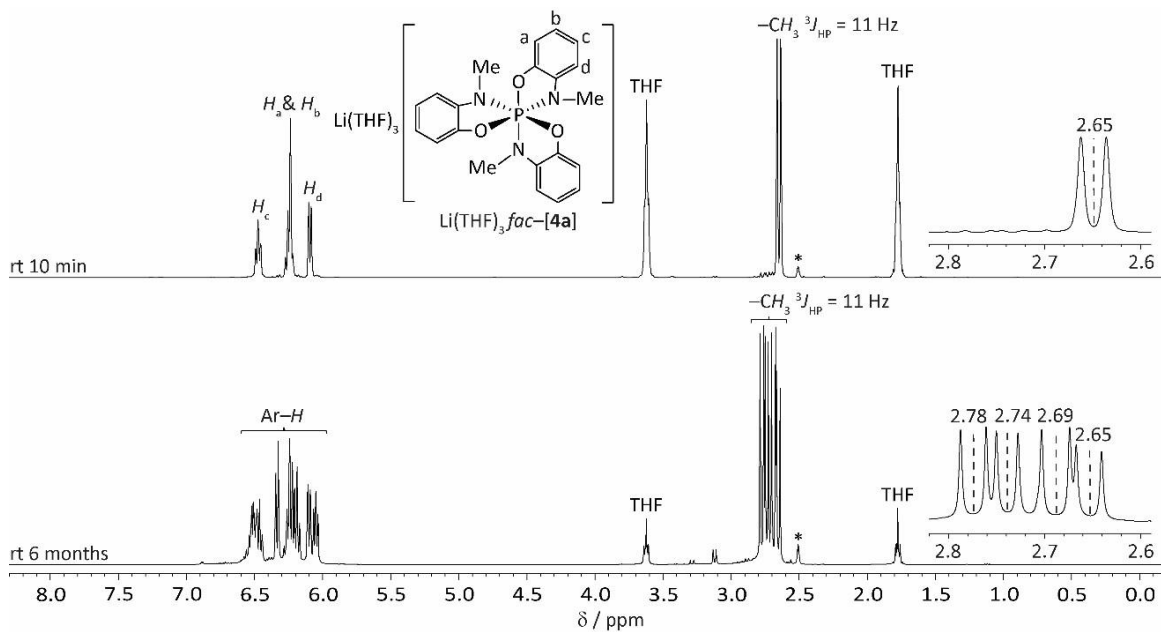


Fig. S6 ^1H NMR spectra (400 MHz, $\text{DMSO}-d_6$, 298 K) of $\text{Li}(\text{THF})_3 \text{ fac-[4a]}$ (* indicates the solvent).

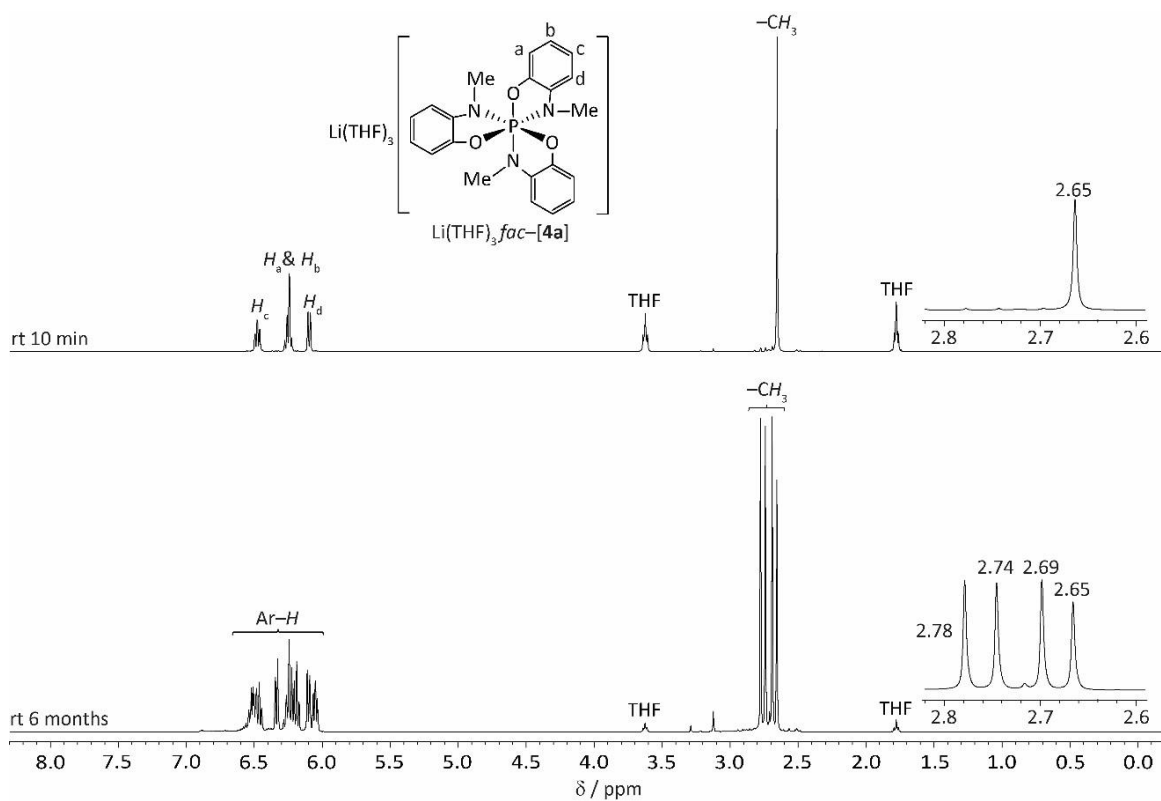


Fig. S7 $^1\text{H}\{^{31}\text{P}\}$ NMR spectra (400 MHz, $\text{DMSO-}d_6$, 298 K) of $\text{Li(THF)}_3 \text{ fac-[4a]}$.

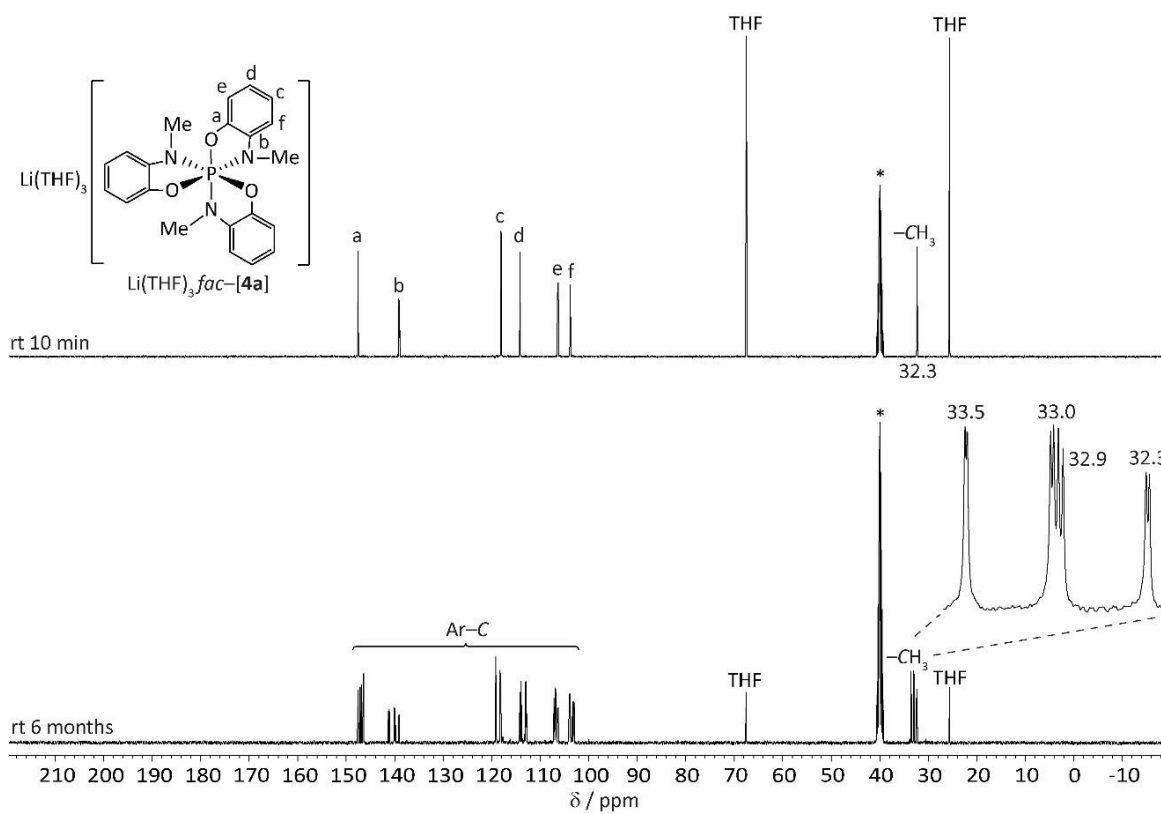


Fig. S8 $^{13}\text{C}\{^1\text{H}\}$ NMR spectra (100.5 MHz, $\text{DMSO-}d_6$, 298 K) of $\text{Li}(\text{THF})_3 \text{ fac-[4a]}$ (* indicates the solvent).

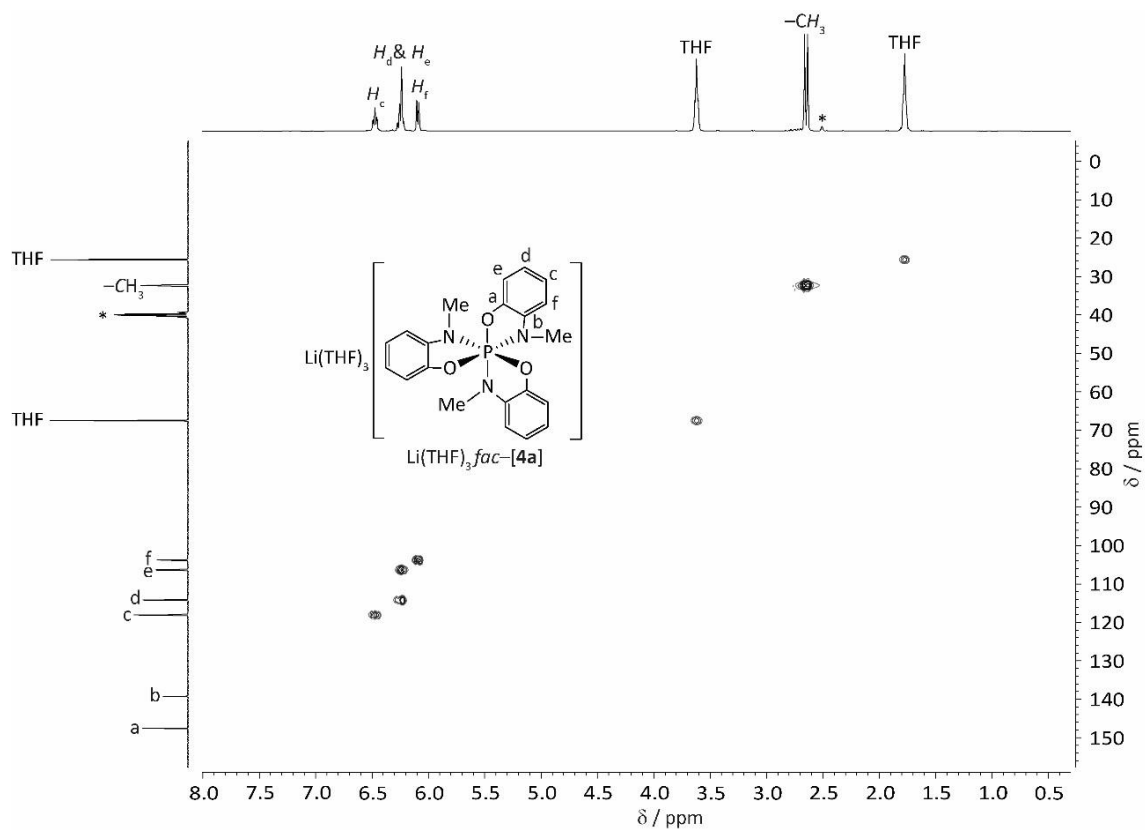


Fig. S9 ^1H - ^{13}C HSQC NMR spectrum (400 MHz for ^1H , $\text{DMSO}-d_6$, 298 K) of $\text{Li(THF)}_3\text{fac-[4a]}$ (The ordinate axis shows the $^{13}\text{C}\{^1\text{H}\}$ NMR spectrum and the abscissa axis shows the ^1H NMR spectrum; * indicates the solvent).

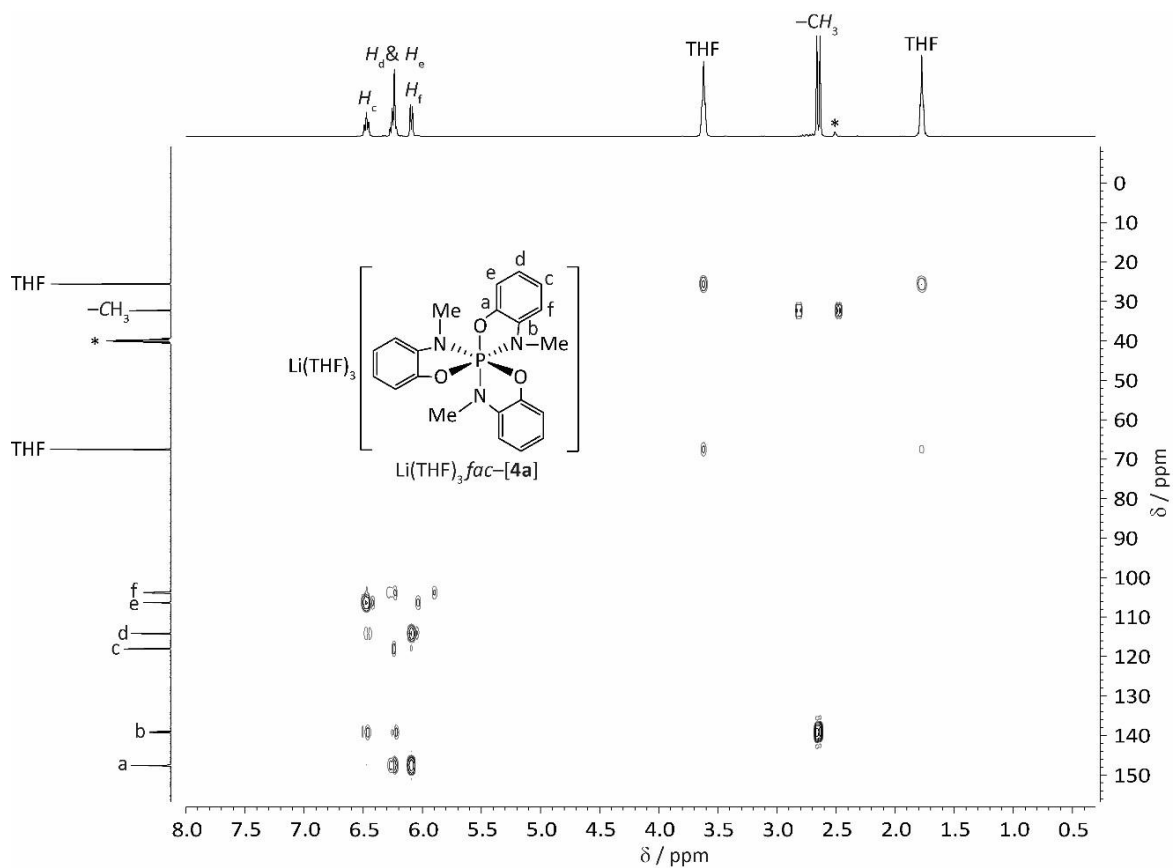


Fig. S10 ^1H - ^{13}C HMBC NMR spectrum (400 MHz for ^1H , $\text{DMSO-}d_6$, 298 K) of $\text{Li}(\text{THF})_3 \text{ fac-[4a]}$ (The ordinate axis shows the $^{13}\text{C}\{^1\text{H}\}$ NMR spectrum and the abscissa axis shows the ^1H NMR spectrum; * indicates the solvent).

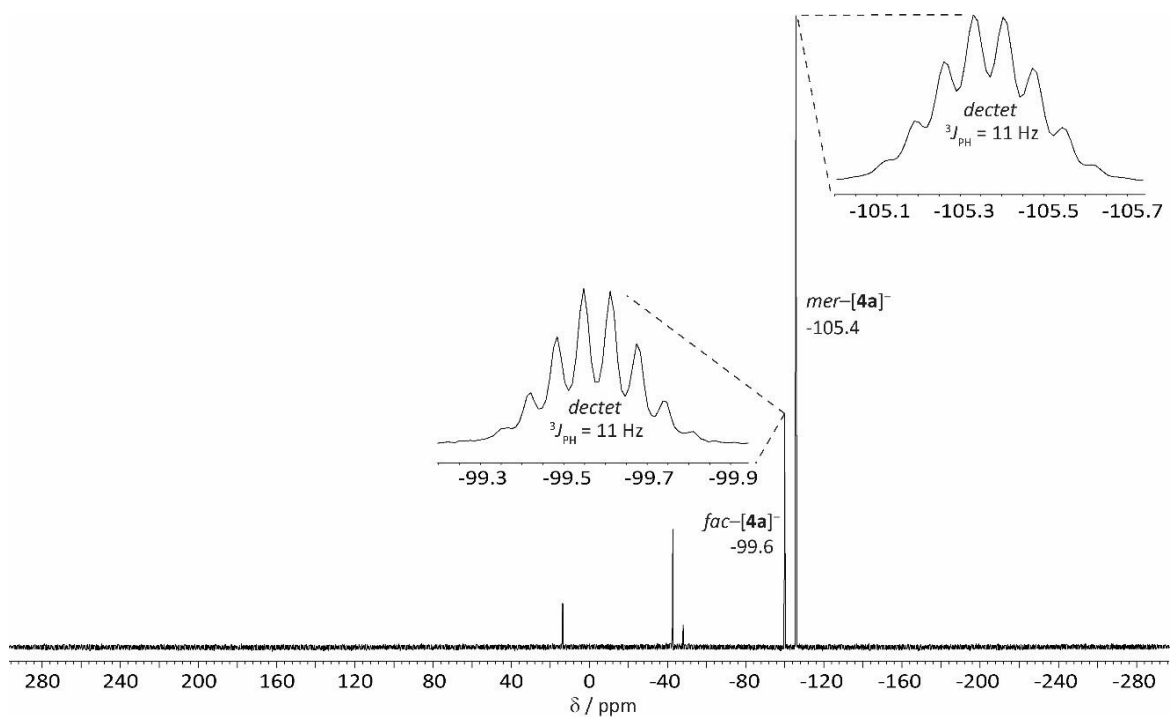


Fig. S11 ^{31}P NMR spectrum (121 MHz, 298 K) of $\text{Li}(\text{THF})_3 \text{ fac-[4a]}$ in $\text{DMSO-}d_6$ over 6 months.

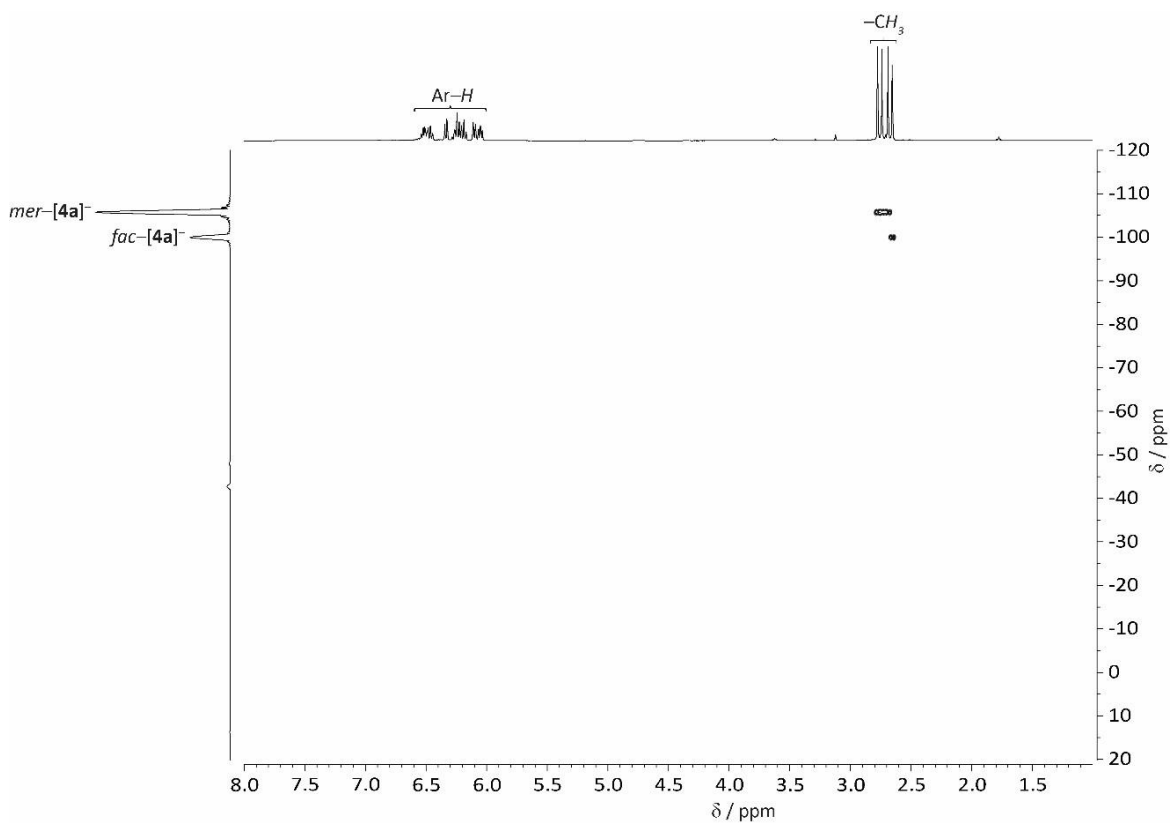


Fig. S12 ^1H - ^{31}P HMBC NMR spectrum (400 MHz for ^1H , $\text{DMSO-}d_6$, 298 K) of $\text{Li}(\text{THF})_3 \text{ fac-[4a]}$ (The ordinate axis shows the $^{31}\text{P}\{^1\text{H}\}$ NMR spectrum and the abscissa axis shows the ^1H NMR spectrum).

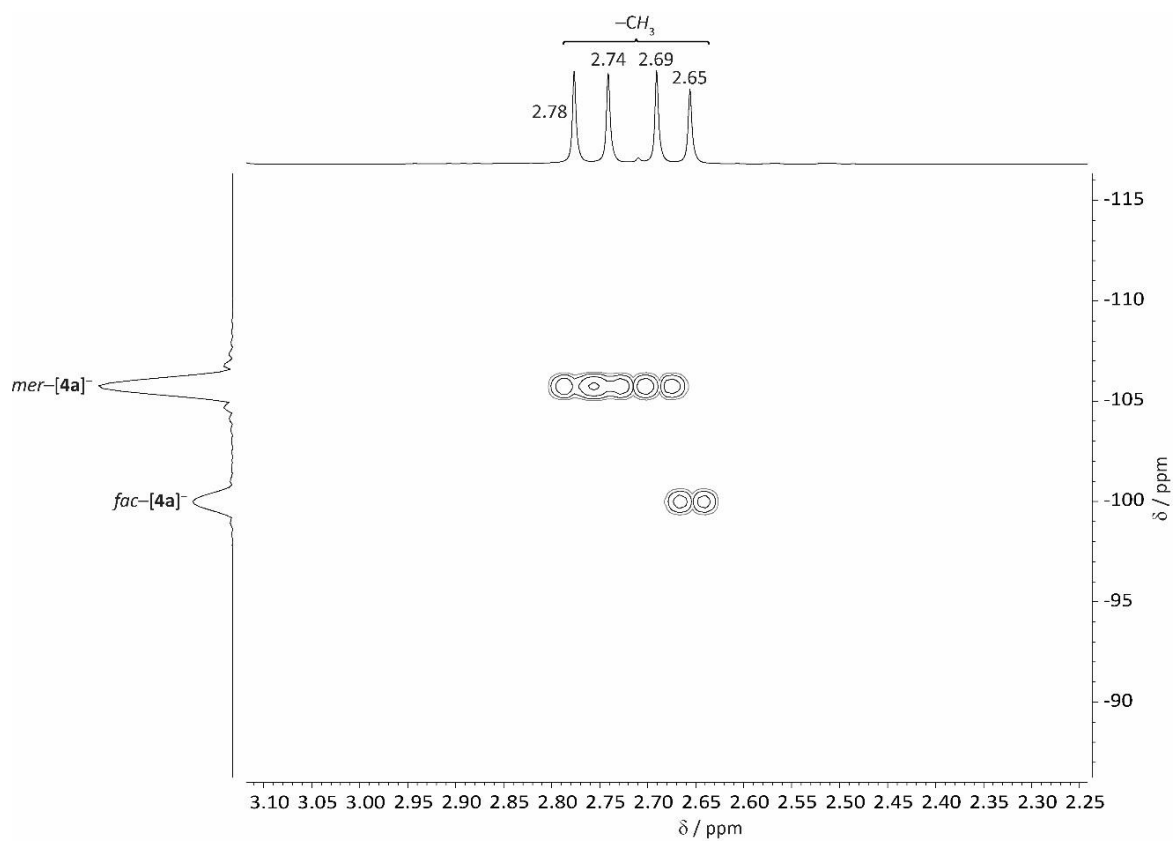


Fig. S13 Partial ^1H - ^{31}P HMBC NMR spectrum (400 MHz for ^1H , $\text{DMSO}-d_6$, 298 K) of $\text{Li}(\text{THF})_3 \text{fac-[4a]}$ (The ordinate axis shows the $^{31}\text{P}\{^1\text{H}\}$ NMR spectrum and the abscissa axis shows the ^1H NMR spectrum).

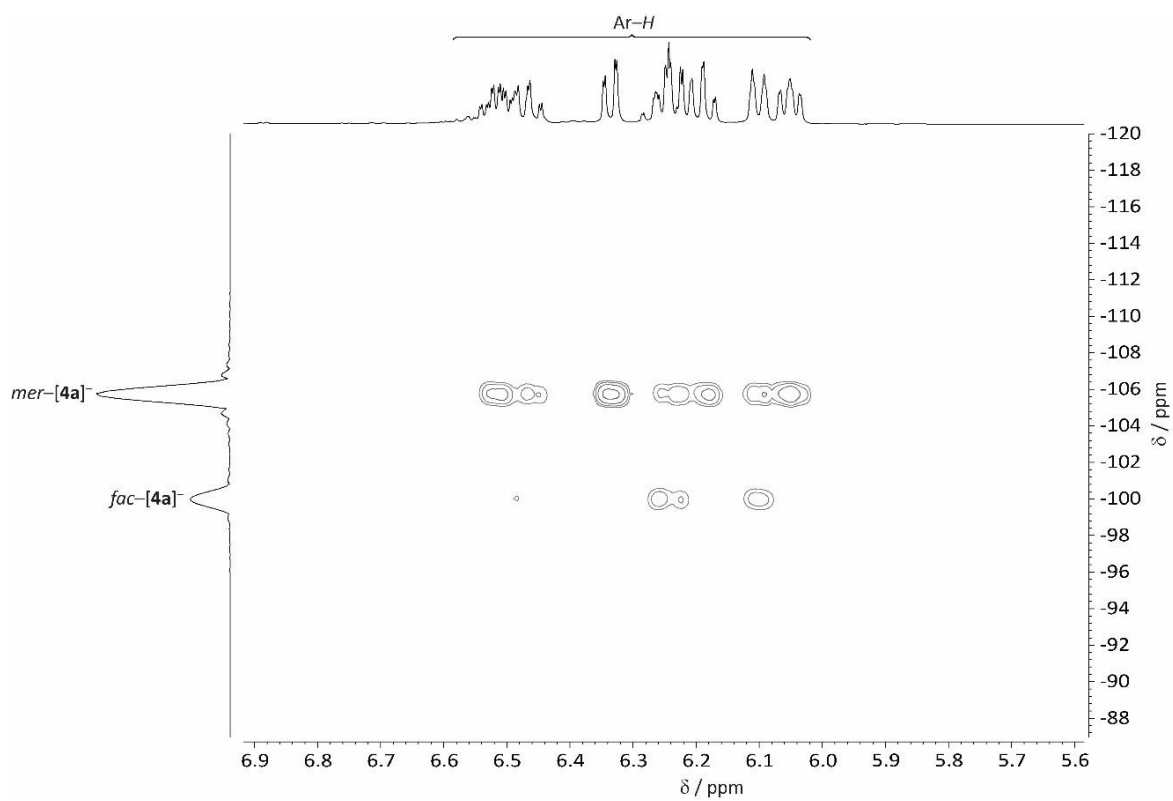


Fig. S14 Partial ^1H - ^{31}P HMBC NMR spectrum (400 MHz for ^1H , $\text{DMSO}-d_6$, 298 K) of $\text{Li}(\text{THF})_3 \text{fac-[4a]}$ (The ordinate axis shows the $^{31}\text{P}\{^1\text{H}\}$ NMR spectrum and the abscissa axis shows the ^1H NMR spectrum).