

Electronic Supporting Information (ESI)

Organomagnesium amide catalyzed cross-dehydrocoupling of organosilanes with amines†

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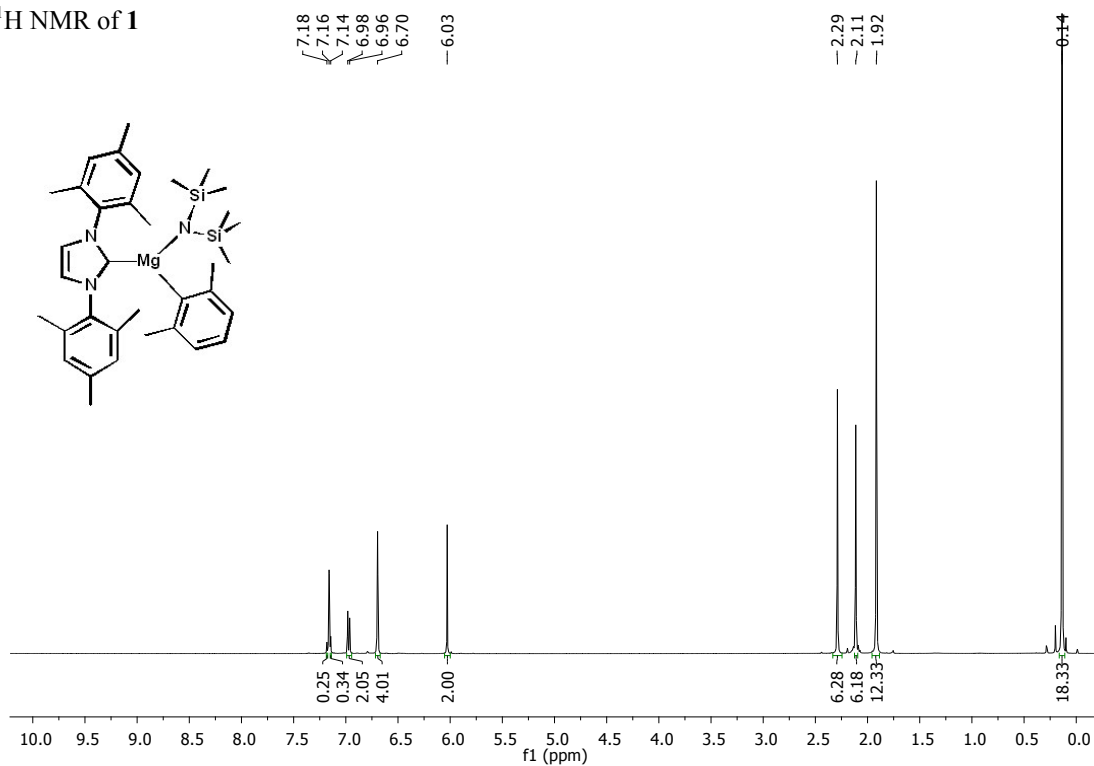
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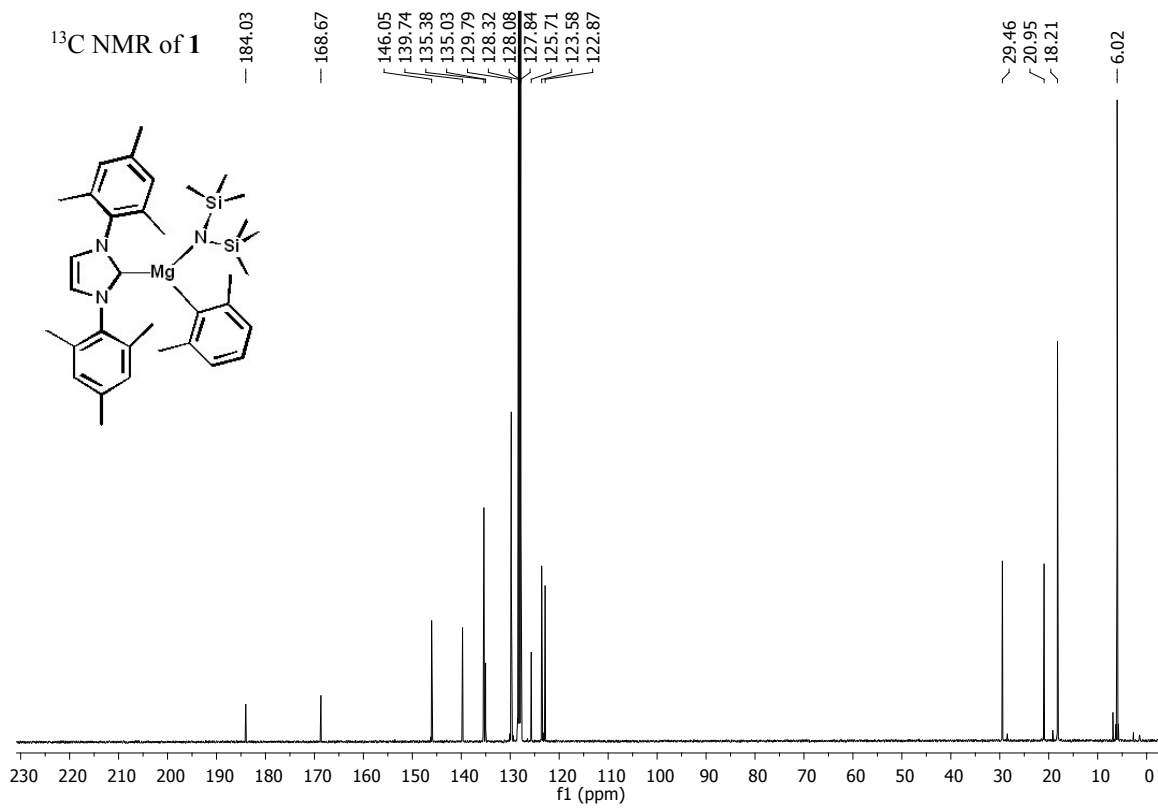
Copy of ^1H , ^{13}C & ^{29}Si NMR spectra of 1 & 2 -----	S2-S4
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Copy of ^1H , ^{13}C & ^{29}Si NMR spectra of 1 & 2:

^1H NMR of 1

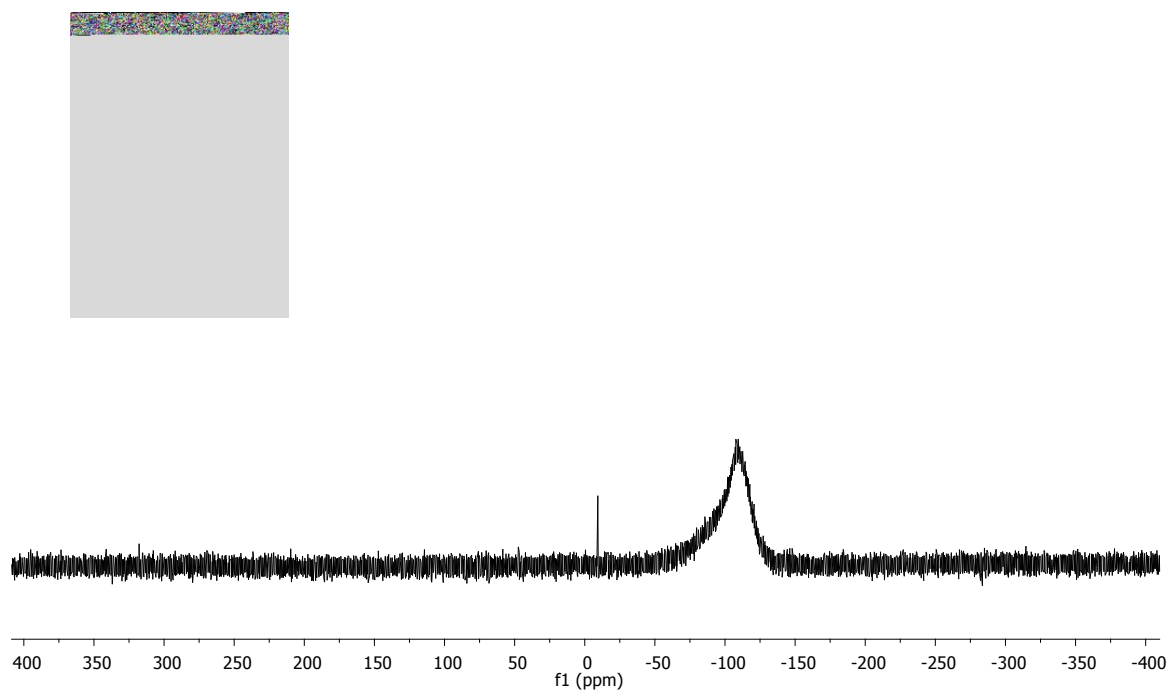


^{13}C NMR of 1

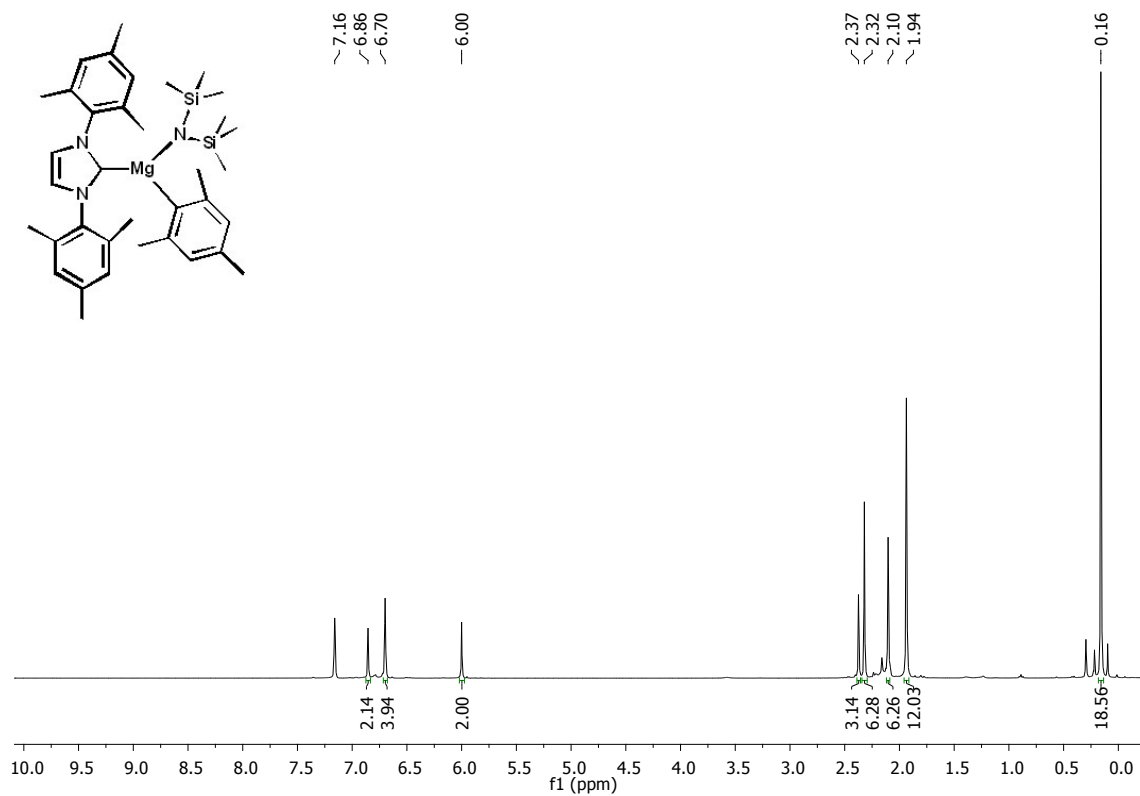


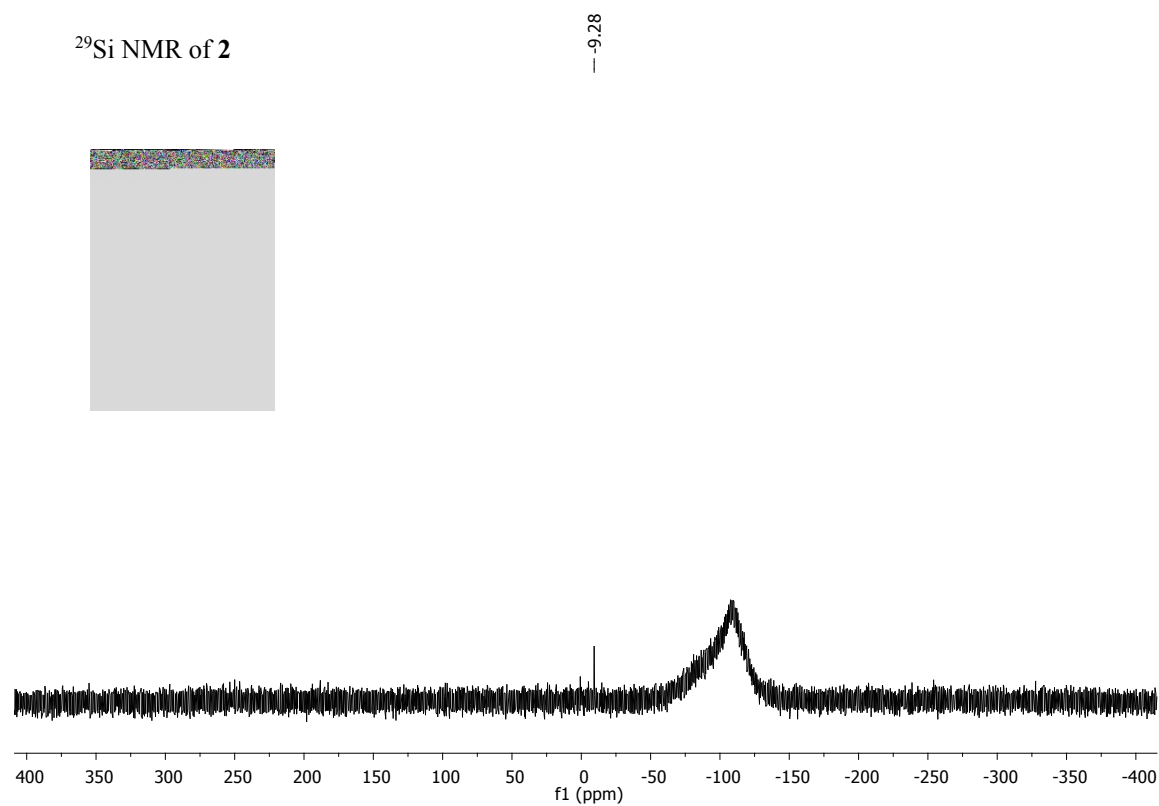
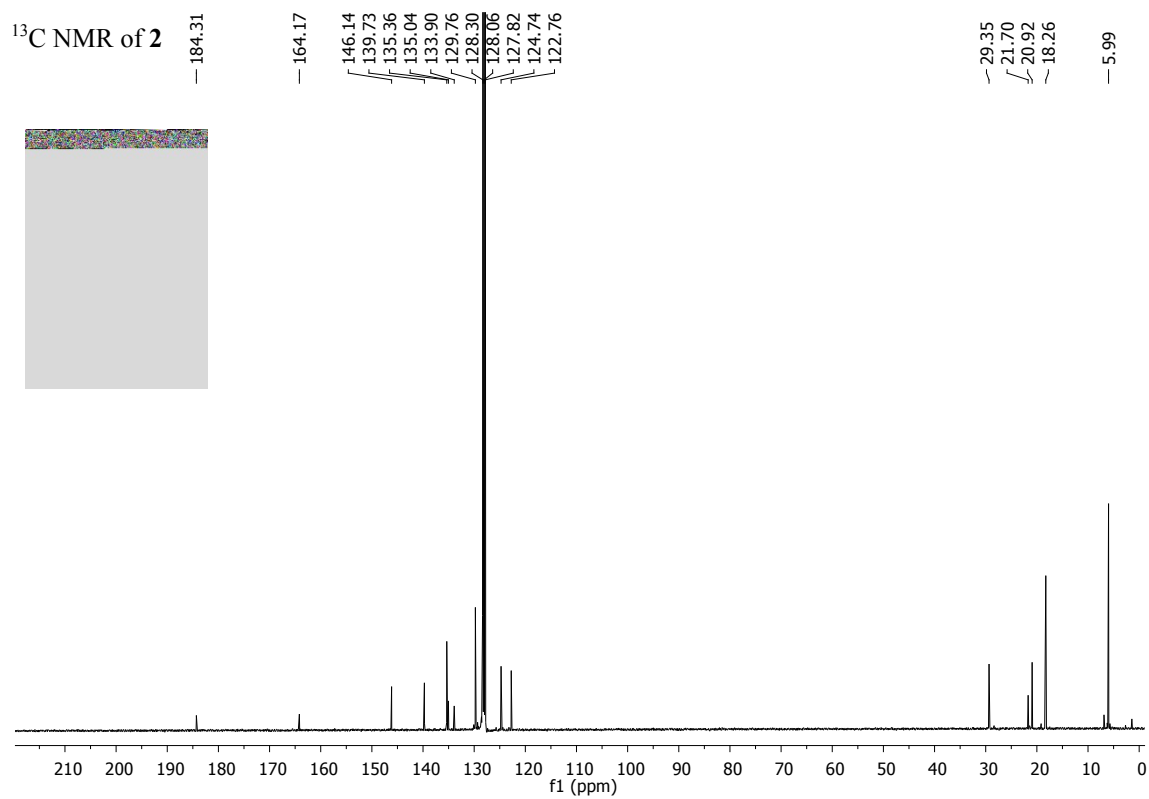
^{29}Si NMR of **1**

— -9.26

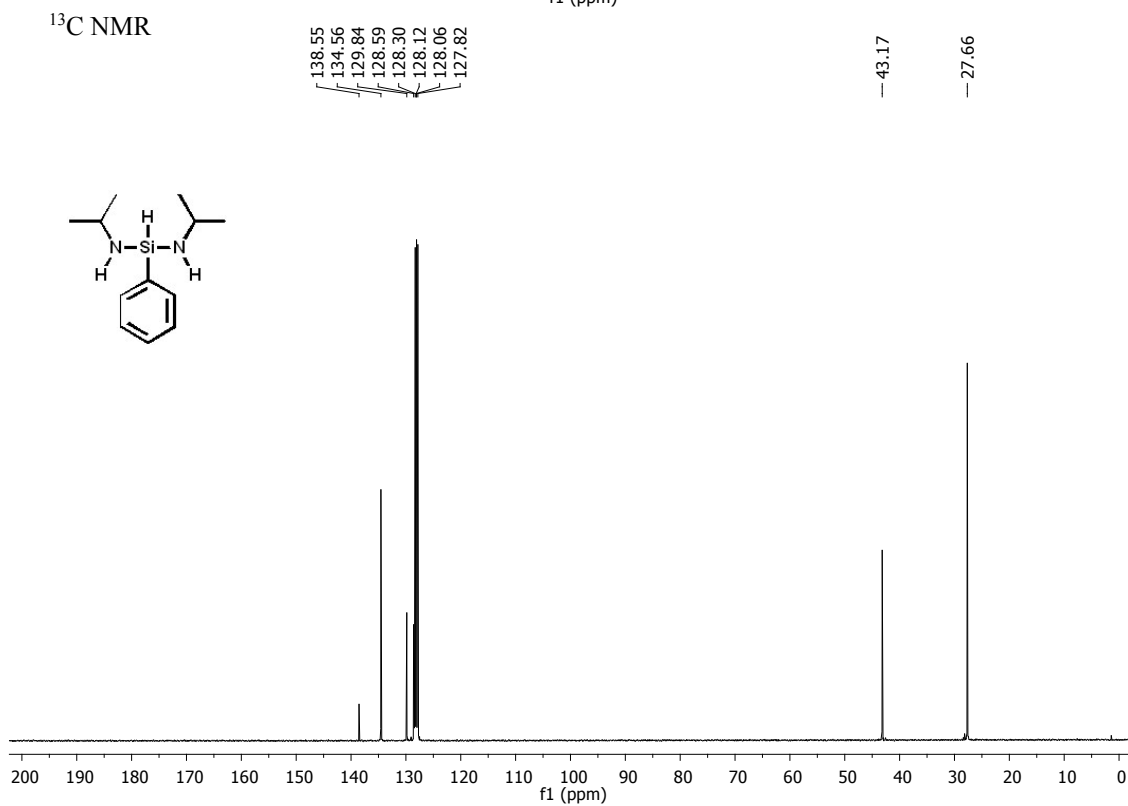
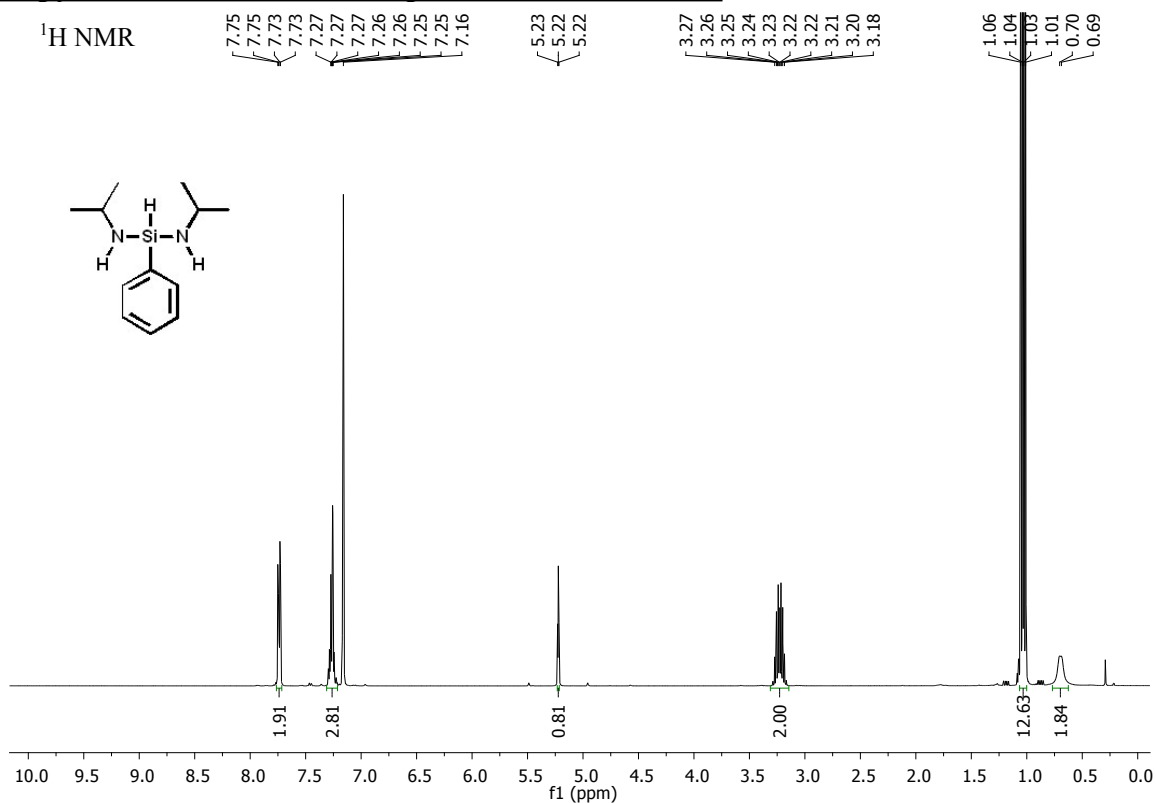


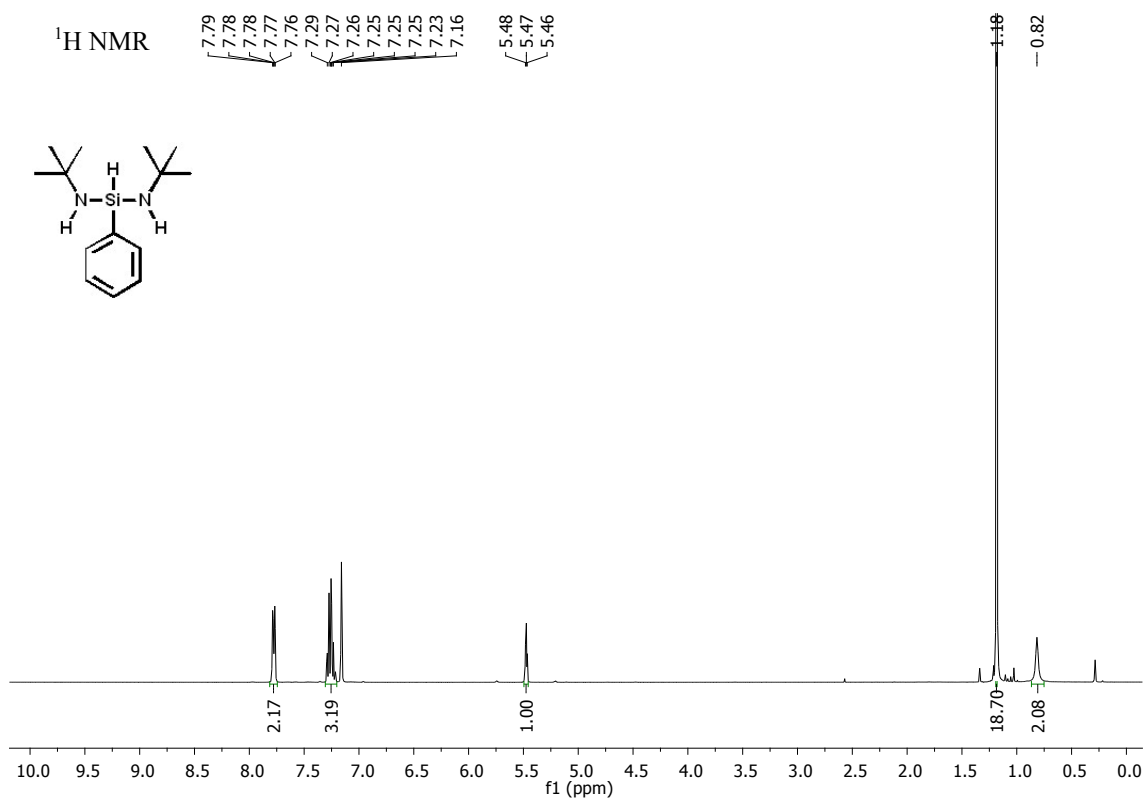
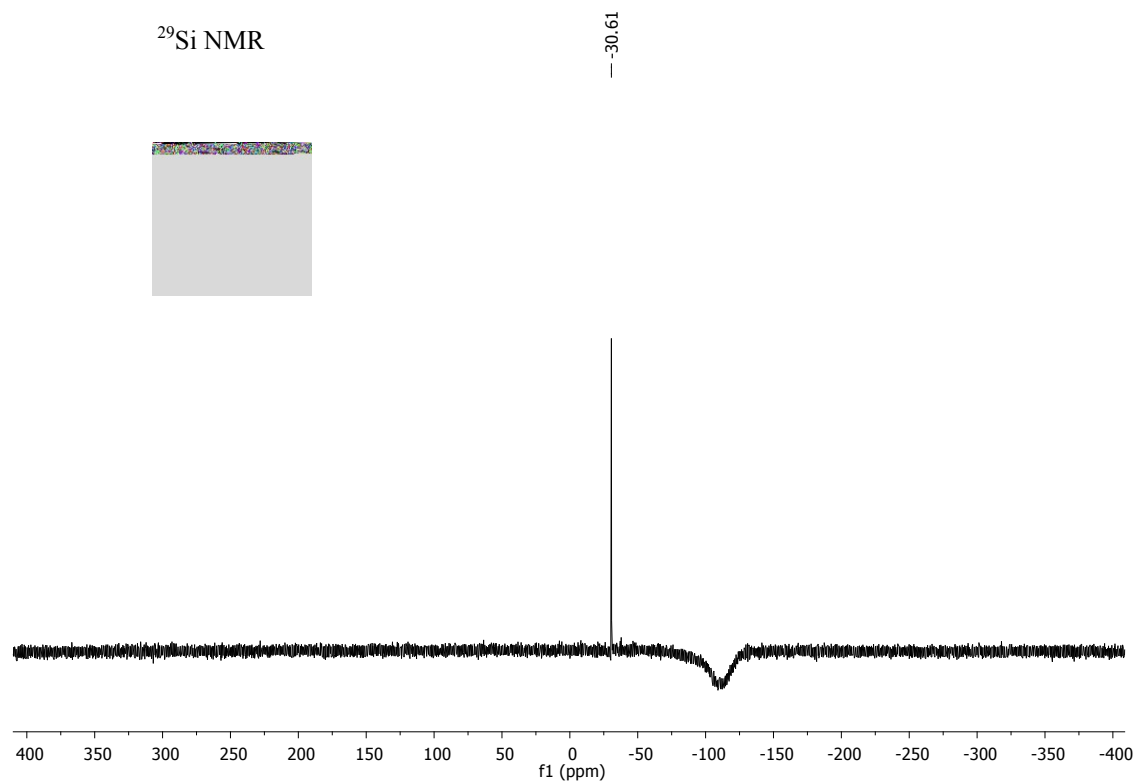
^1H NMR of **2**

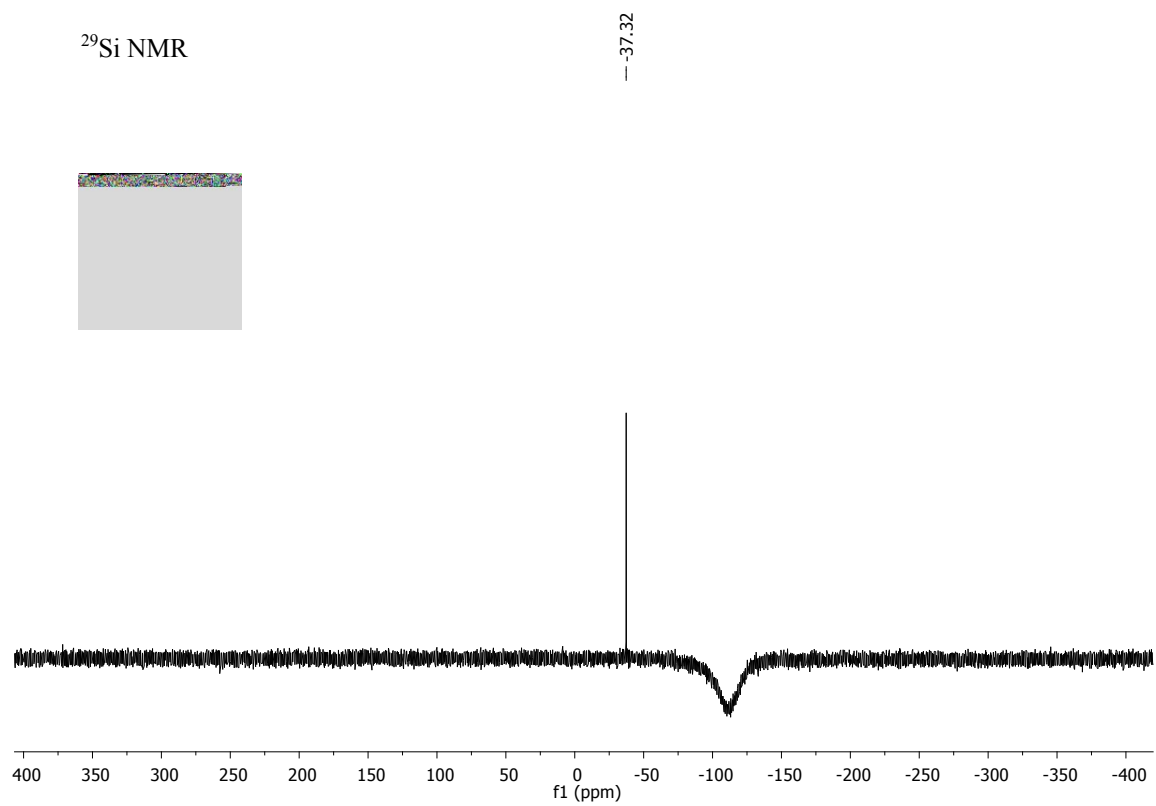
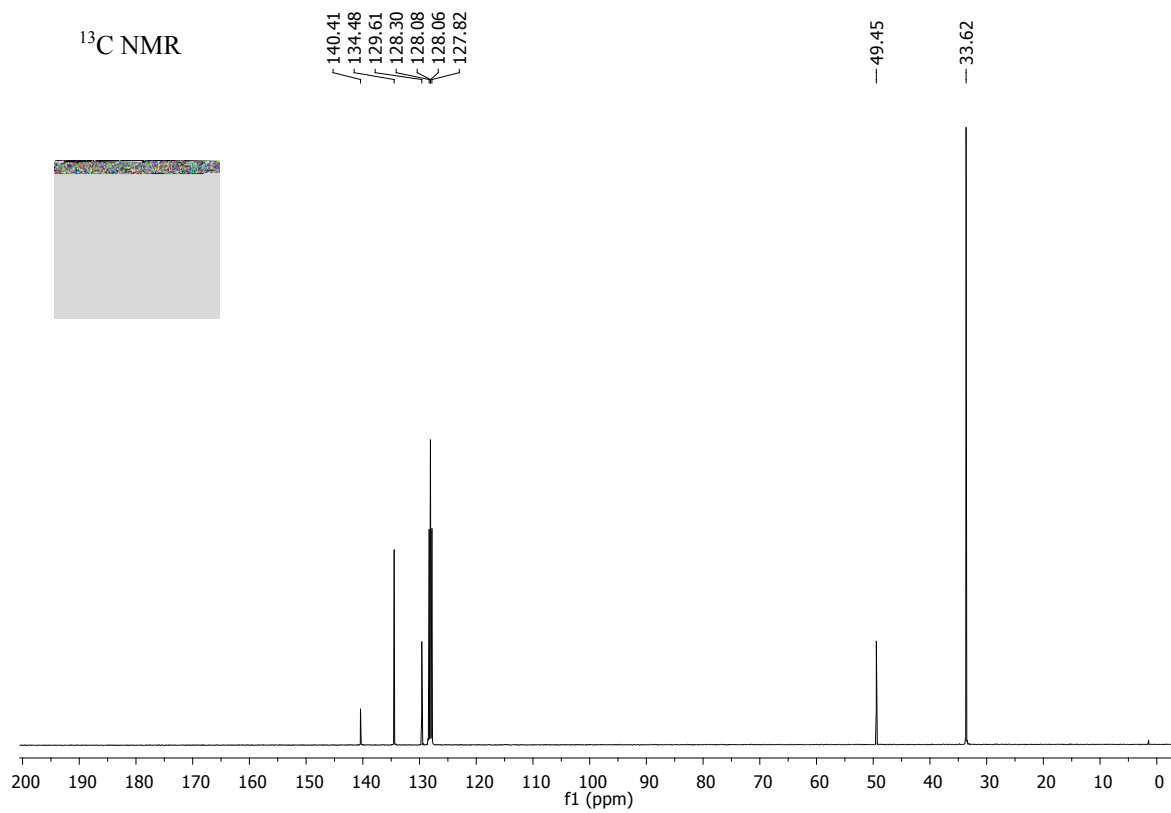


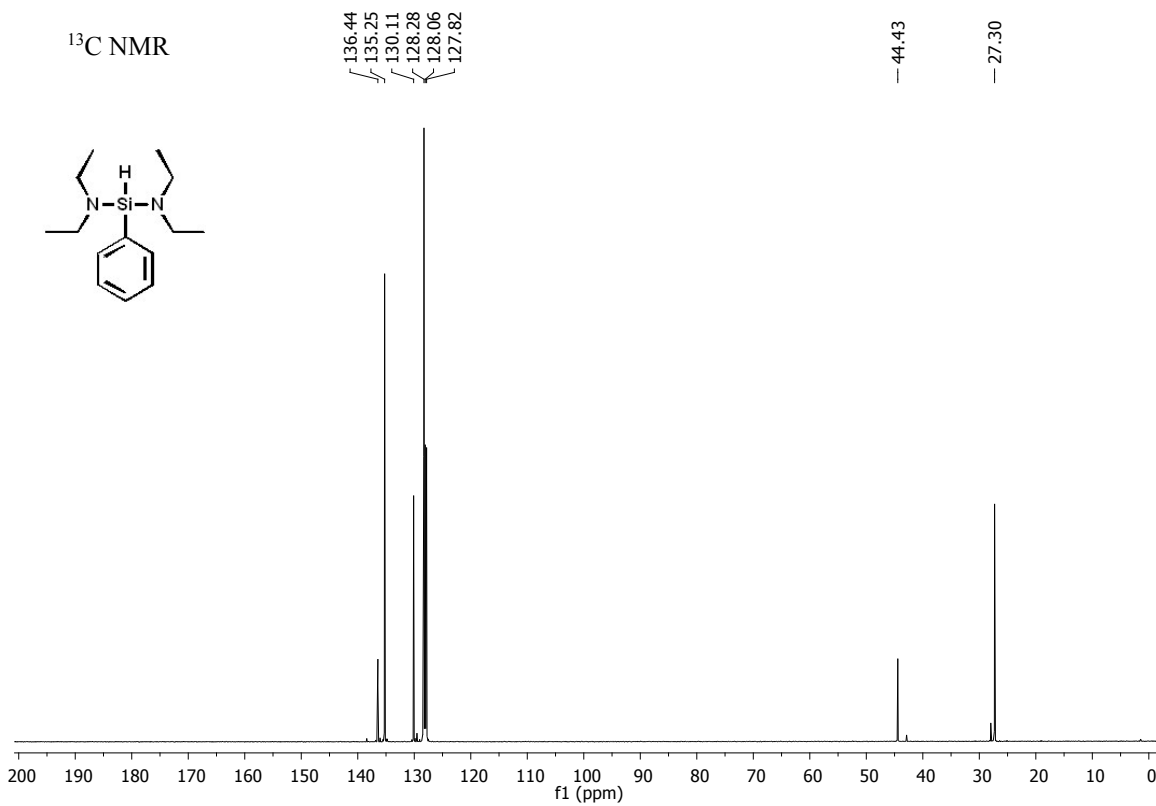
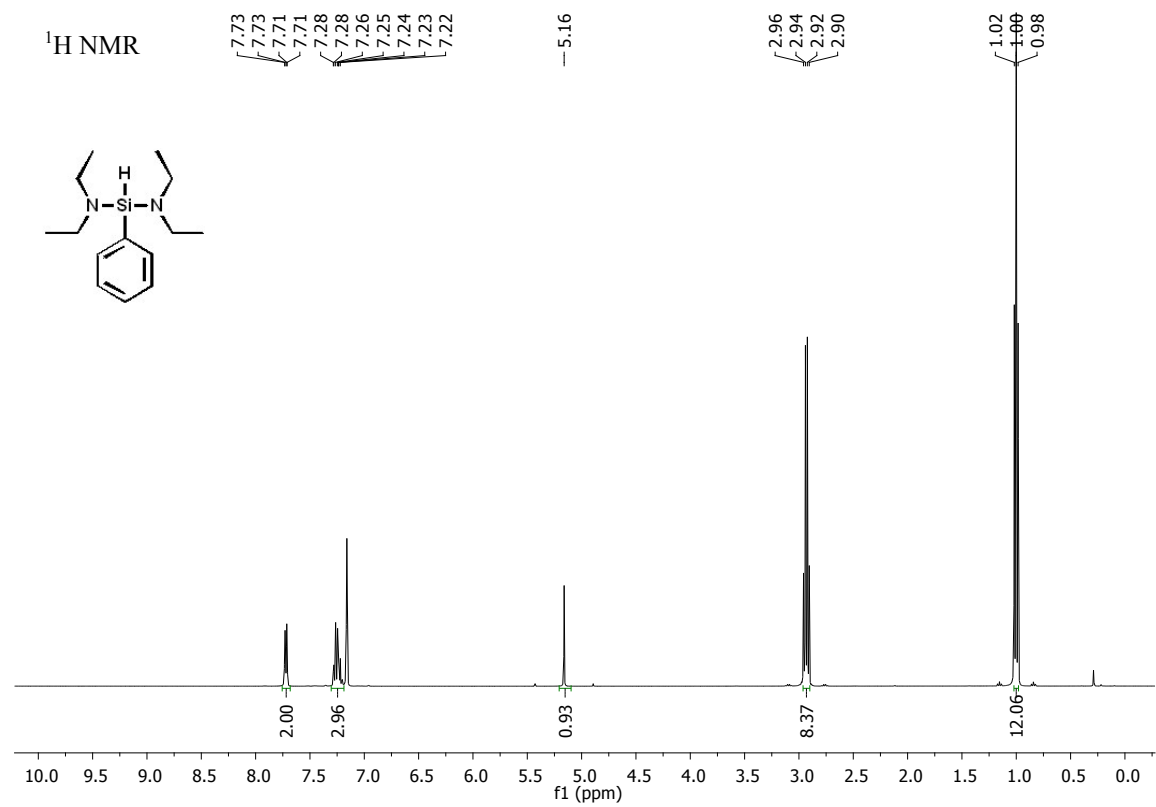


Copy of ^1H , ^{13}C & ^{29}Si NMR spectra of aminosilanes:

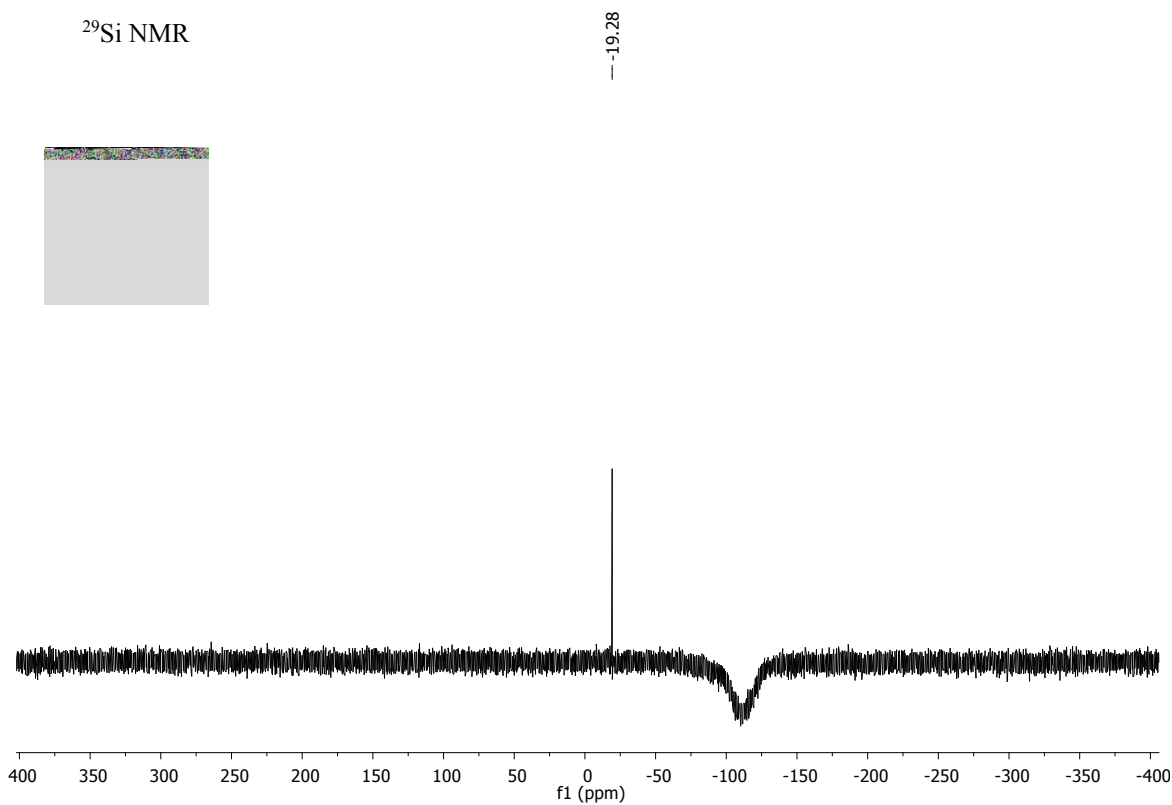




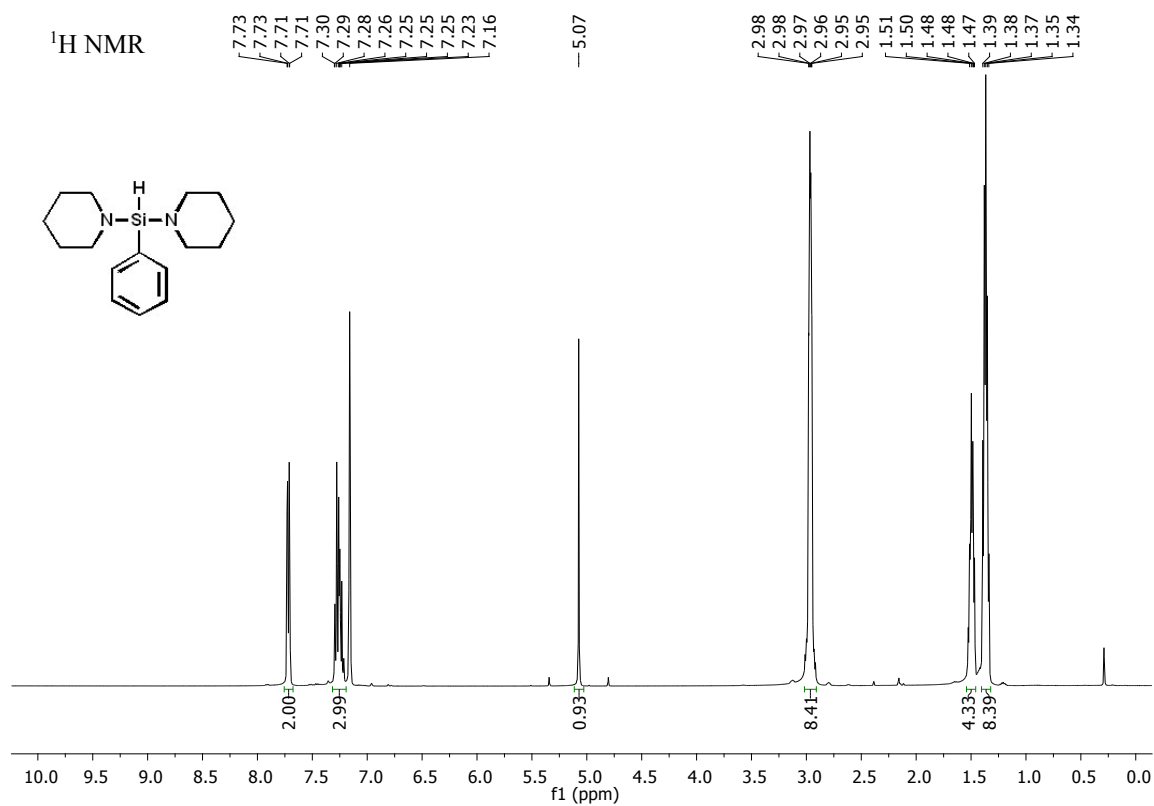


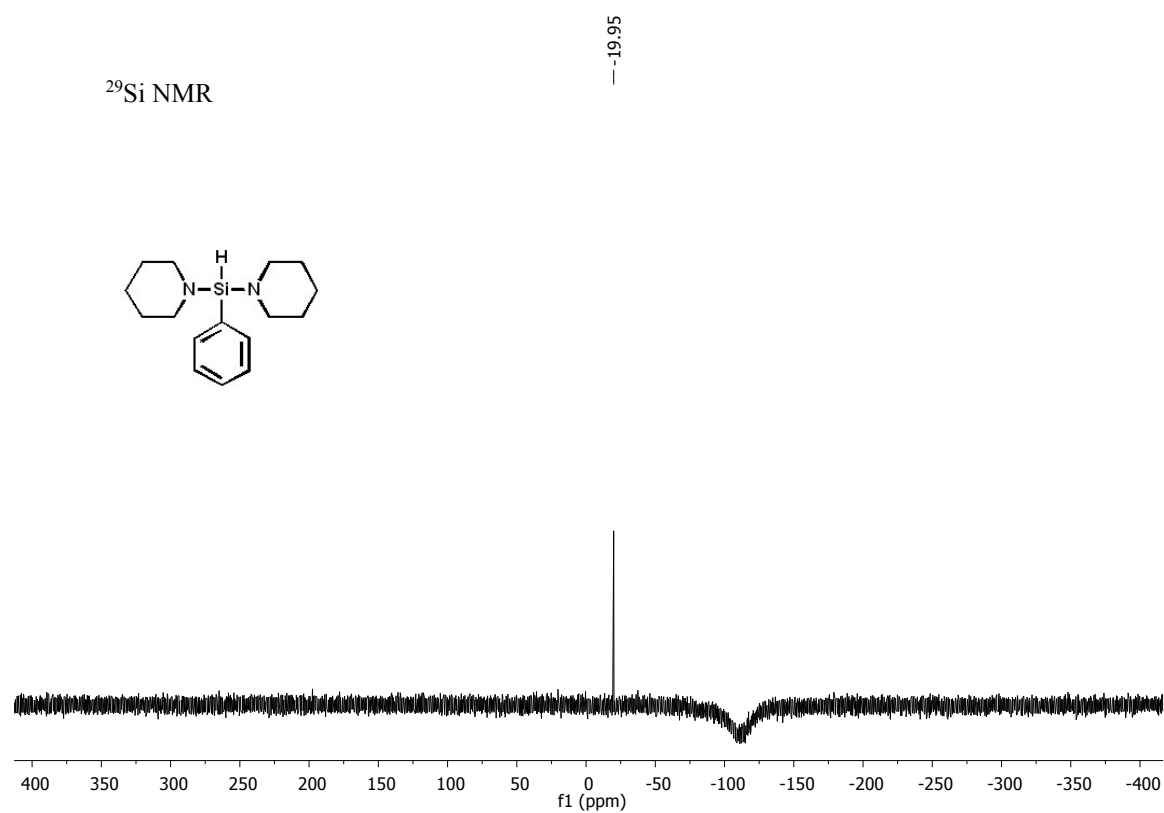
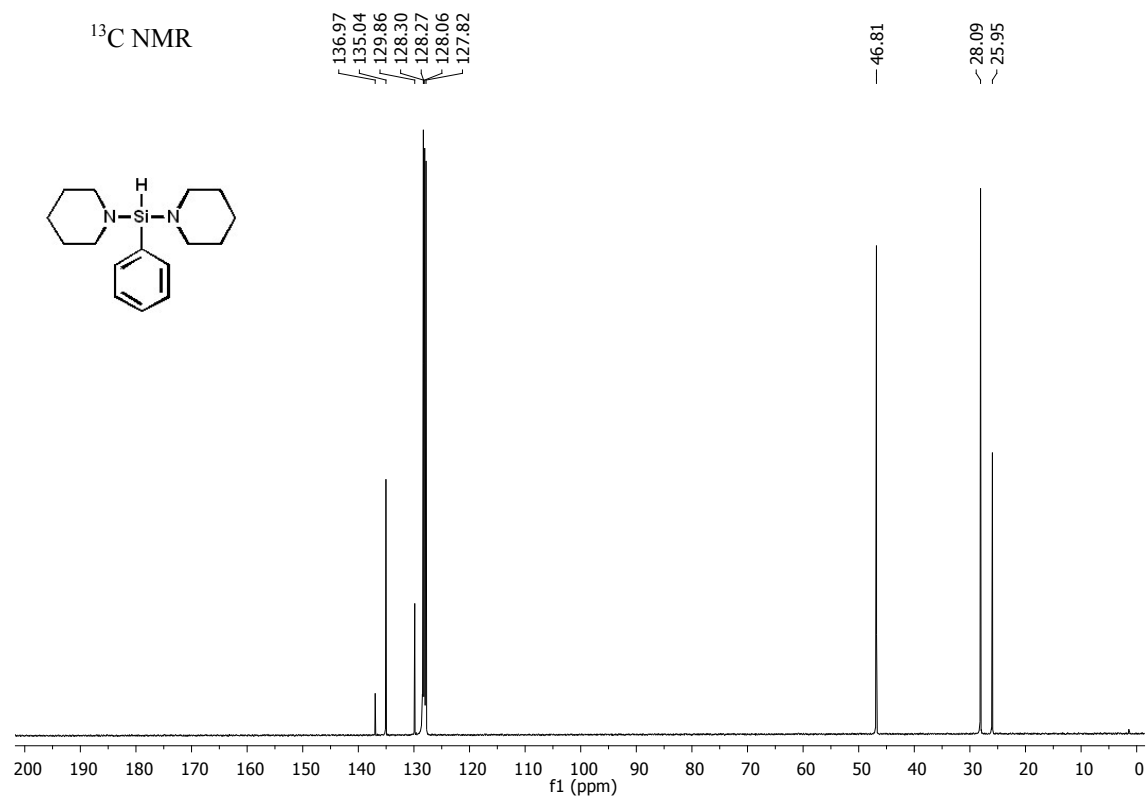


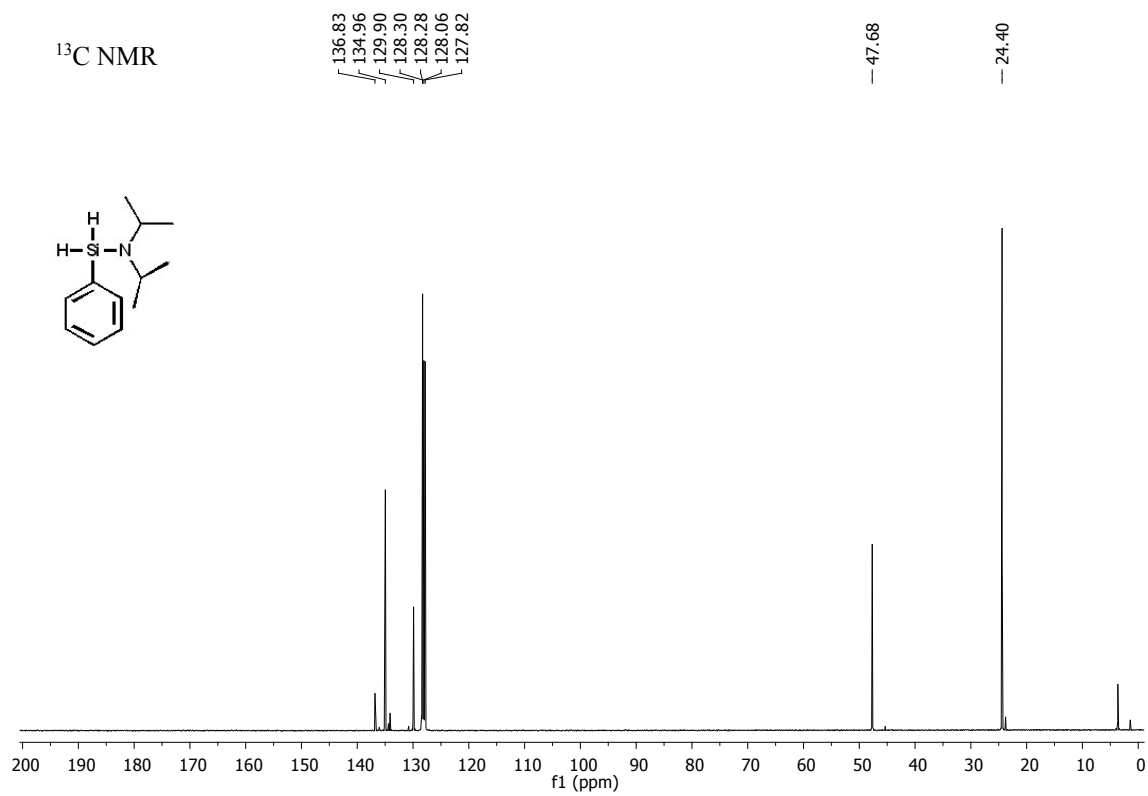
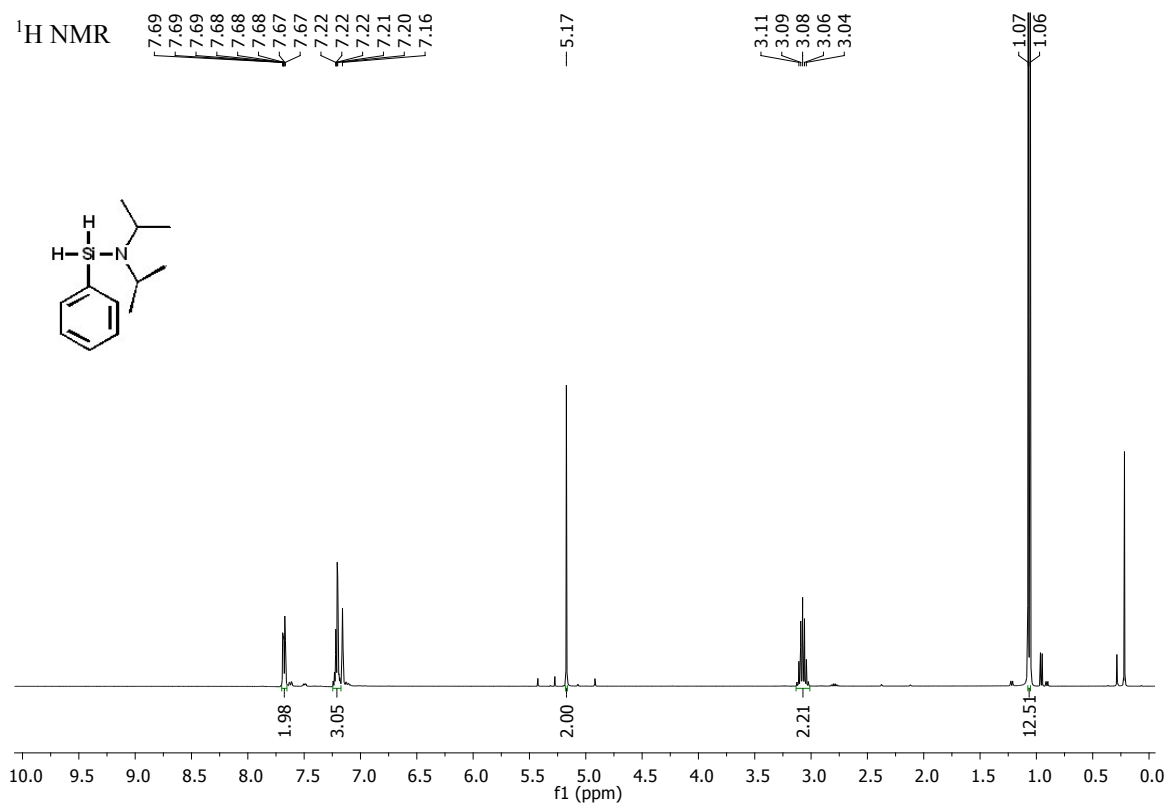
^{29}Si NMR



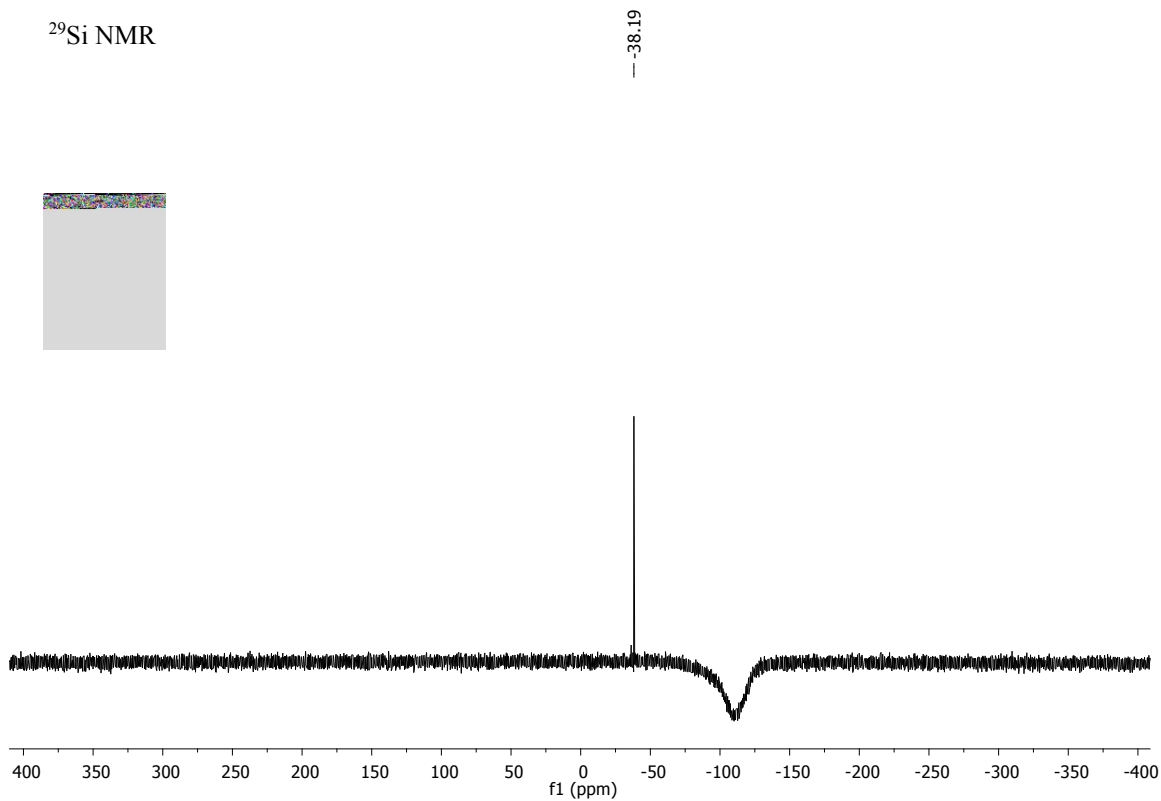
^1H NMR



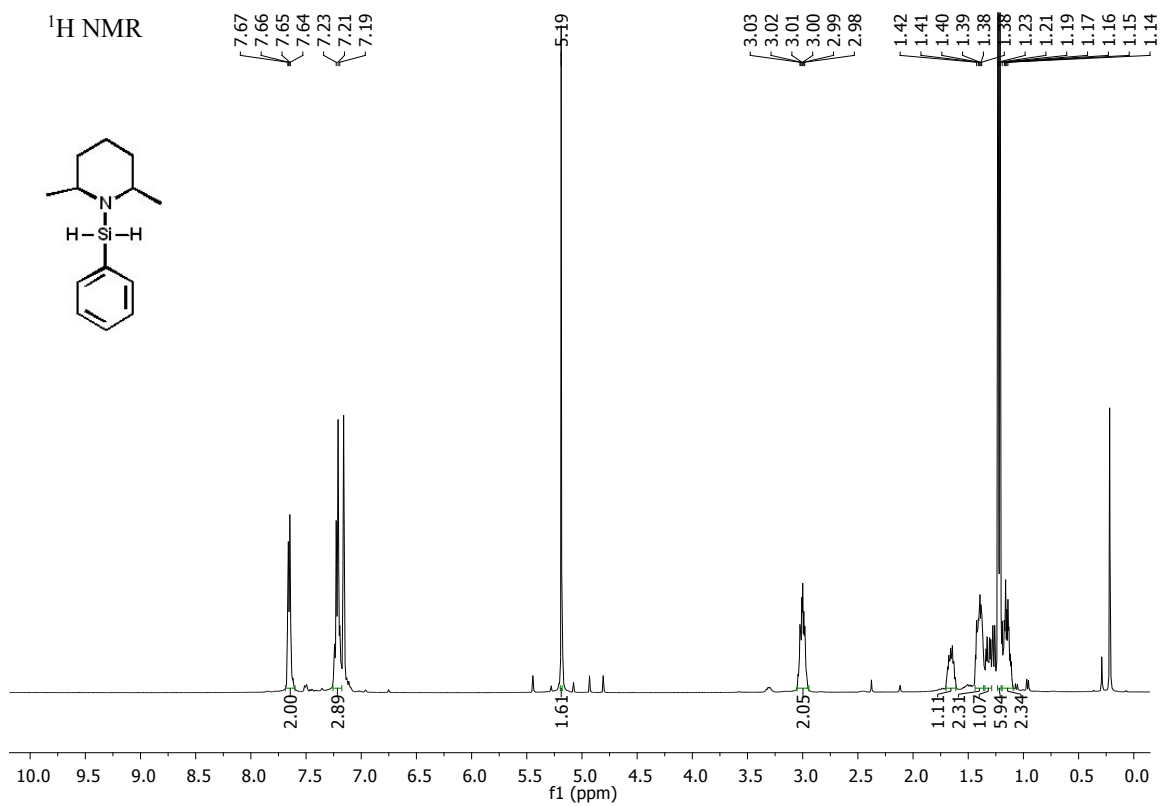




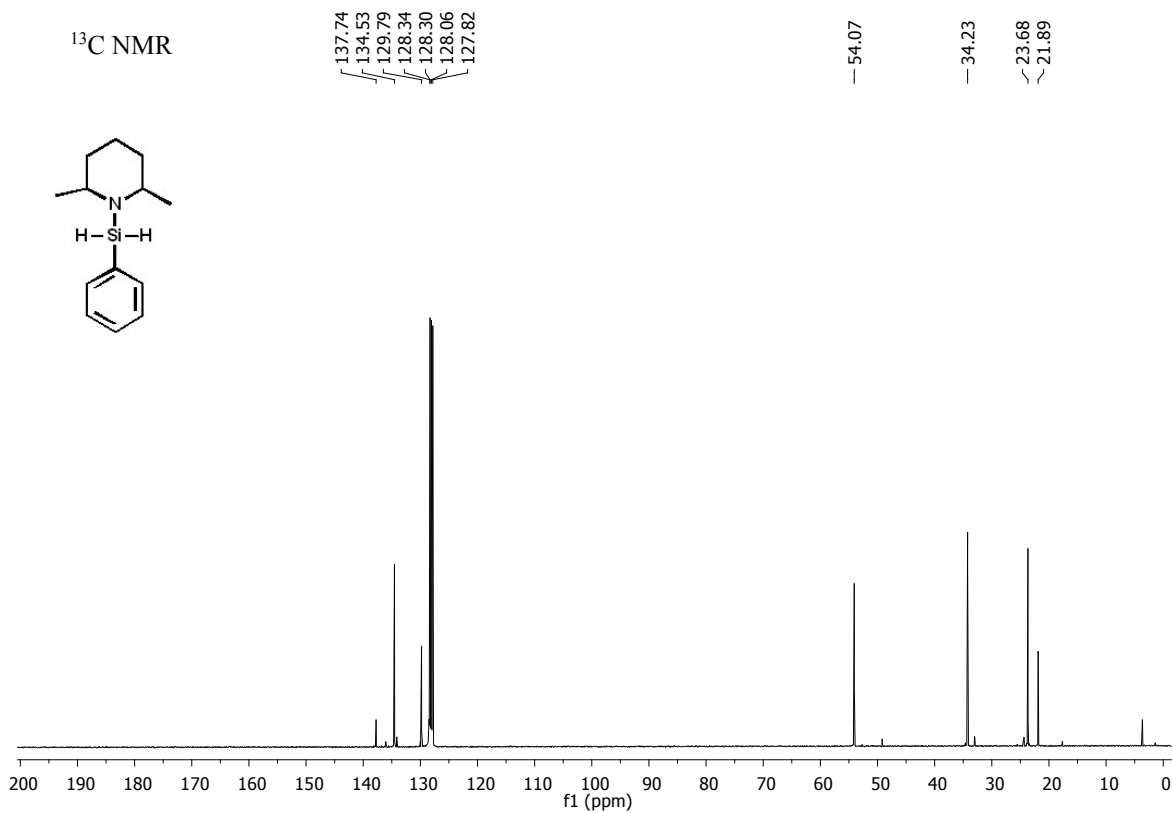
^{29}Si NMR



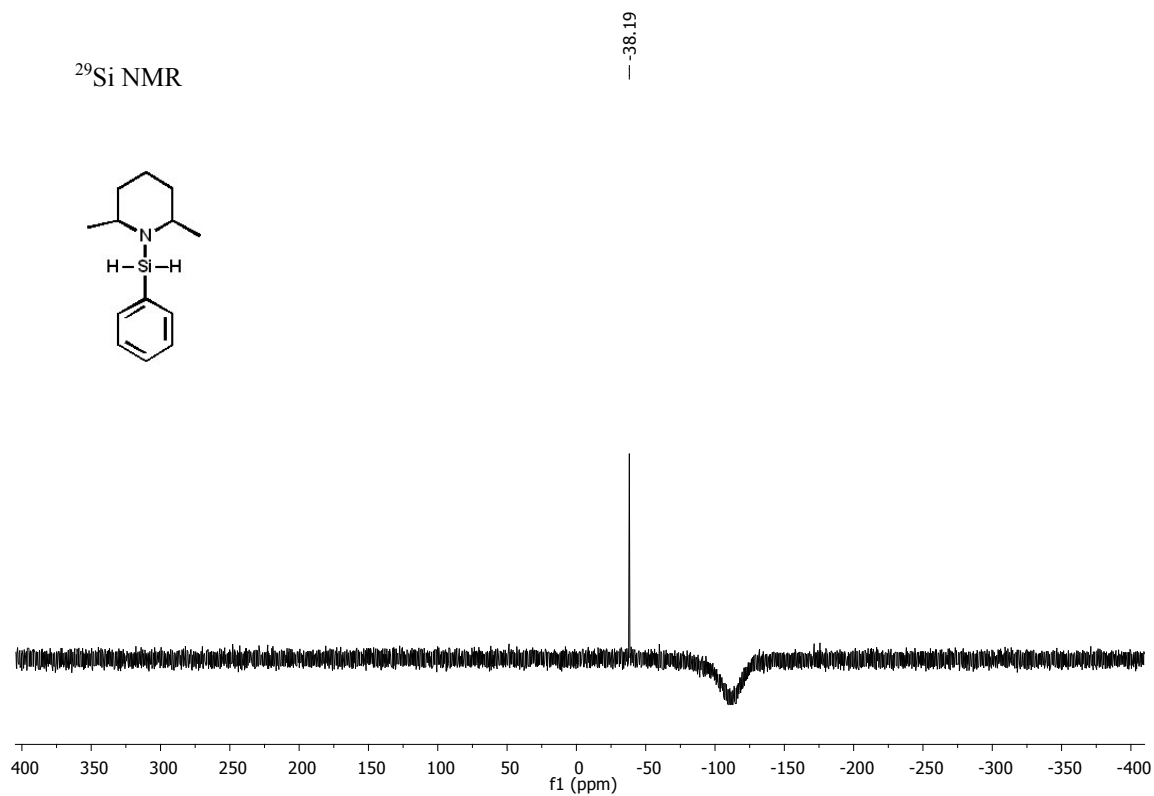
^1H NMR



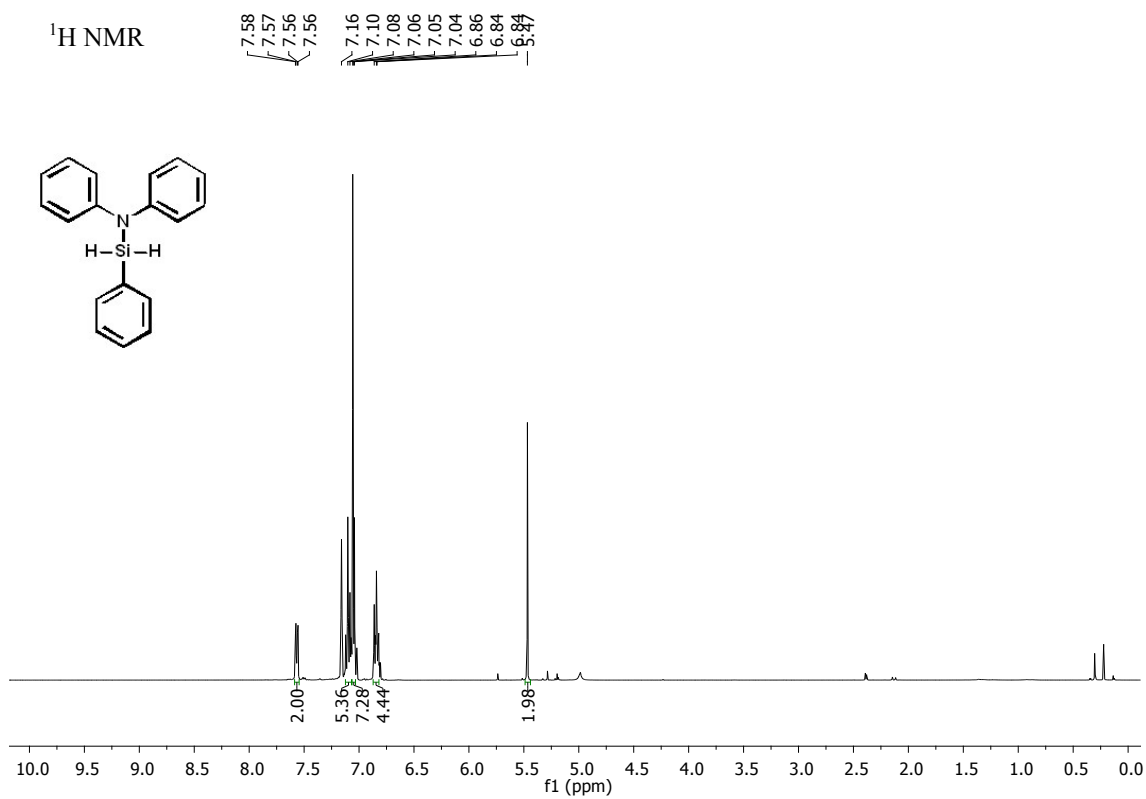
^{13}C NMR



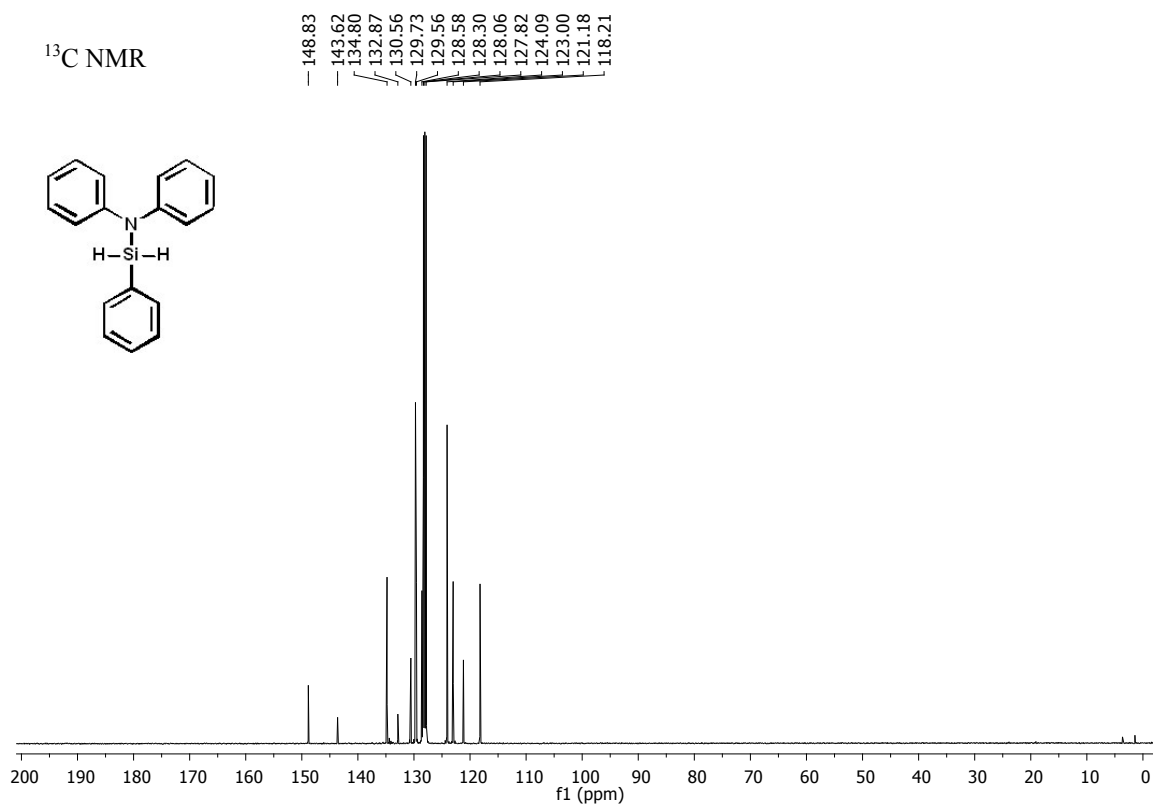
^{29}Si NMR

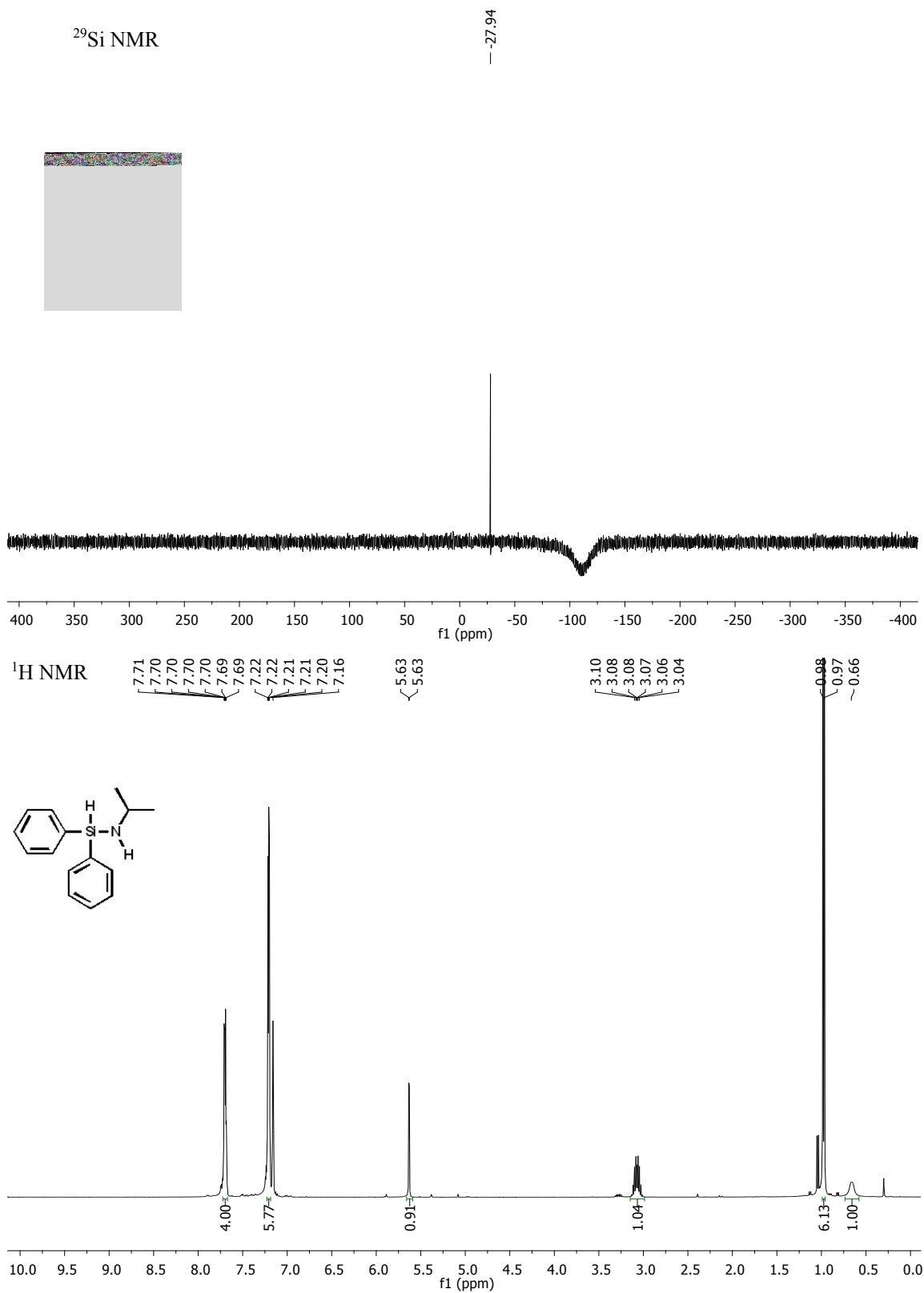


¹H NMR

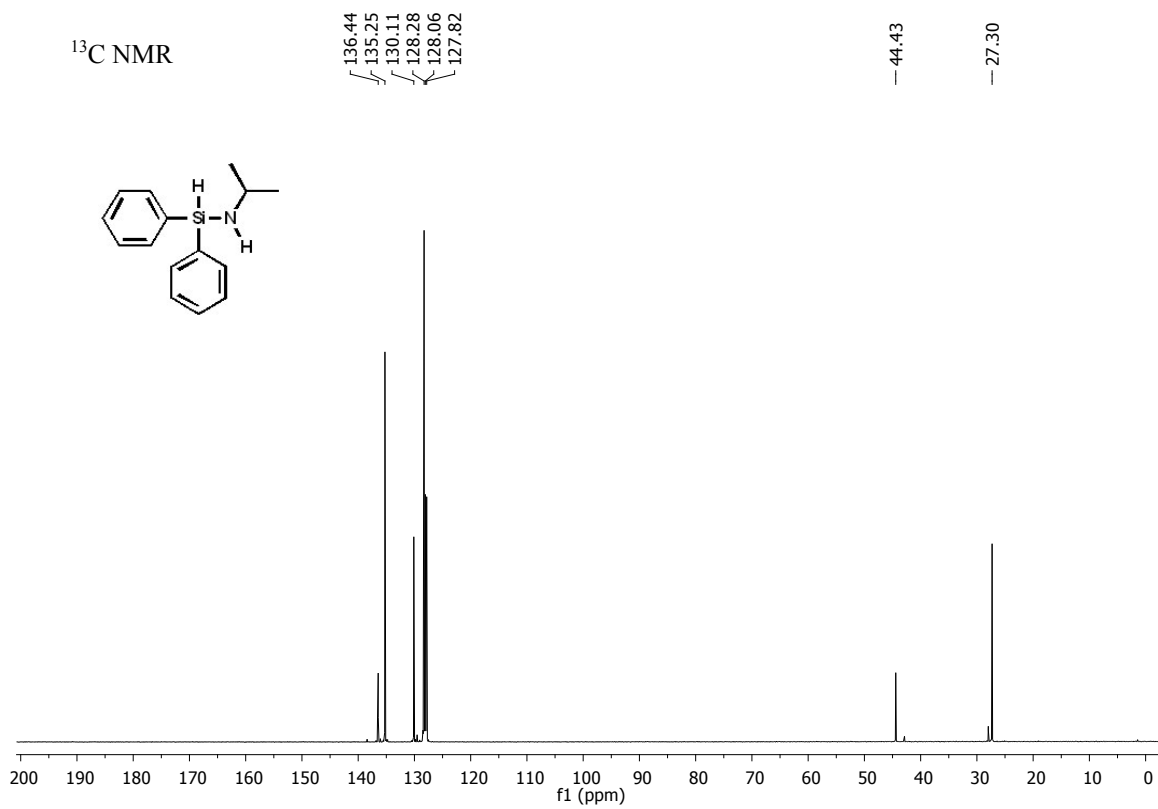


¹³C NMR

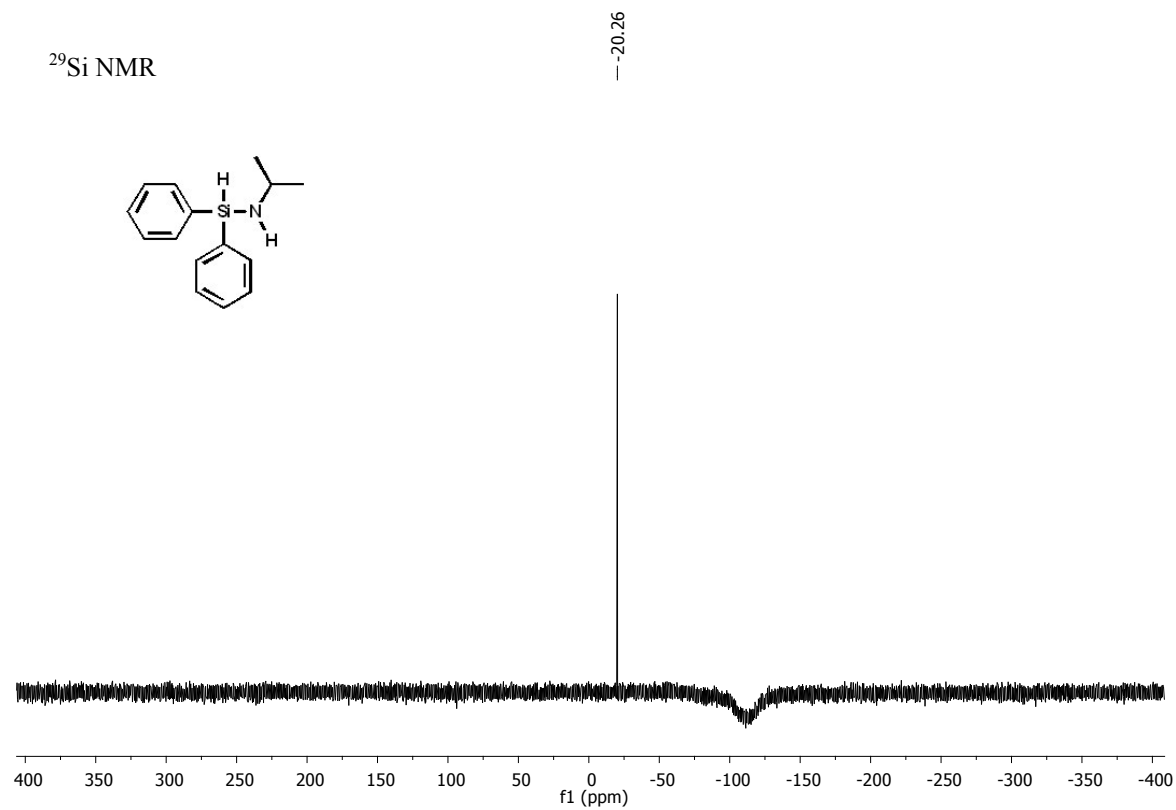


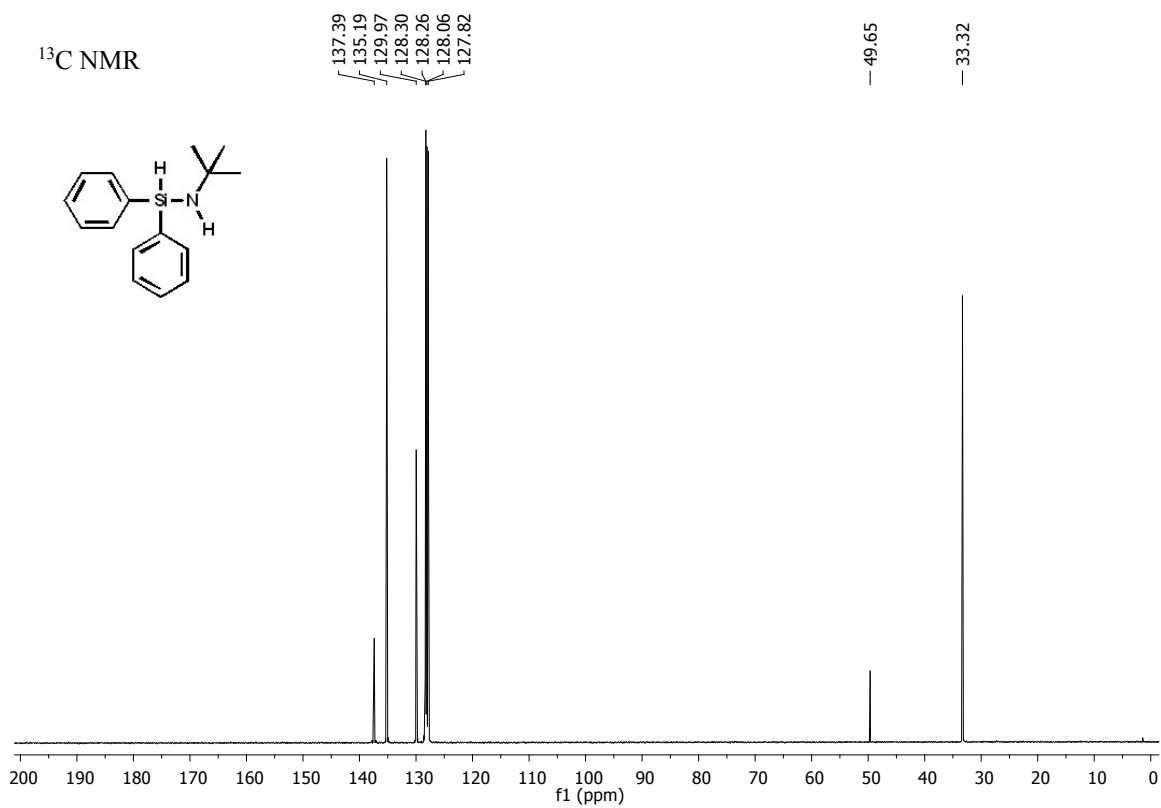
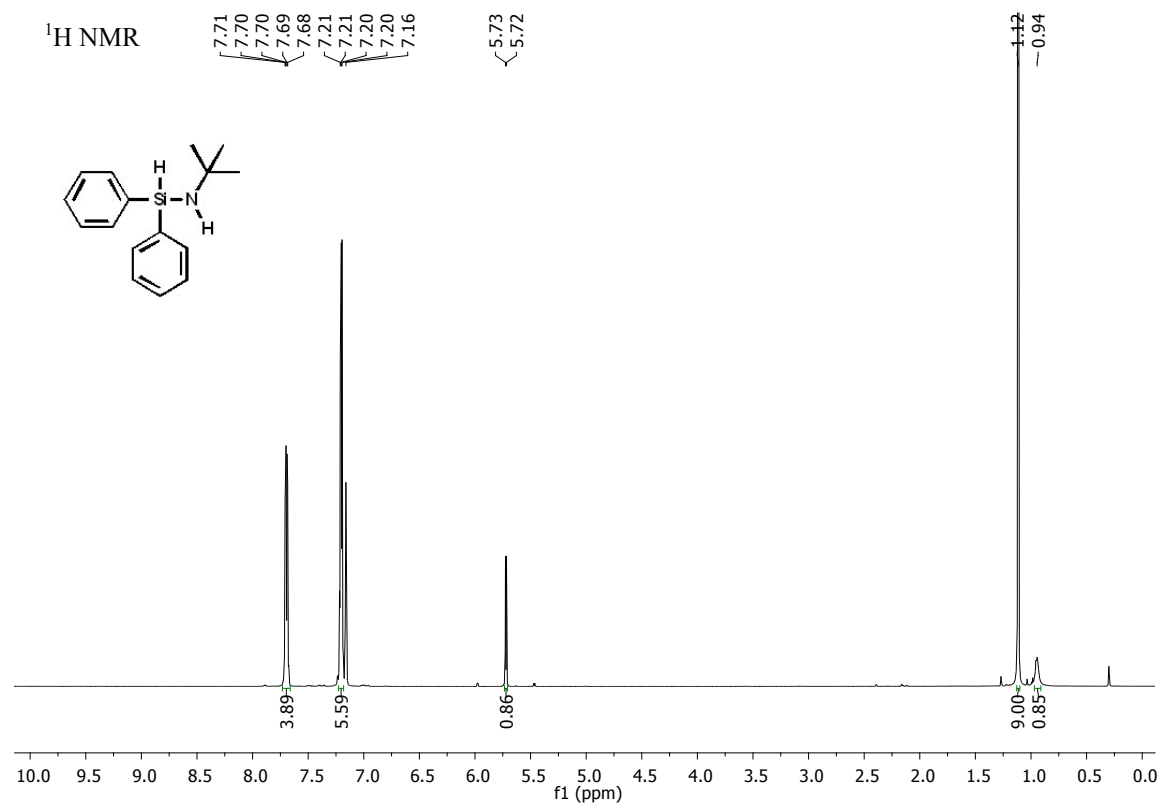


^{13}C NMR



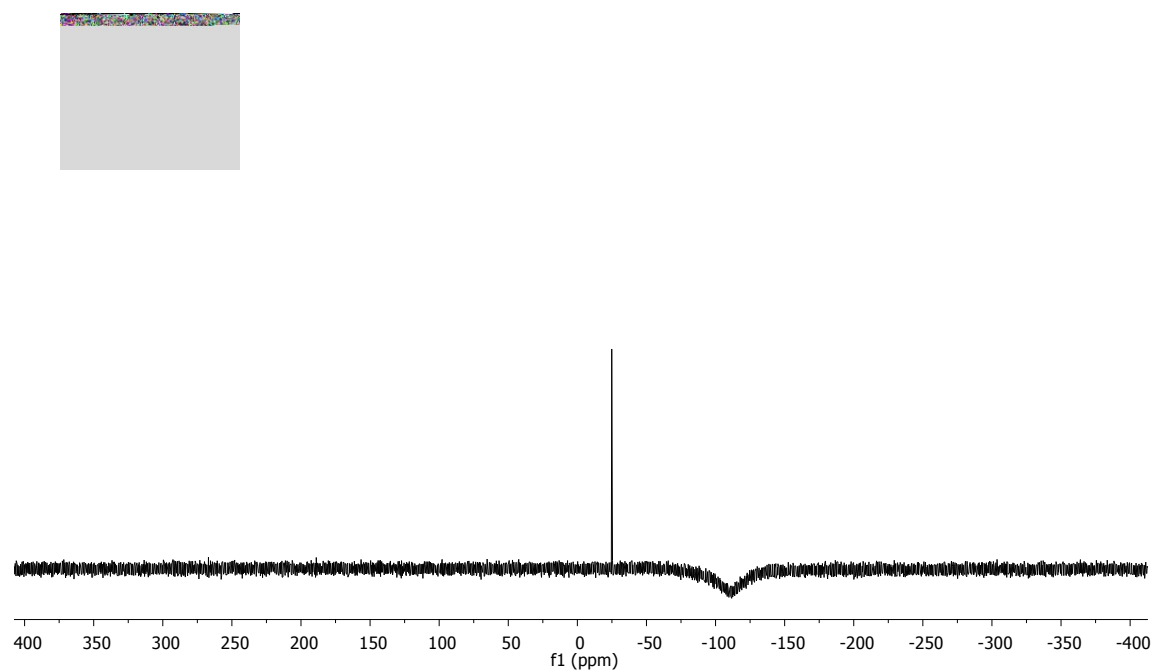
^{29}Si NMR



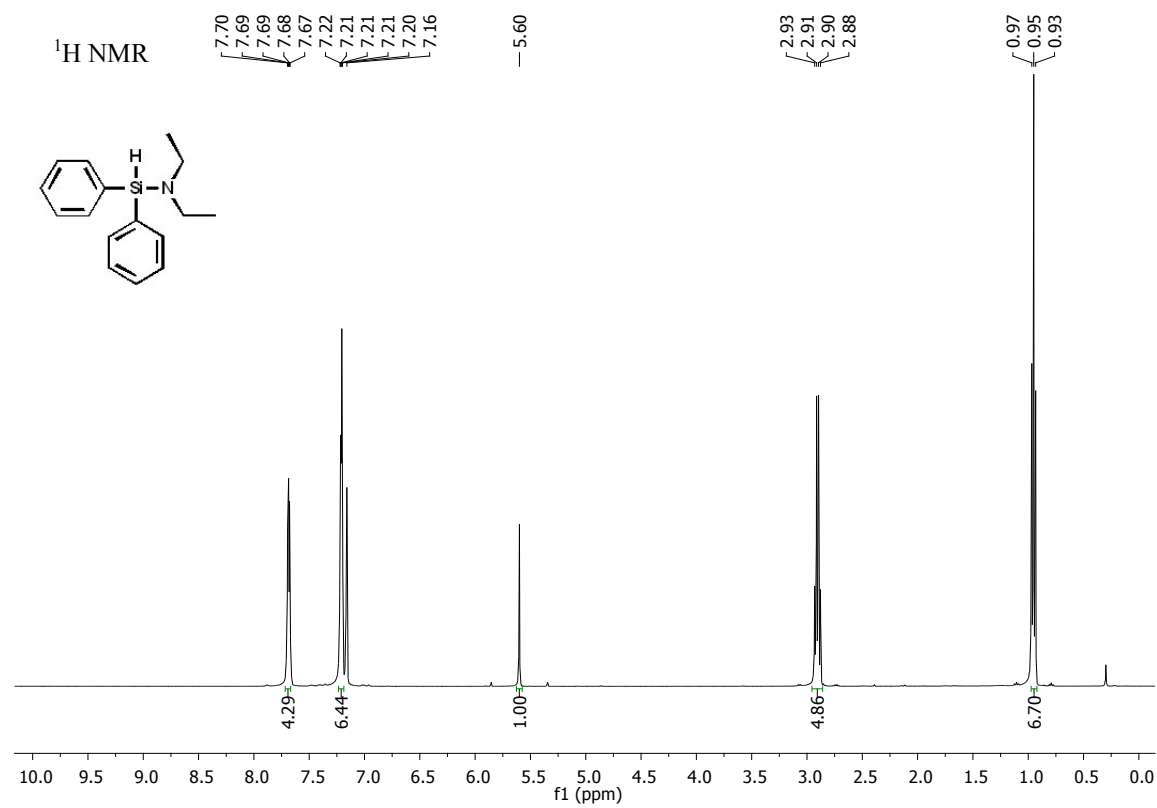


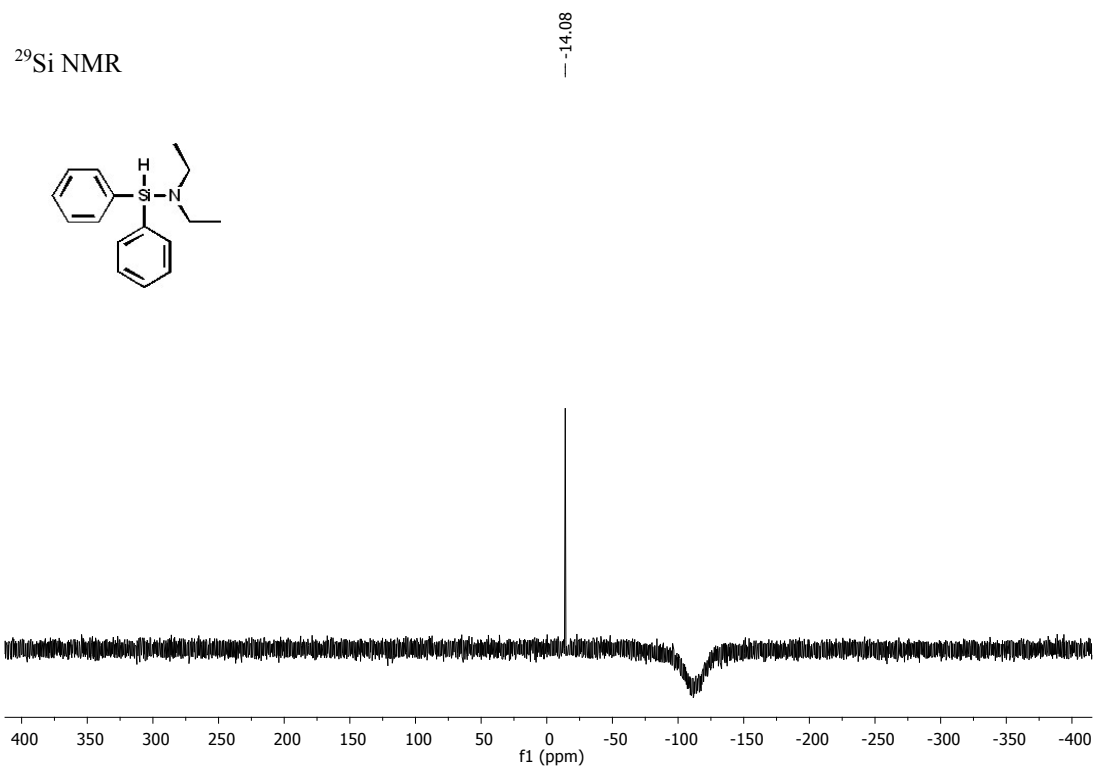
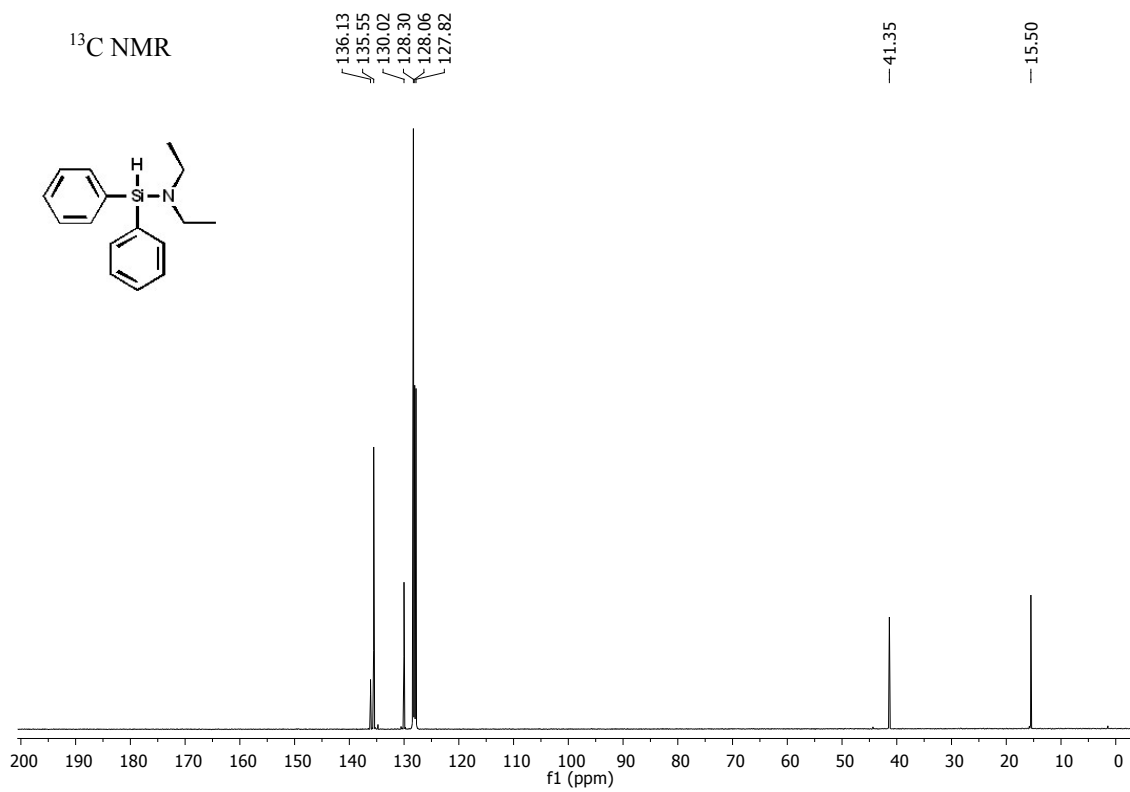
²⁹Si NMR

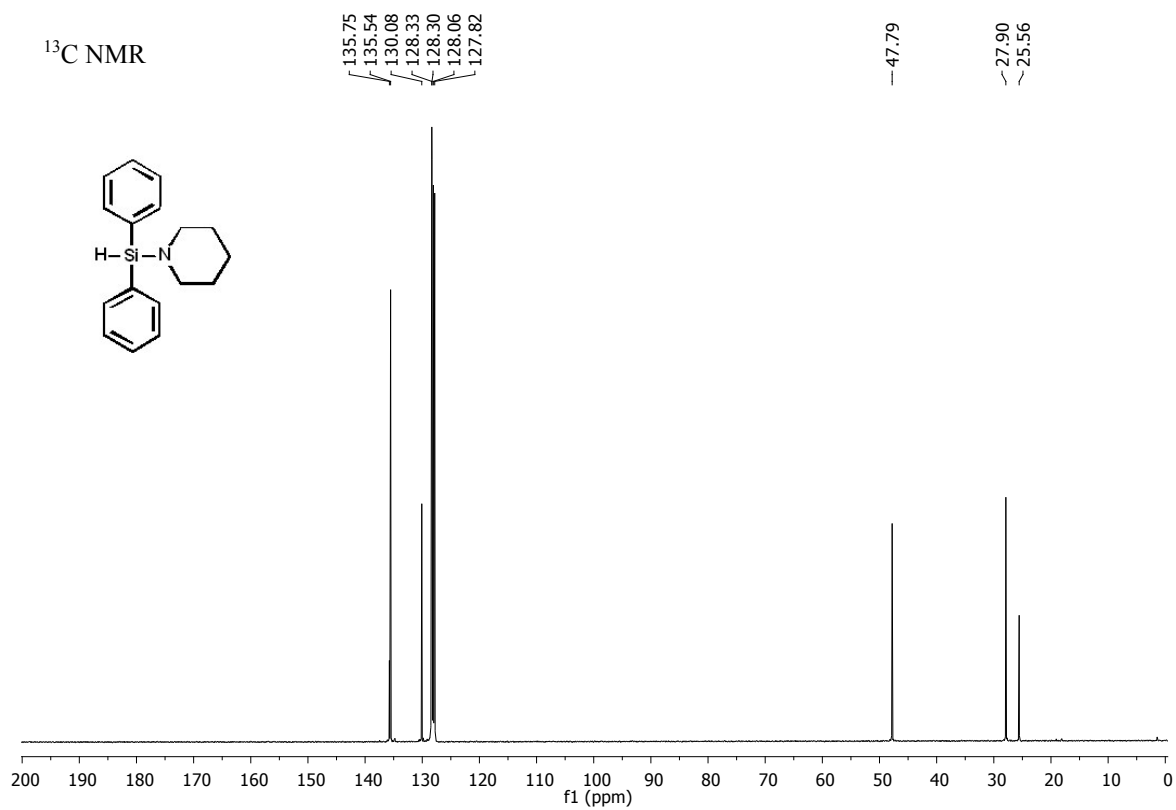
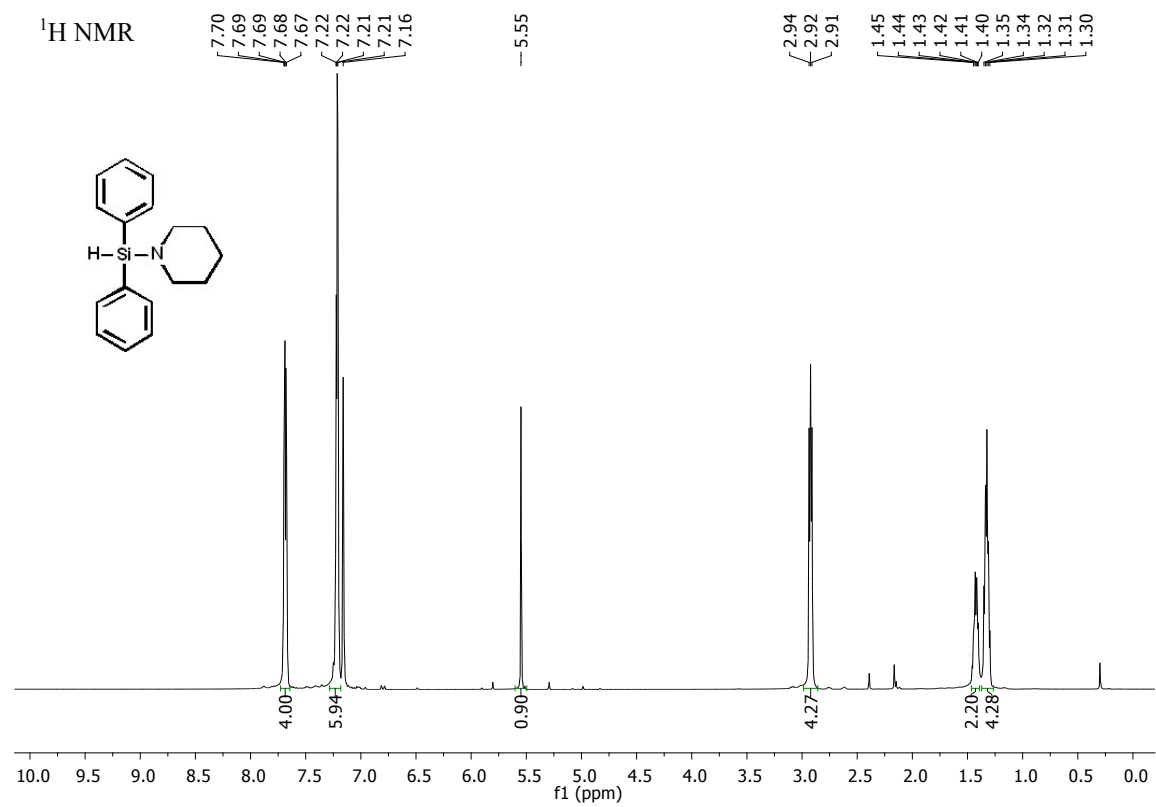
—24.88

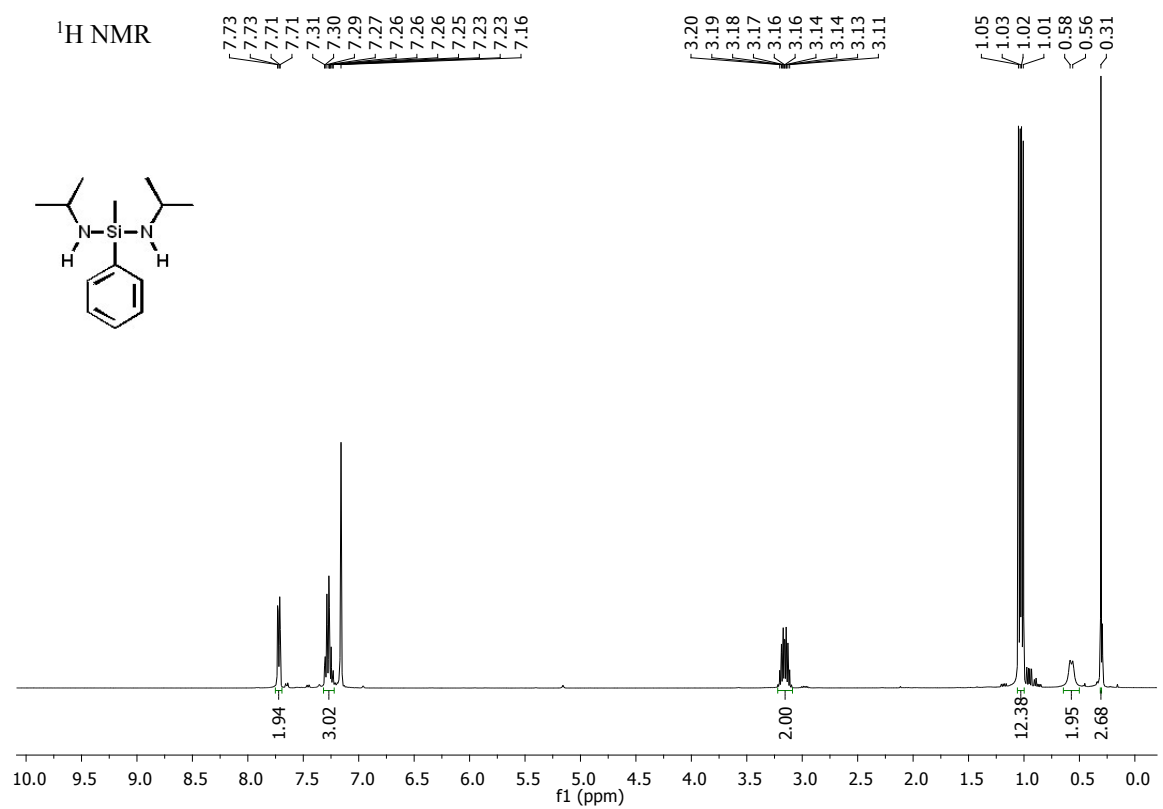
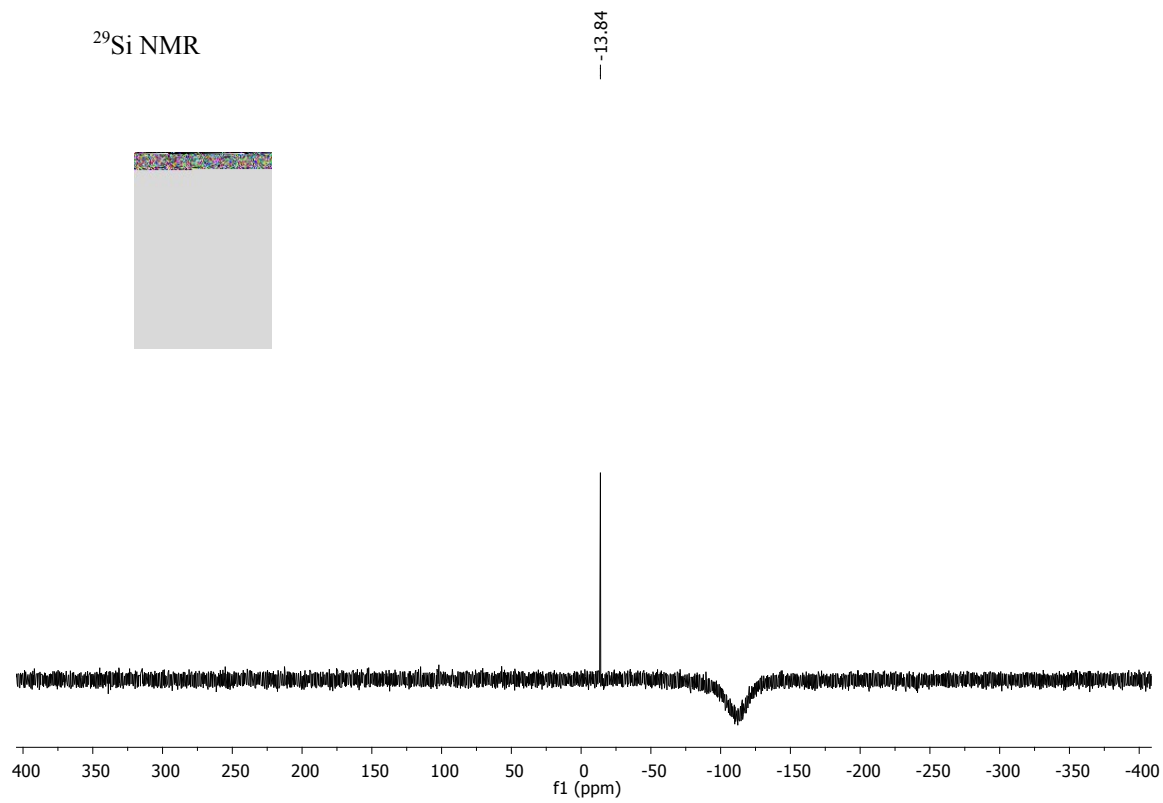


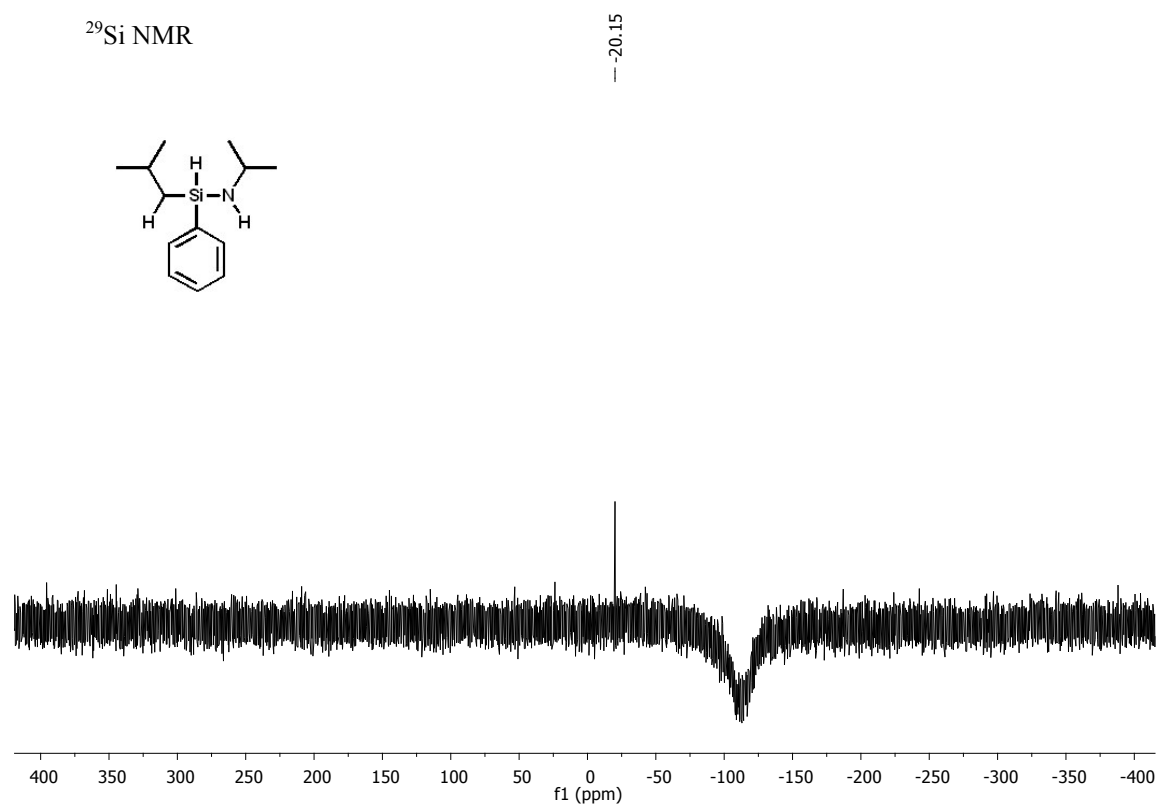
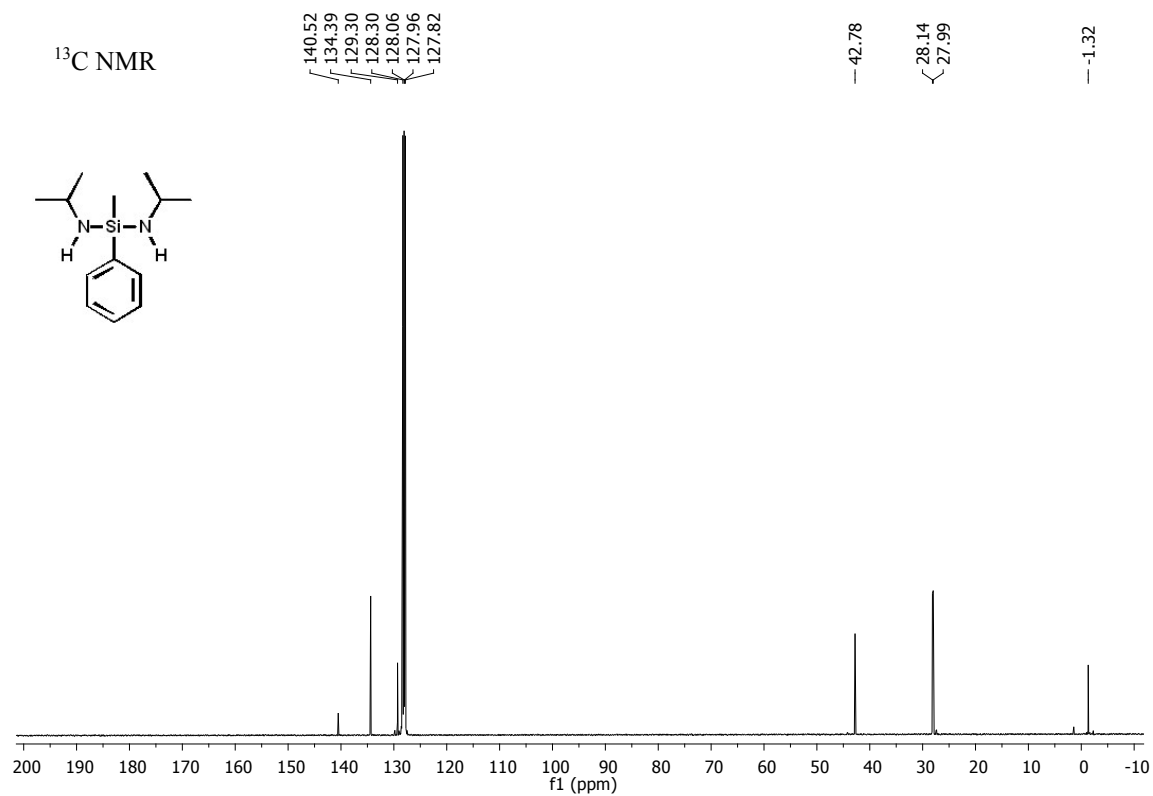
¹H NMR



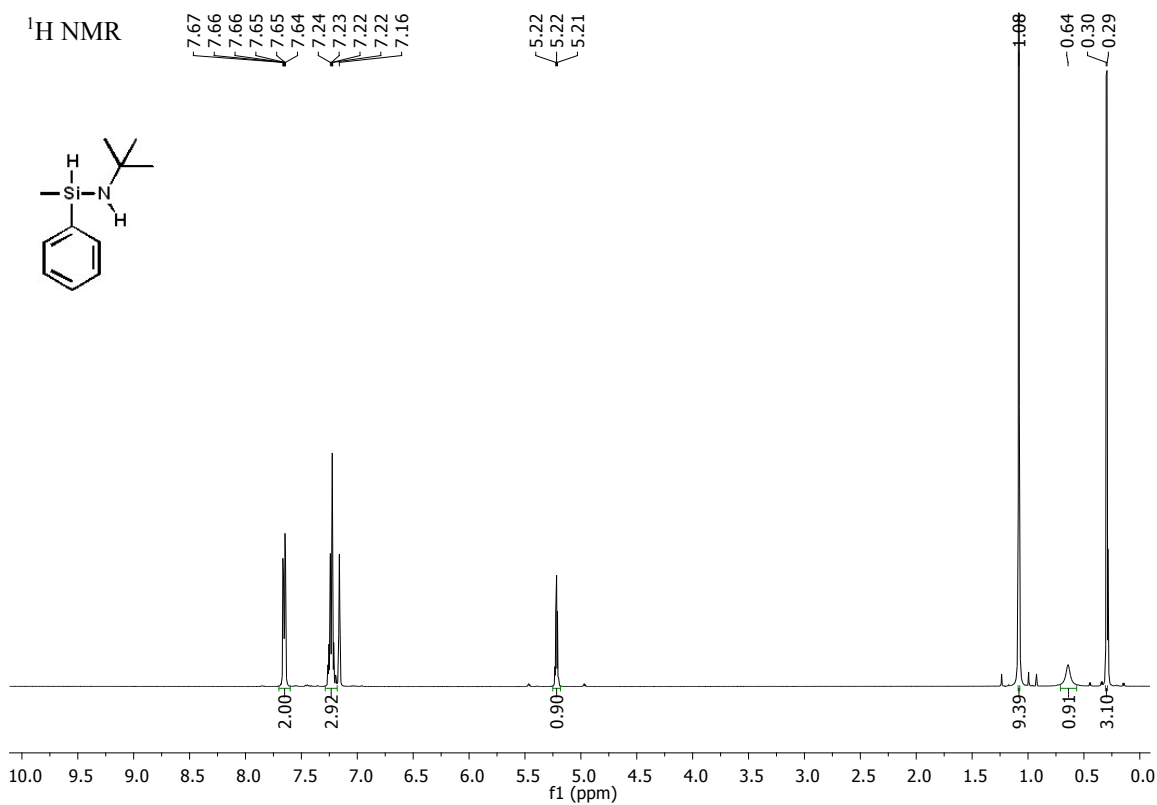




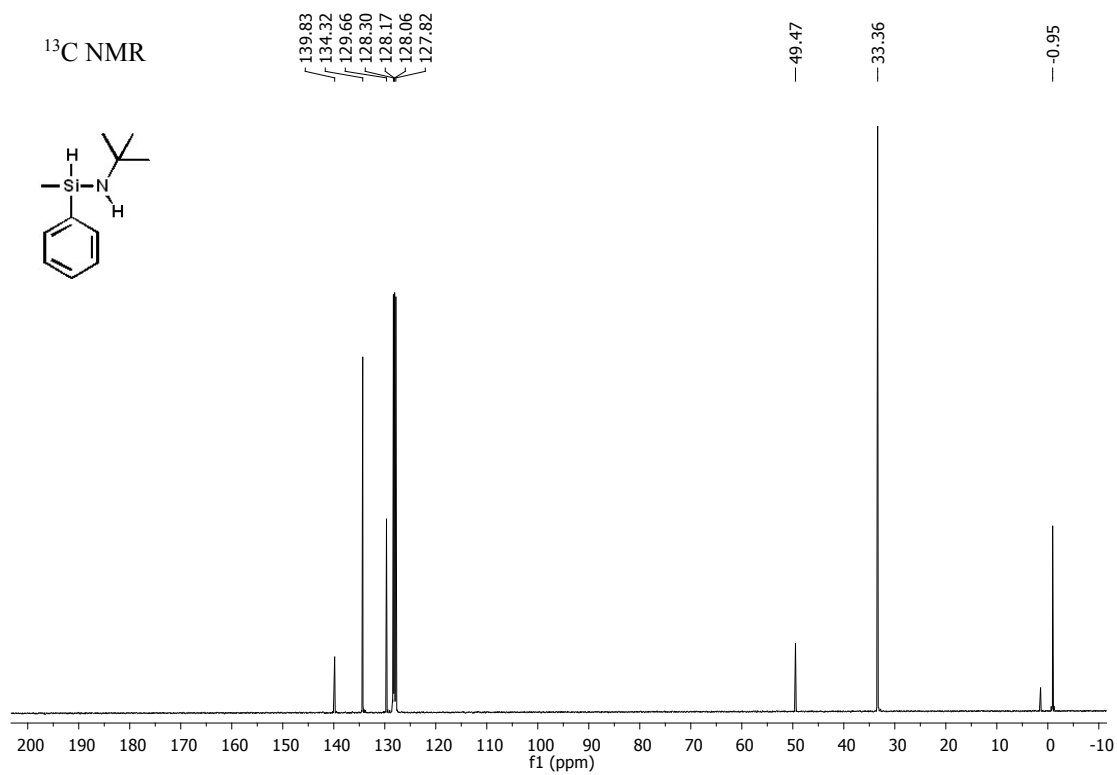




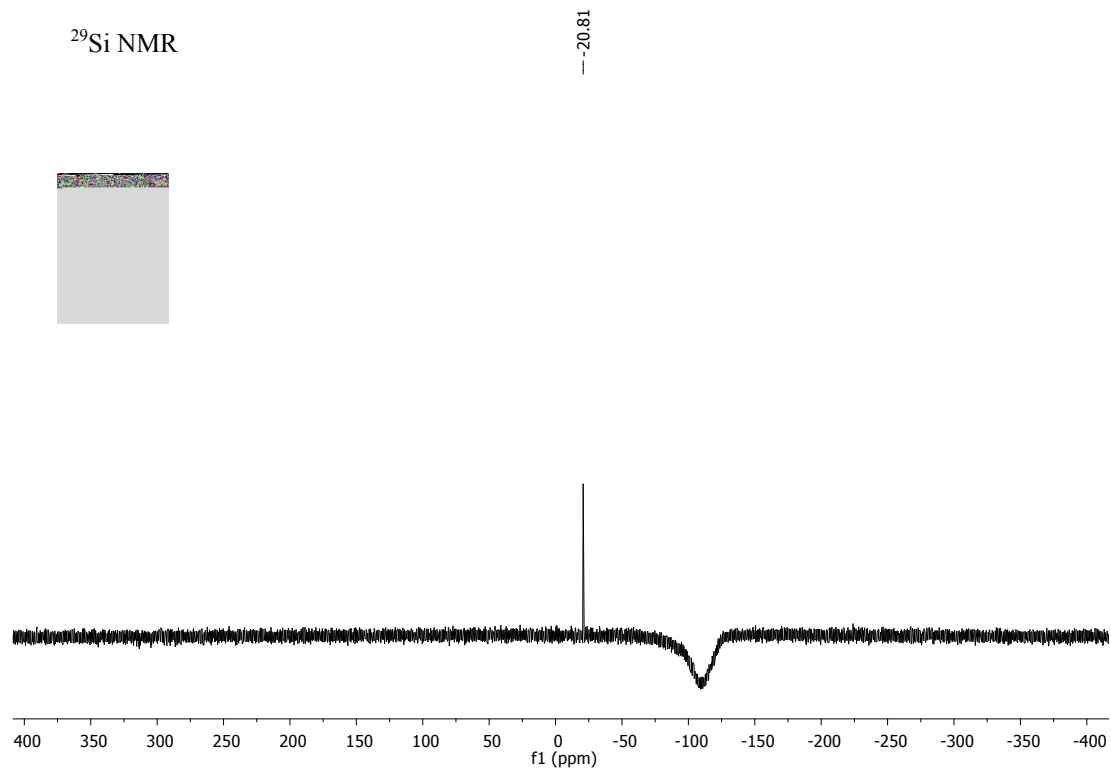
¹H NMR



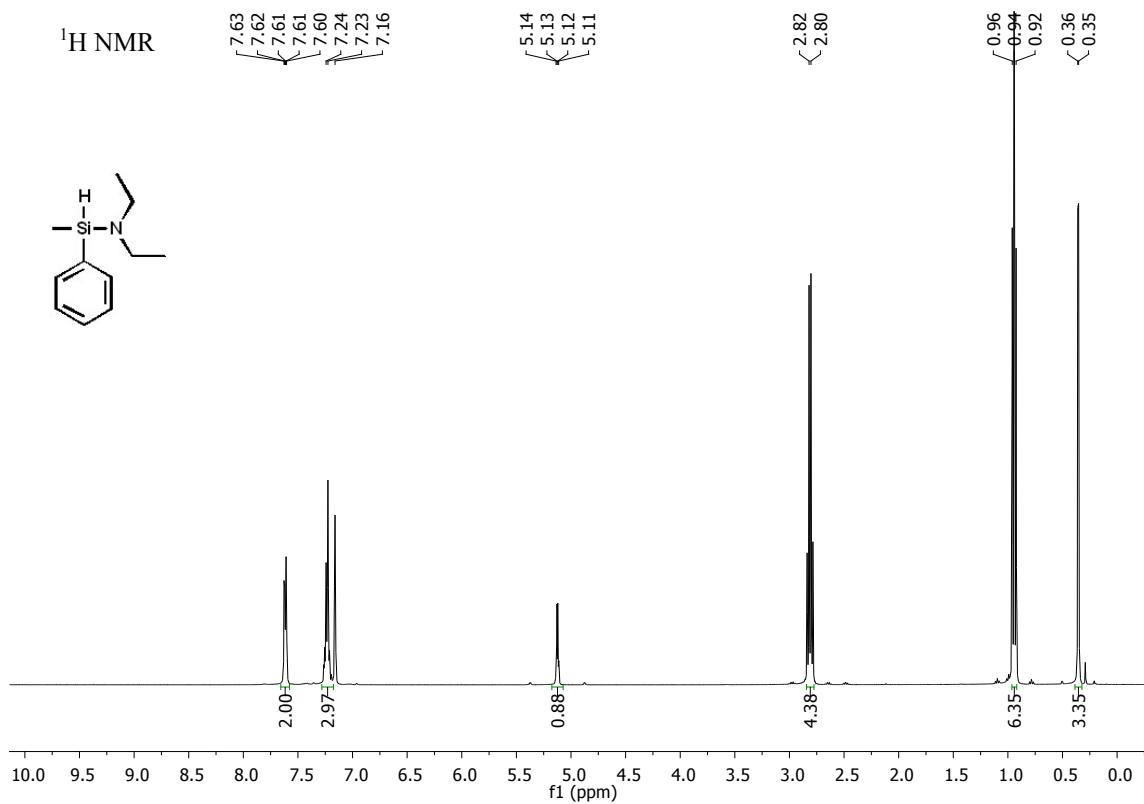
¹³C NMR



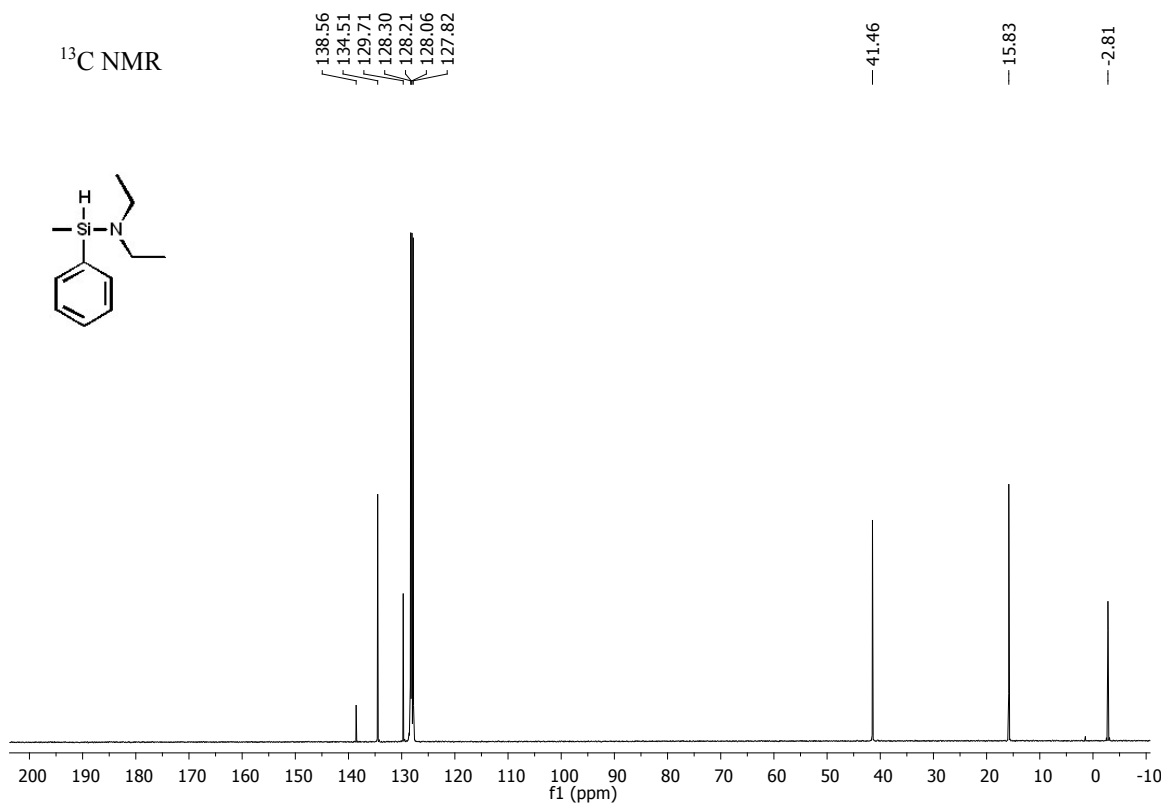
^{29}Si NMR



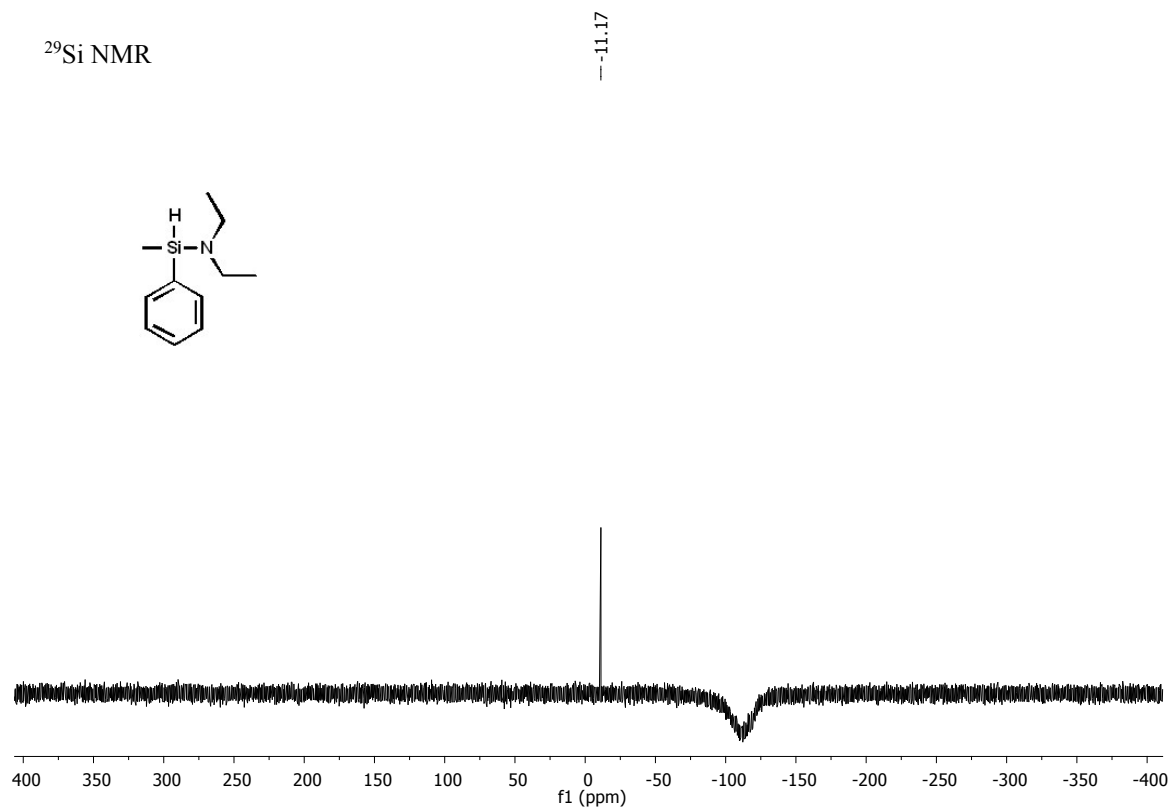
^1H NMR

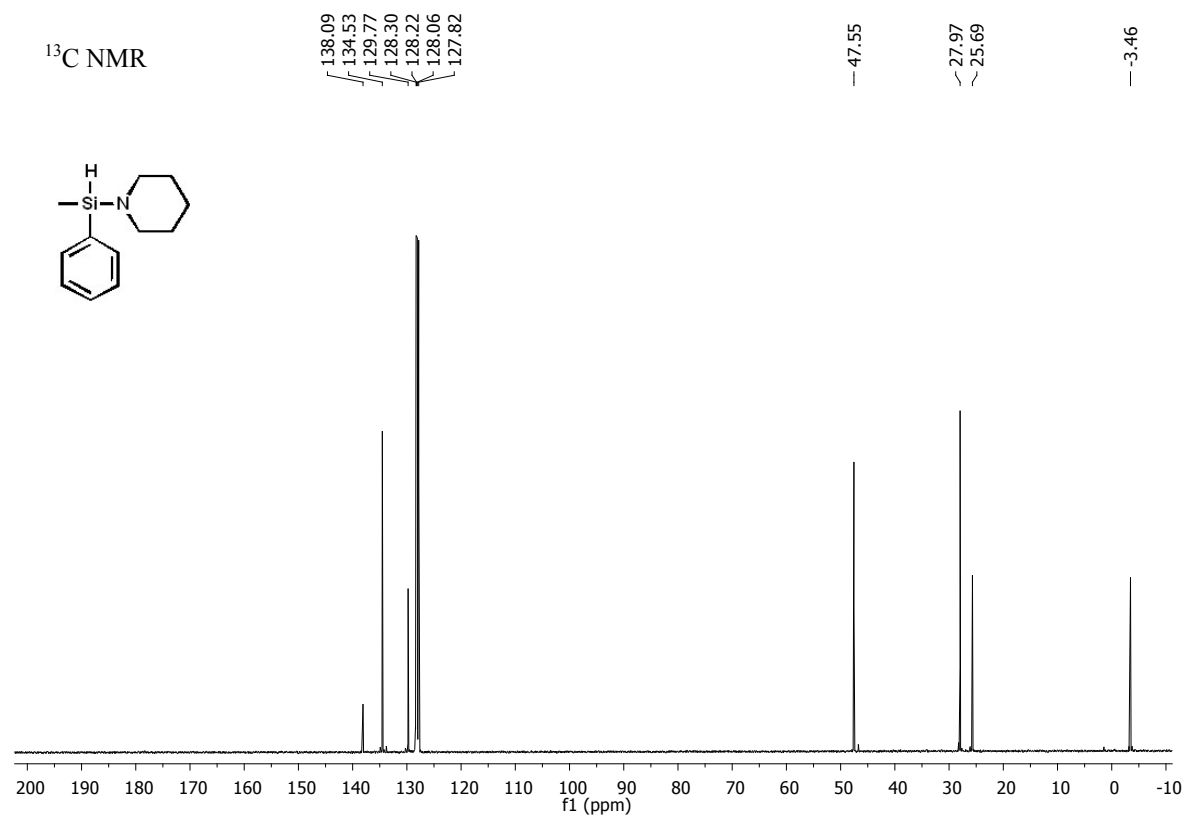
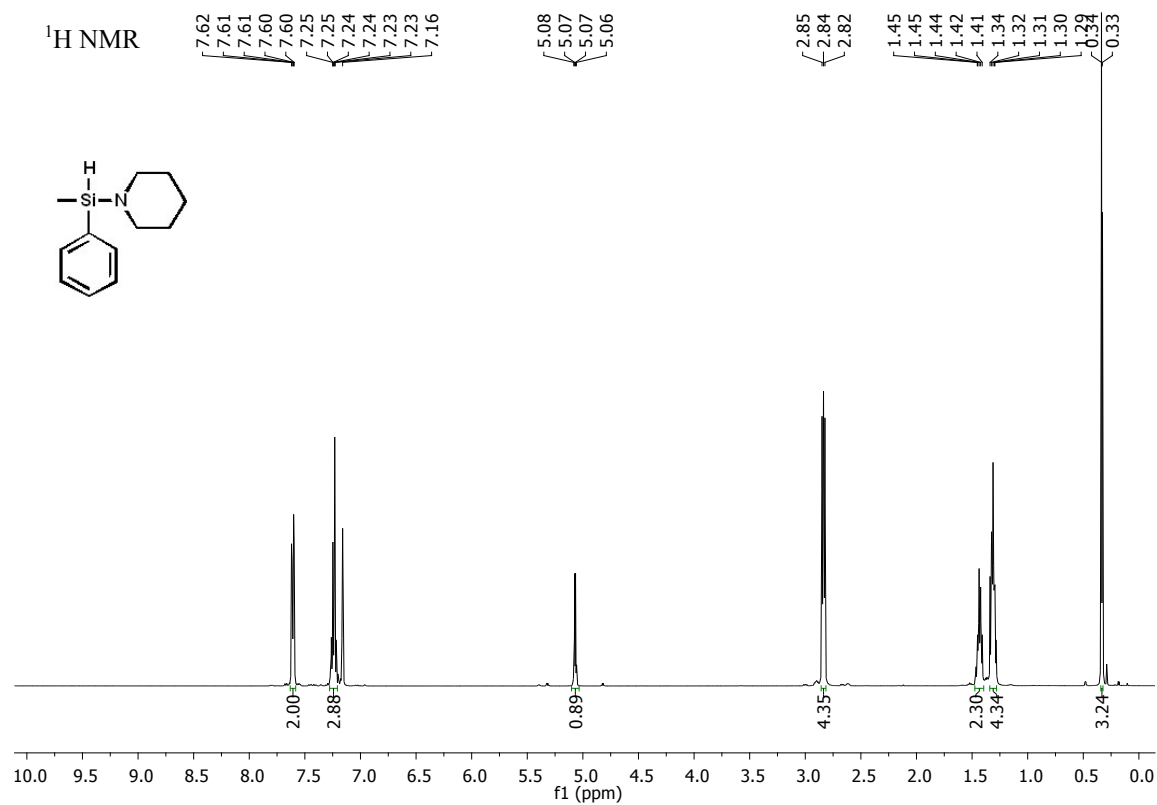


^{13}C NMR



^{29}Si NMR





^{29}Si NMR

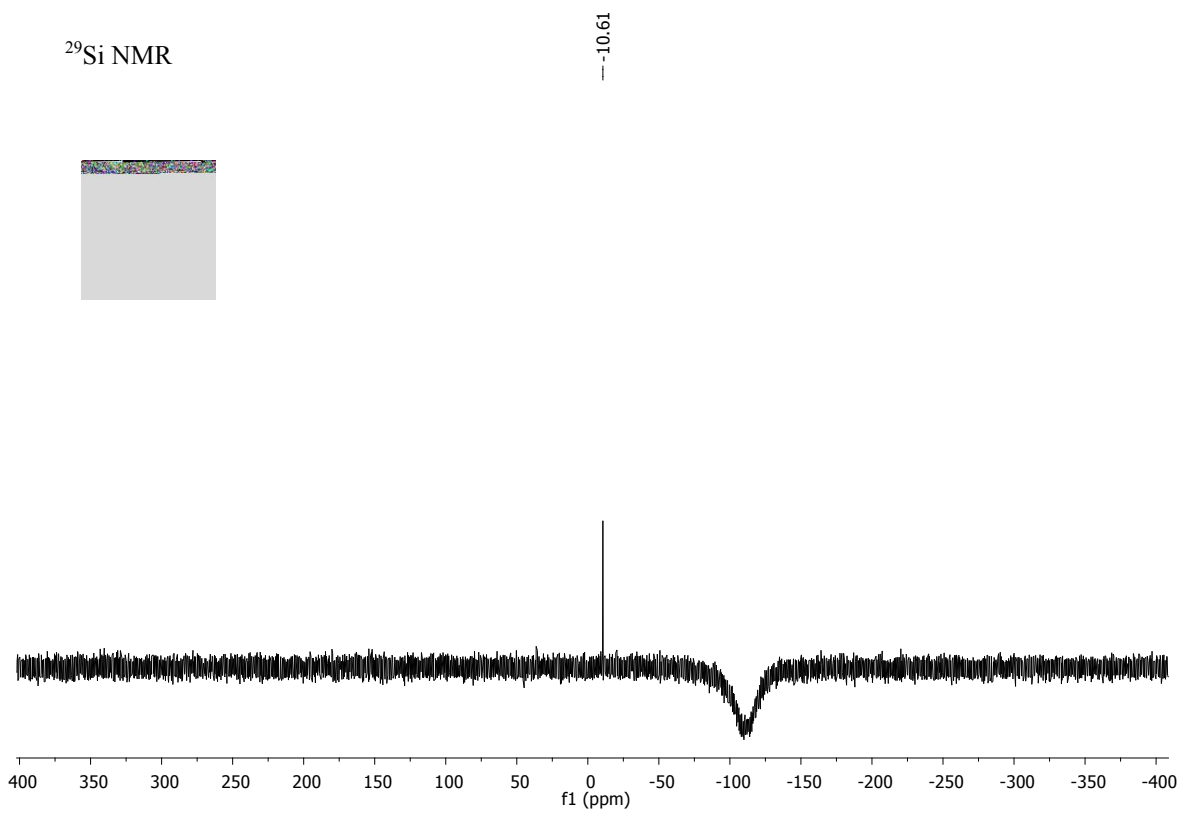


Table S1. Crystal data and structure refinement details of compounds **1** and **2**

	1	2
CCDC	1517518	1517519
Empirical formula	C ₄₂ H ₅₉ MgN ₃ Si ₂	C ₃₆ H ₅₃ N ₃ MgSi ₂
Formula weight	686.41	608.30
Crystal system	Triclinic	Triclinic
Space group	$P\bar{1}$	$P\bar{1}$
a/Å	9.4770(4)	9.1877(3)
b/Å	12.9087(6)	11.2375(3)
c/Å	18.2858(8)	18.5799(5)
α /°	72.347(3)	85.9850(10)
β /°	76.154(3)	77.6490(10)
γ /°	89.454(3)	1830.86(9)
Volume/Å ³	2064.92(16)	1830.86(9)
Z	2	2
ρ_{calc} /cm ³	1.104	1.103
μ /mm ⁻¹	0.132	0.141
F(000)	744.0	660.0
Crystal size /mm ³	0.16 × 0.14 × 0.1	0.19 × 0.14 × 0.1
2 θ range for data collection/°	7.738 to 60.226	4.49 to 61.14
Index ranges	-13 ≤ h ≤ 12, -18 ≤ k ≤ 17, -25 ≤ l ≤ 25	-12 ≤ h ≤ 13, -15 ≤ k ≤ 16, -24 ≤ l ≤ 26
Reflections collected	31544	29800
Independent reflections	11856 [R _{int} = 0.0443, R _{sigma} = 0.0567]	10999 [R _{int} = 0.0416, R _{sigma} = 0.0518]
Data/restraints/parameters	11856/369/576	10999/0/394
Goodness-of-fit on F ²	1.032	1.040
R1, wR2 [I ≥ 2 σ (I)]	R ₁ = 0.0500,	R ₁ = 0.0505, wR ₂ = 0.1344

	$wR_2 = 0.1242$	
R1, wR2 [all data]	$R_1 = 0.0836, wR_2 = 0.1436$	$R_1 = 0.0736, wR_2 = 0.1521$
Largest diff. peak/hole/eÅ ⁻³	0.39/-0.51	0.73/-0.38