

Phosphorus-31 to Lithium-7 Nuclear Spin Coupling in Lithiated Organophosphorus Compounds

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^{31}P and ^7Li n.m.r. spectra with $J(^{31}\text{P}-^7\text{Li}) = ca. 50 \text{ Hz}$ show $\text{LiCH}(\text{PPh}_2)_2$ and $\text{LiCH}[\text{PPh}_2][\text{P}(\text{S})\text{PPh}_2]$ to be covalent monomeric, and LiPPh_2 to be covalent dimeric in Et_2O solution at low temperatures.

The reaction between alkyl-lithiums and a range of organophosphorus compounds with an active hydrogen can yield synthetically important reagents whose structures in solution are not yet established, but which may in principle be ionic, or covalent with the lithium attached either to carbon or to phosphorus.¹⁻³ Furthermore, since most known alkyl-lithiums exist as polymers in solution,⁴ this possibility must also be considered for the lithiated organophosphorus species. We now report ^7Li ($I = \frac{3}{2}$, abundance 92.6%) and ^{31}P n.m.r. spectra, together with $^{31}\text{P}\{^7\text{Li}, ^1\text{H}\}$ decoupling experiments from several of these species in ethereal solution at temperatures below 240 K, which show the first reported examples of well resolved $^{31}\text{P}-^7\text{Li}$ spin coupling, and demonstrate the presence of covalent monomers and dimers.

The ^7Li spectrum [JEOL FX90Q, Figure 1(a)] of a 0.5 M solution of Ph_2PLi (from Ph_2PH and BuLi) in Et_2O at 200 K is a well resolved 1:2:1 triplet [$\delta(^7\text{Li}) 0.5 \text{ p.p.m. rel. to molar LiBr in D}_2\text{O}$] showing each lithium to be coupled to two phosphorus nuclei with $J(^{31}\text{P}-^7\text{Li}) 45 \text{ Hz}$; the proton-decoupled ^{31}P spectrum [Figure 1(b)] is a 1:2:3:4:3:2:1 septet [$\delta(^{31}\text{P}) -36 \text{ p.p.m.}$, $J(^{31}\text{P}-^7\text{Li}) 45 \text{ Hz}$] indicating that each phosphorus is associated symmetrically with two lithium

atoms. The major species present is thus covalent, and in view of earlier molecular weight measurements⁵ can be assigned the dimeric structure (1) in which the two Ph_2P units are joined by a pair of bridging lithium atoms. (On a statistical basis the abundance of dimers containing only one ^7Li is <14%.) The ^7Li ($\delta 0.3 \text{ p.p.m.}$) spectrum of $(\text{Ph}_2\text{P})_2\text{CHLi}$ [from $(\text{Ph}_2\text{P})_2\text{CH}_2$ and BuLi] under the same conditions showed that each lithium is coupled to two phosphorus atoms, and the ^{31}P spectrum was a 1:1:1:1 quartet [$\delta(^{31}\text{P}) -3.7 \text{ p.p.m.}$] with $J(^{31}\text{P}-^7\text{Li}) 46 \text{ Hz}$. This species is therefore a covalent monomer in Et_2O . Each lithium ($\delta 0.5 \text{ p.p.m.}$) in $[\text{Ph}_2\text{P}(\text{S})][\text{Ph}_2\text{P}]\text{CHLi}$ {from $[\text{Ph}_2\text{P}(\text{S})][\text{Ph}_2\text{P}]\text{CH}_2$ and BuLi } was similarly found to be coupled to only one phosphorus, and the ^{31}P spectrum of this species is shown in Figure 2(a). The $^{31}\text{P}\{^7\text{Li}, ^1\text{H}\}$ decoupling experiment [Figure 2(b)] confirms the assignments: $\text{Ph}_2\text{P}(\text{S})$, $\delta(^{31}\text{P}) +45.6 \text{ p.p.m.}$, $J(^{31}\text{P}-^7\text{Li}) <4 \text{ Hz}$; Ph_2P , $\delta(^{31}\text{P}) -17.4 \text{ p.p.m.}$, $J(^{31}\text{P}-^7\text{Li}) 54 \text{ Hz}$; $^2J(^{31}\text{P}-^{31}\text{P}) 166 \text{ Hz}$. This species is therefore also a covalent monomer in Et_2O .

These results do not establish unequivocally whether the lithium is attached to carbon or to phosphorus in the last two compounds, but for $(\text{Ph}_2\text{P})_2\text{CHLi}$ we tentatively favour the structure (2a) rather than (2b), and for $[\text{Ph}_2\text{P}(\text{S})][\text{Ph}_2\text{P}]\text{CHLi}$ the structure (3a) rather than (3b) because (i) in each the magnitude of $J(^{31}\text{P}-^7\text{Li})$ is very close to that in (1) in which P-Li bonds must be present; (ii) (2b) and (3b) are substituted alkyl-lithiums which can be expected to exist as

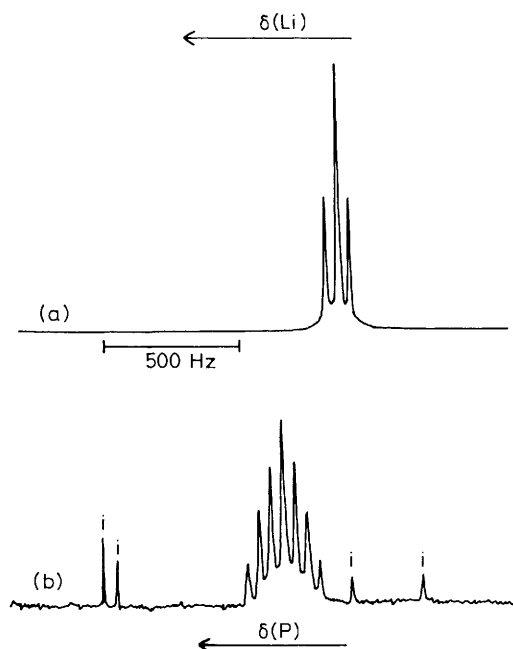


Figure 1. N.m.r. spectra of Ph_2PLi in Et_2O at 200 K. (a) ^7Li at 34.8 MHz. (b) ^{31}P at 36.2 MHz. Peaks marked i are impurities.

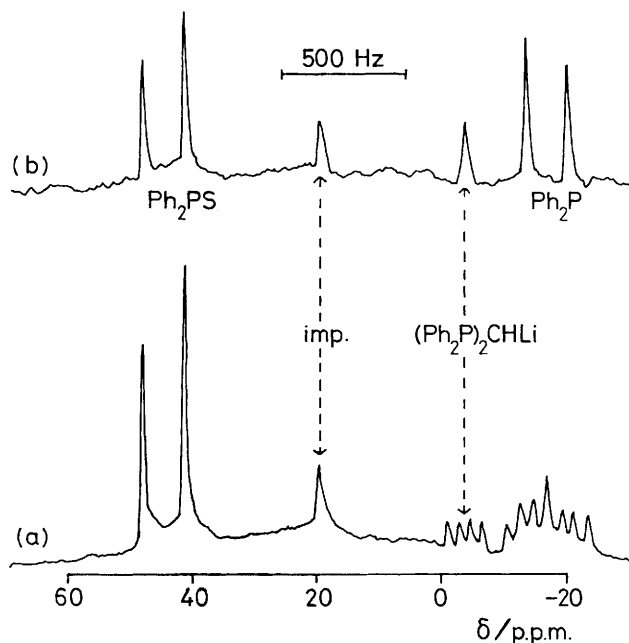
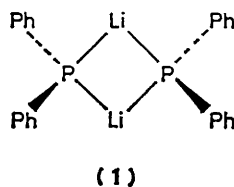
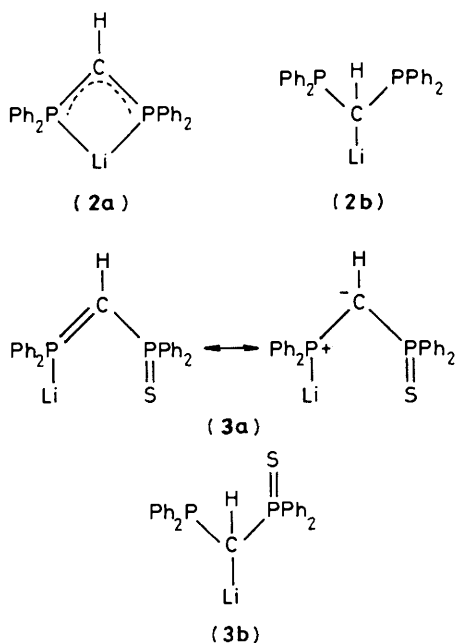


Figure 2. Proton-decoupled ^{31}P n.m.r. spectra at 24.2 MHz of $[\text{Ph}_2\text{P}(\text{S})][\text{Ph}_2\text{P}]\text{CHLi}$ in Et_2O at 200 K. (a) Normal spectrum showing ^7Li coupling to the Ph_2P -resonance only. (b) With ^7Li decoupled at 23.2 MHz to give an AB pattern. Time sharing of the receiver and the ^7Li irradiating field was used to avoid frequency interference.



dimers or higher polymers; and (iii) it is unlikely that in (3b) 2J [$^{31}\text{P}(\text{S})$ - ^7Li] would be zero, although this possibility cannot

be completely excluded in view of the wide variability of coupling constants involving phosphorus.

We have also used ^7Li and ^{31}P n.m.r. spectroscopy to examine several other solutions of lithiated organophosphorus compounds including $(\text{Ph}_2\text{P})_3\text{CLi}$ which contained several different species showing ^{31}P - ^7Li spin coupling, and $(\text{Me}_2\text{P})_2\text{CHLi}$, $(\text{Me}_2\text{P})(\text{Ph}_2\text{P})\text{CHLi}$, and $[\text{Ph}_2\text{P}(\text{S})]_2\text{CHLi}$ which have not yet given satisfactory results. However, the spectra obtained depend upon the conditions, and indeed we have found ^7Li and ^{31}P n.m.r. spectroscopy a valuable guide to reagent quality in synthetic work.

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