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Figure 6. Total electron density surfaces (0.02 e/a.u.³) for **1** (a) and **2** (b). The same orientation of systems is maintained as in Figure 1 and Figure 4.



Figure 7. Li(SC₆H₅)·Li(CH₂MeNCH₂CH₂NMe₂) (distances in Å)



Figure 8. Li₂(SC₆H₄)·Li(CH₂MeNCH₂CH₂NMe₂) (distances in Å)



Figure 9. Li₂{(2-S-3-SiMe₃-C₆H₃)·Li(CH₂MeNCH₂CH₂NMe₂) (distances in Å)