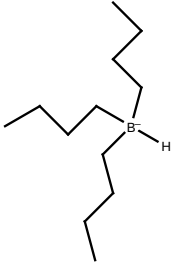
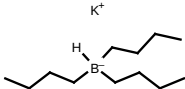
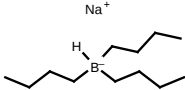
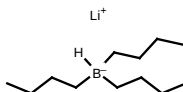


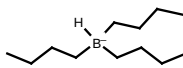
Query

	Query	Results	Date
1. Query	 Search as: As drawn	5 substances	2011-11-15 08h:06m:33s (EST)

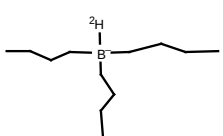
Reaxys ID 4006677 View in Reaxys		1/5
		Chemical Name: potassium-selectride=Kalium-selectrid; nBu3BHK Molecular Formula: C ₁₂ H ₂₈ B*K CAS Registry Number: 67940-39-0 Molecular Weight: 222.264 Linear Structure Formula: K ⁽¹⁺⁾ *BH(C ₄ H ₉) ₃ ⁽¹⁻⁾ =KBH(C ₄ H ₉) ₃ Type of Substance: acyclic; Coordination compound InChIKey: DMVRIUSSUUTYFE-UHFFFAOYSA-N Note:
NMR Spectroscopy (1)		
1 of 1	Description	NMR
	Comment	11B
	Brown; Krishnamurthy ; Journal of Organometallic Chemistry; vol. 156; (1978); p. 111,117, View in Reaxys	
IR Spectroscopy (1)		
Description		References
IR		Brown; Krishnamurthy ; Journal of Organometallic Chemistry; vol. 156; (1978); p. 111,117, View in Reaxys

Reaxys ID 4157405 View in Reaxys		2/5
		Chemical Name: sodium tri-n-butylhydroborate; Natrium-tri-n-butylborhydrid; Natrium-tributyl-boranat Molecular Formula: C ₁₂ H ₂₈ B*Na Molecular Weight: 206.155 Linear Structure Formula: Na ⁽¹⁺⁾ *HB(C ₄ H ₉) ₃ ⁽¹⁻⁾ =NaBH(C ₄ H ₉) ₃ Type of Substance: acyclic; Coordination compound InChIKey: INBGIUCSCRWOHB-UHFFFAOYSA-N Note:
Melting Point (2)		
Melting Point [°C]	References	
62	Binger,P. et al. ; Justus Liebigs Annalen der Chemie; vol. 717; (1968); p. 21 - 40, View in Reaxys	
62	Binger, P.; benedikt, G.; Rotermund, G. W.; Koester, R. ; Liebigs Annalen der Chemie; vol. 717; (1968); p. 21 - 40, View in Reaxys ; vol. B: B-Verb.8; 4, page 66 - 72 ; (from Gmelin) , View in Reaxys	
NMR Spectroscopy (2)		
1 of 2	Nucleus	11B
	Solvents	toluene; tetrahydrofuran
	Comment	Signals given
	Hubbard, John L. ; Journal of the Chemical Society, Chemical Communications; (1989); p. 1639 - 1640 ; (from Gmelin) , View in Reaxys	
2 of 2	Description	NMR
	Comment	11B
	Brown,H.C. et al. ; Journal of the American Chemical Society; vol. 100; (1978); p. 3343 - 3349, View in Reaxys	
IR Spectroscopy (3)		
Description	Solvent	References
IR		Brown,H.C. et al. ; Journal of the American Chemical Society; vol. 100; (1978); p. 3343 - 3349, View in Reaxys ; Binger,P. et al. ; Justus Liebigs Annalen der Chemie; vol. 717; (1968); p. 21 - 40, View in Reaxys
Bands	cyclohexane	Binger, P.; benedikt, G.; Rotermund, G. W.; Koester, R. ; Liebigs Annalen der Chemie; vol. 717; (1968); p. 21 - 40 ; (from Gmelin) , View in Reaxys
Bands	cyclohexane	vol. B: B-Verb.8; 4, page 66 - 72 ; (from Gmelin) , View in Reaxys

Reaxys ID 4007576 View in Reaxys		3/5	
		Chemical Name: LiHB(s-Bu)3; Lithium-tri-n-butylborhydrid; LiH(n-Bu)3B; n-Bu3BHLi; LiBHBu3 Molecular Formula: C ₁₂ H ₂₈ B*Li CAS Registry Number: 67335-72-2 Molecular Weight: 190.106 Linear Structure Formula: Li ⁽¹⁺⁾ *C ₁₂ H ₂₈ B ⁽¹⁻⁾ Type of Substance: acyclic; Coordination compound InChIKey: XEGHKBNMKSZTJS-UHFFFAOYSA-N Note:	
NMR Spectroscopy (3)			
1 of 3	Nucleus	11B	
	Coupling Nuclei	1H	
	Solvents	hexane; tetrahydrofuran	
	Comment	Signals given	
	Biffar, Werner; Noth, Heinrich; Sedlak, Dieter; Organometallics; vol. 2; (1983); p. 579 - 585 ; (from Gmelin) , View in Reaxys		
2 of 3	Nucleus	11B	
	Coupling Nuclei	1H	
	Solvents	tetrahydrofuran; pentane	
	Comment	Signals given	
	Brown, Herbert C.; Kramer, Gary W.; Hubbard, John L.; Krishnamurthy, S.; Journal of Organometallic Chemistry; vol. 188; (1980); p. 1 - 10 ; (from Gmelin) , View in Reaxys		
3 of 3	Description	NMR	
	Comment	11B	
	Brown,H.C. et al.; Journal of the American Chemical Society; vol. 100; (1978); p. 3343 - 3349, View in Reaxys		
IR Spectroscopy (2)			
Description	Solvent	Temperature [°C]	References
Bands	tetrahydrofuran	25	Brown, Herbert C.; Kramer, Gary W.; Hubbard, John L.; Krishnamurthy, S.; Journal of Organometallic Chemistry; vol. 188; (1980); p. 1 - 10 ; (from Gmelin) , View in Reaxys
IR			Brown,H.C. et al.; Journal of the American Chemical Society; vol. 100; (1978); p. 3343 - 3349, View in Reaxys

Reaxys ID 4126433 View in Reaxys		4/5
		Chemical Name: tributyl-hydrido-borate(1-); Tri-n-butylborhydrid-Anion; tri-n-butylborohydride Molecular Formula: C ₁₂ H ₂₈ B Molecular Weight: 183.165 Linear Structure Formula: C ₁₂ H ₂₈ B ⁽¹⁻⁾ Type of Substance: acyclic InChIKey: NMXXHSNUFXSPTG-UHFFFAOYSA-N Note:
Derivative (2)		
Comment	References	
Li-Salz: B.: B(n-Bu) ₃ , LiAlH ₄ , Tri-aethylendiamin; IR; 11B-NMR (Tab.I)	Brown et al.; Journal of Organic Chemistry; vol. 44; (1979); p. 5004, View in Reaxys	

Li ⁺ -Salz: aus n-Bu ₃ B, LiAlH(OMe) ₃	Brown et al. ; Journal of Organometallic Chemistry; vol. 166; (1979); p. 271,274, 275, 276, 278, View in Reaxys
Further Information (2)	
Description	References
Further information	Brown et al. ; Journal of Organometallic Chemistry; vol. 166; (1979); p. 271,274, 275, 276, 278, View in Reaxys
Further information	Brown et al. ; Journal of Organic Chemistry; vol. 44; (1979); p. 5004, View in Reaxys

Reaxys ID 17140631 View in Reaxys		5/5
	Molecular Formula: C ₁₂ H ₂₈ B ⁺ K Molecular Weight: 223.256 Linear Structure Formula: $K^{(1+)}B^{(2)}H((CH_2)_3CH_3)_3^{(1-)}=K(B^{(2)}H((CH_2)_3CH_3)_3)$ Type of Substance: Isotope or isotope containing compound InChIKey: DMVRIUSSUUTYFE-WSCBMPOBSA-N Note:	