

Supporting Information for

***Synthesis and characterization of functionalized
molecular and macromolecular double-decker
silsesquioxane systems***

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E-mail: Bogdan.Marciniak@amu.edu.pl

Copies of ¹H, ¹³C, ²⁹Si NMR spectra for all compounds

... S -2-

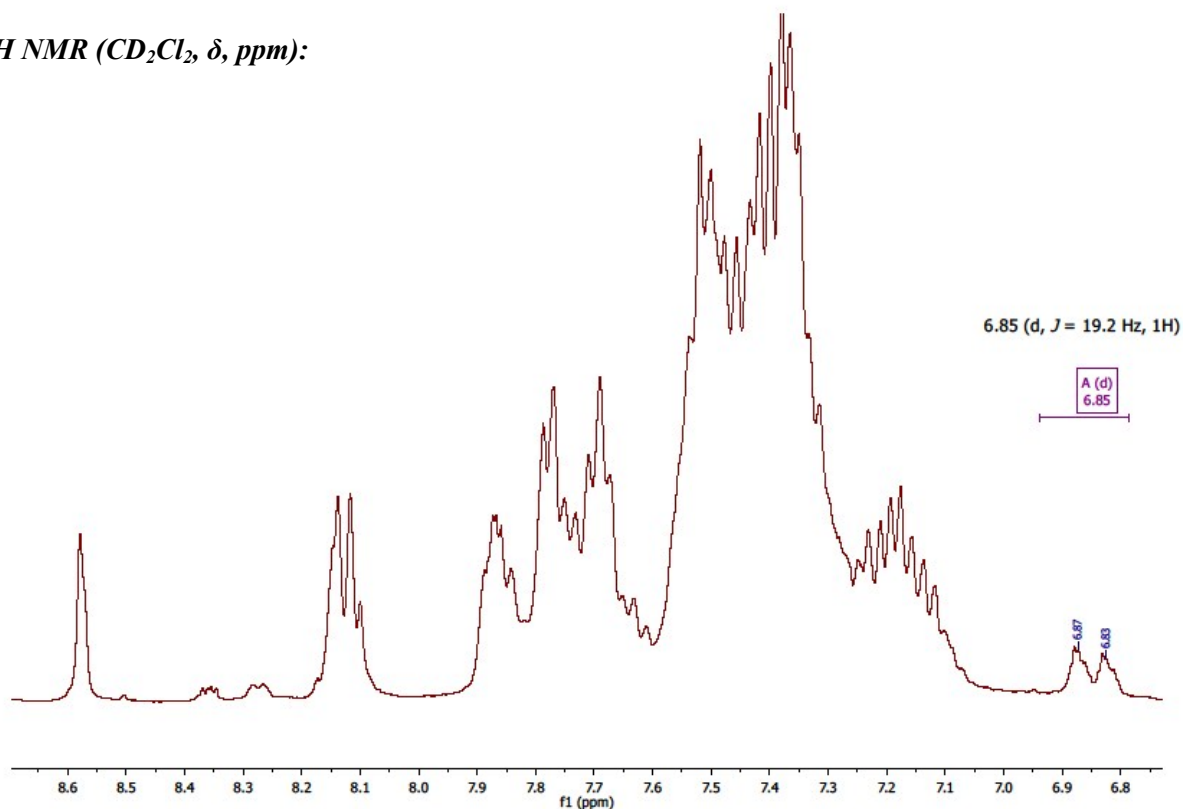
GPC Calibration Curve

... S -24-

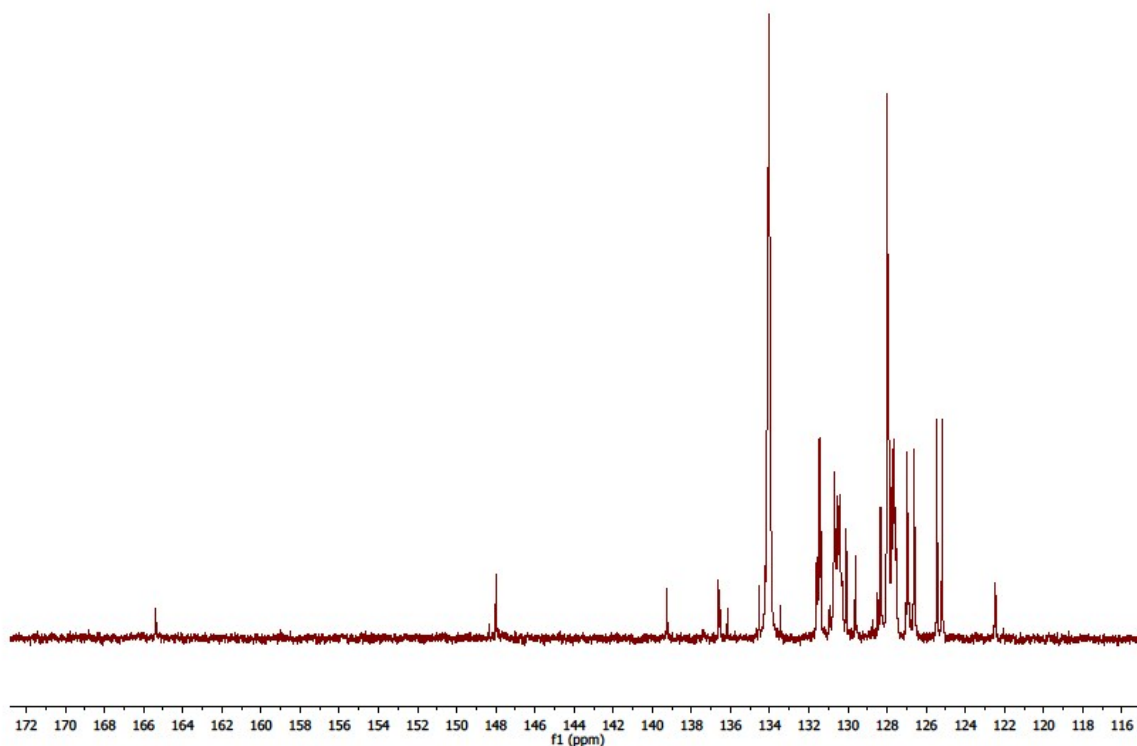
^1H , ^{13}C , ^{29}Si NMR spectra of the isolated molecular compounds (1-8). The spectra are presented in selected regions due to their relevance and clarity according to chemical structure.

1: *trans di[9,19-(E)-4-(9-anthracenyl)styrylphenyl]-1,3,5,7,11,13,15,17-octaphenylpentacyclo-[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane*

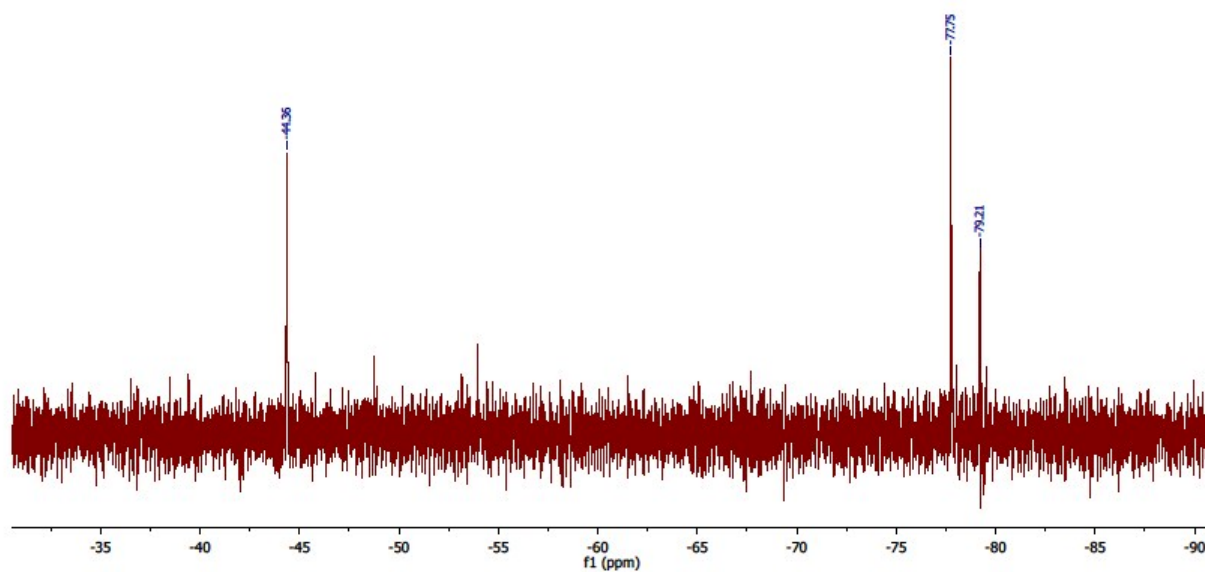
^1H NMR (CD_2Cl_2 , δ , ppm):



^{13}C NMR (CD_2Cl_2 , δ , ppm):

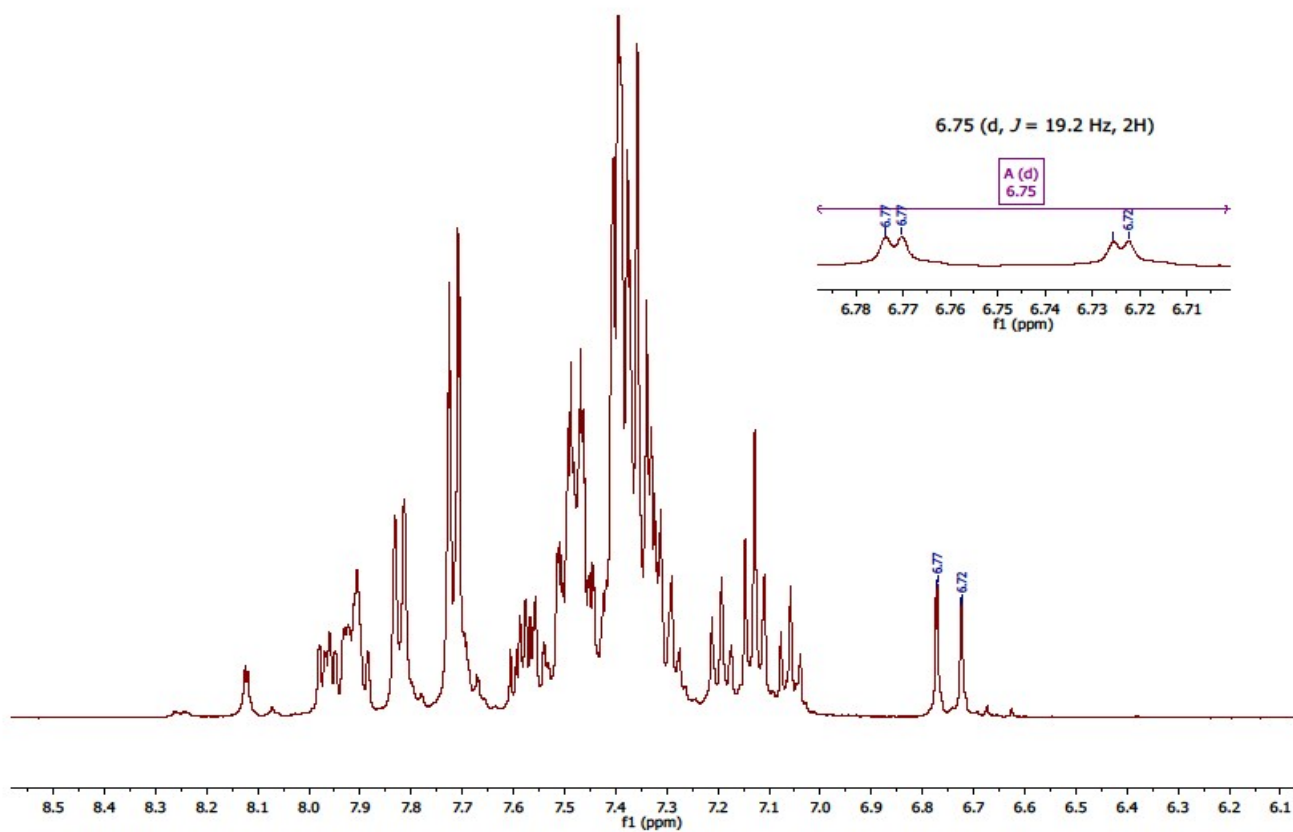


^{29}Si NMR (CD_2Cl_2 , δ , ppm):

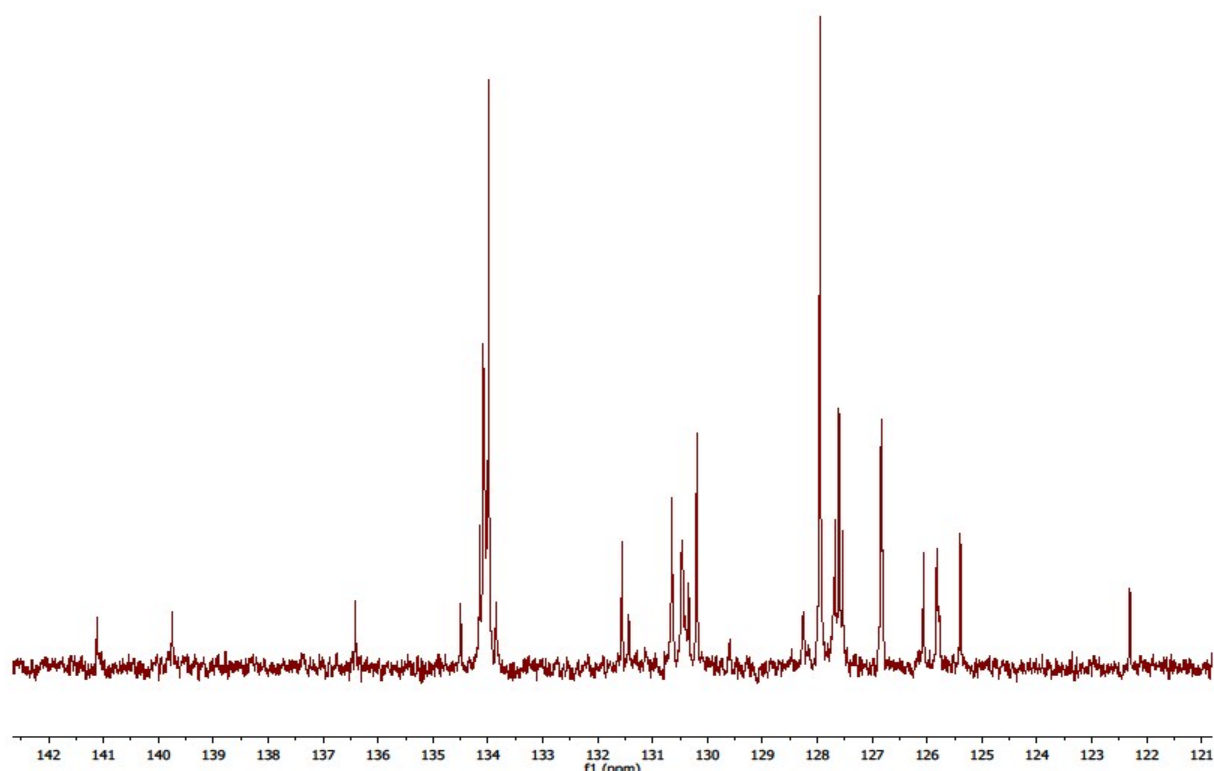


2: *trans* di[9,19-(*E*)-4-(1-naphthyl)styrylphenyl]-1,3,5,7,11,13,15,17-octaphenylpentacyclo-[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane

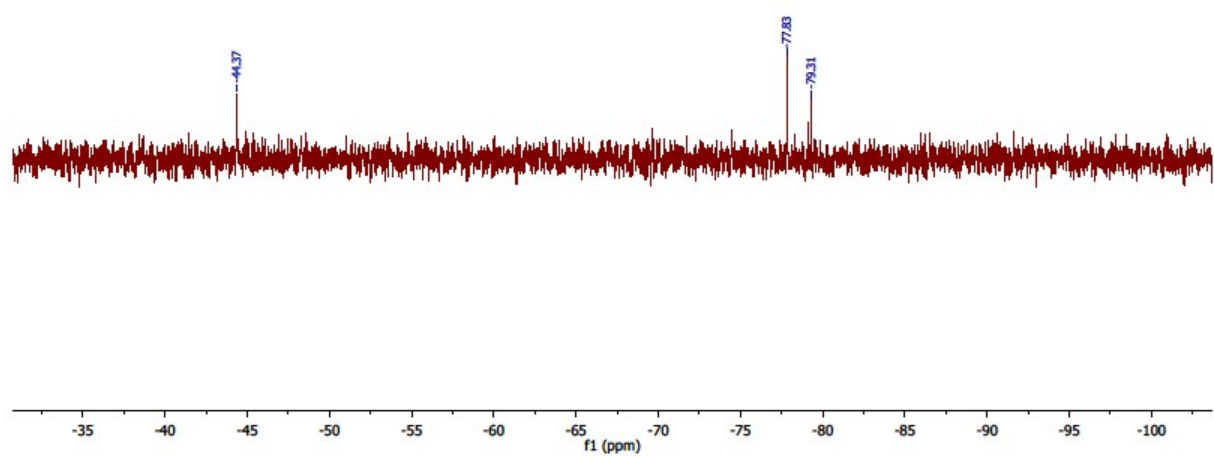
^1H NMR (CD_2Cl_2 , δ , ppm):



^{13}C NMR (CD_2Cl_2 , δ , ppm):

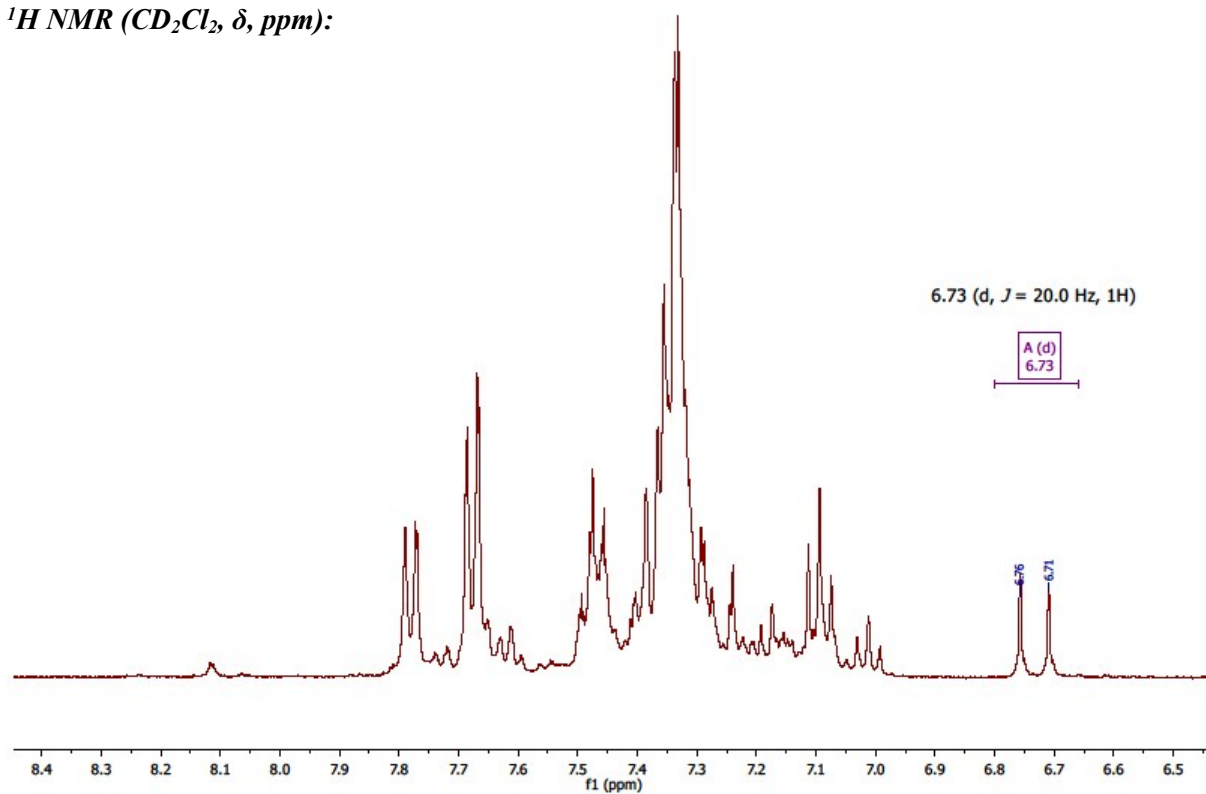


^{29}Si NMR (CD_2Cl_2 , δ , ppm):

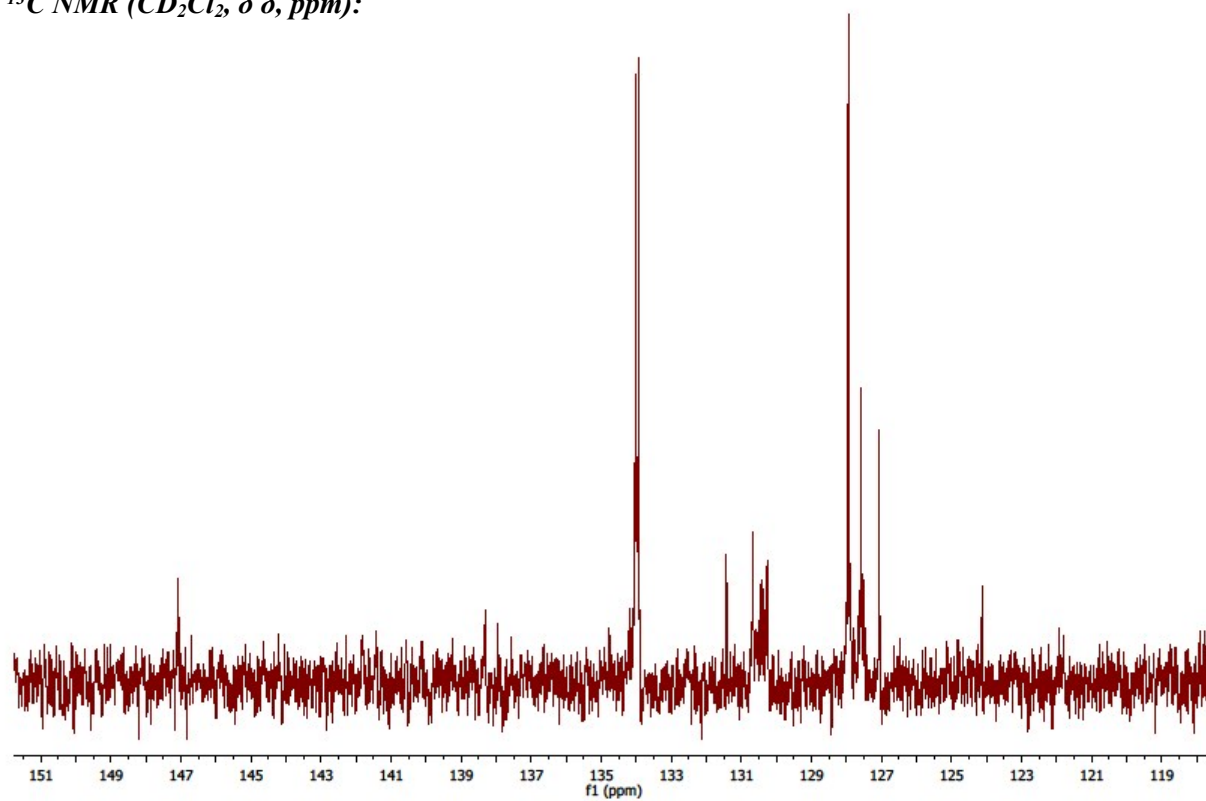


3: *trans* di[9,19-(*E*)-4-(2,3,4,5,6-pentafluorophenyl)styryl-phenyl]-1,3,5,7,11,13,15,17-octaphenyl-pentacyclo[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane:

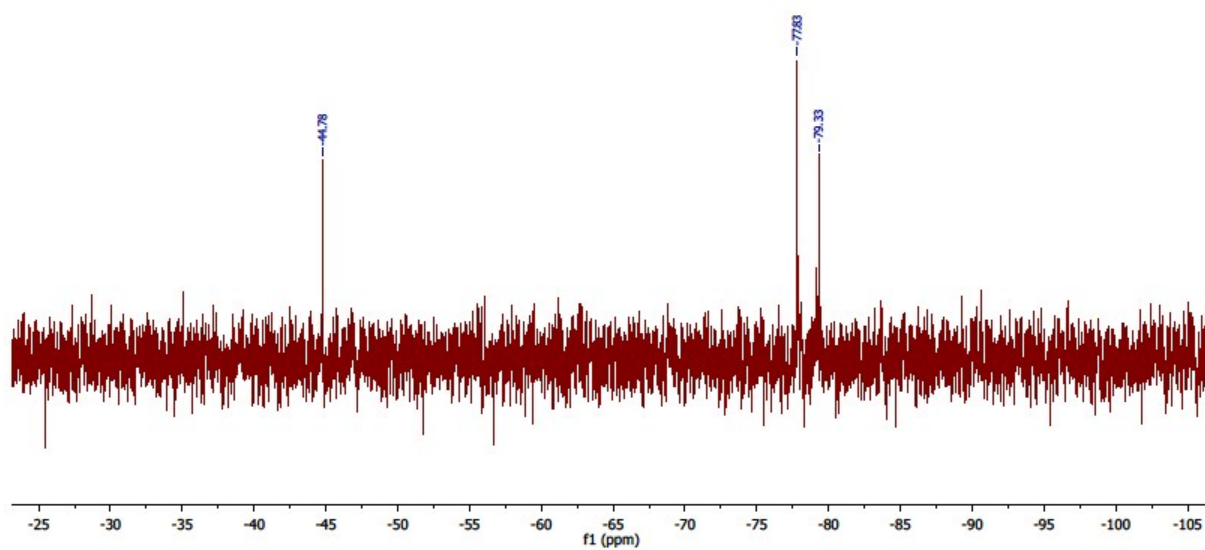
¹H NMR (CD₂Cl₂, δ, ppm):



¹³C NMR (CD₂Cl₂, δ, ppm):

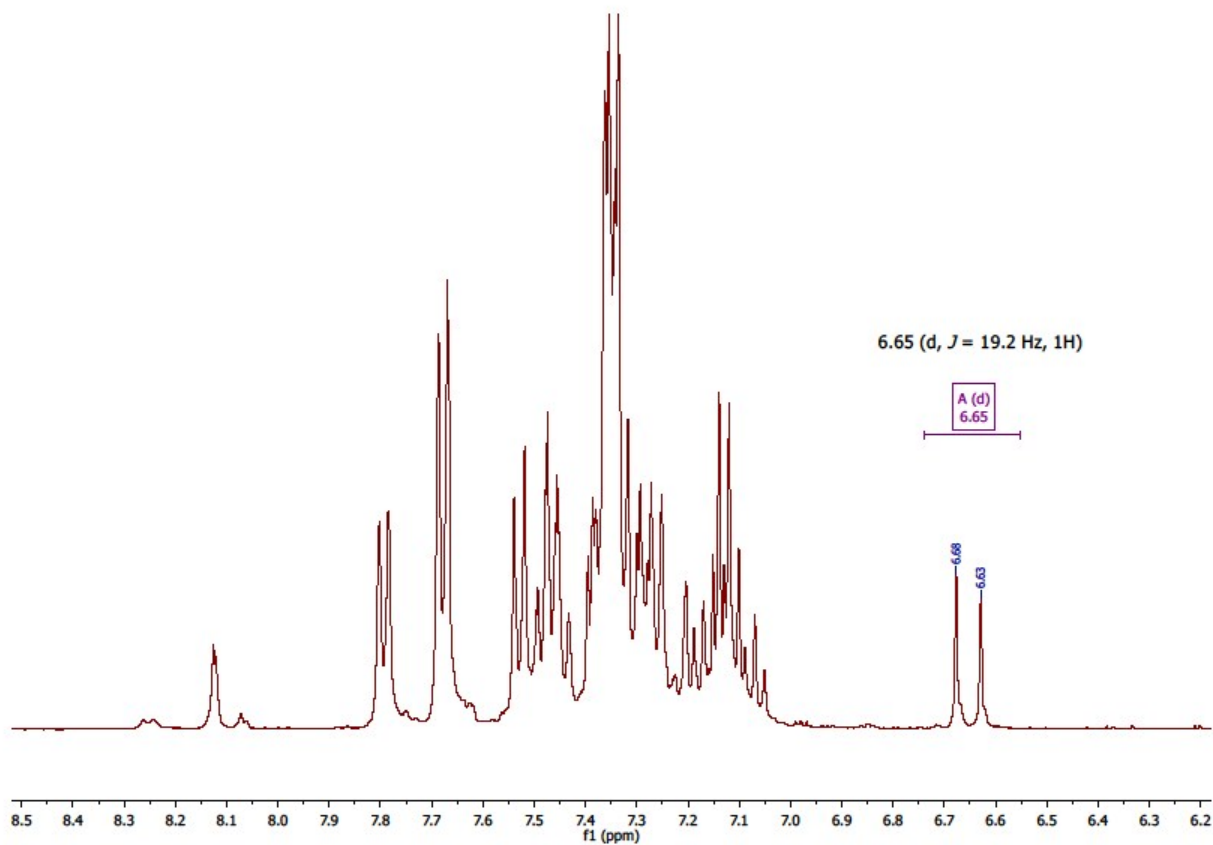


^{29}Si NMR (CD_2Cl_2 , δ , ppm):

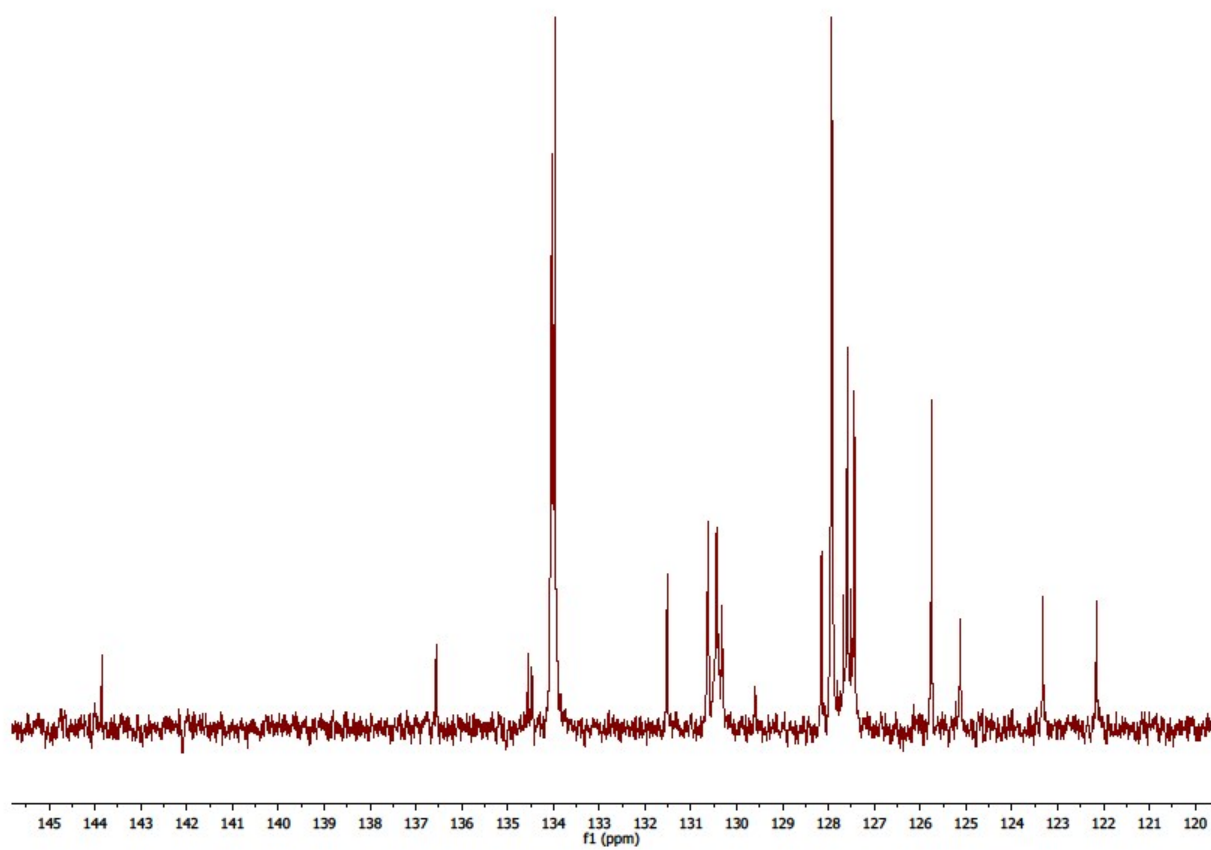


4: *trans* di[9,19-(*E*)-4-(2-thienyl)styryl]phenyl]-1,3,5,7,11,13,15,17octaphenylpentacyclo-[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane

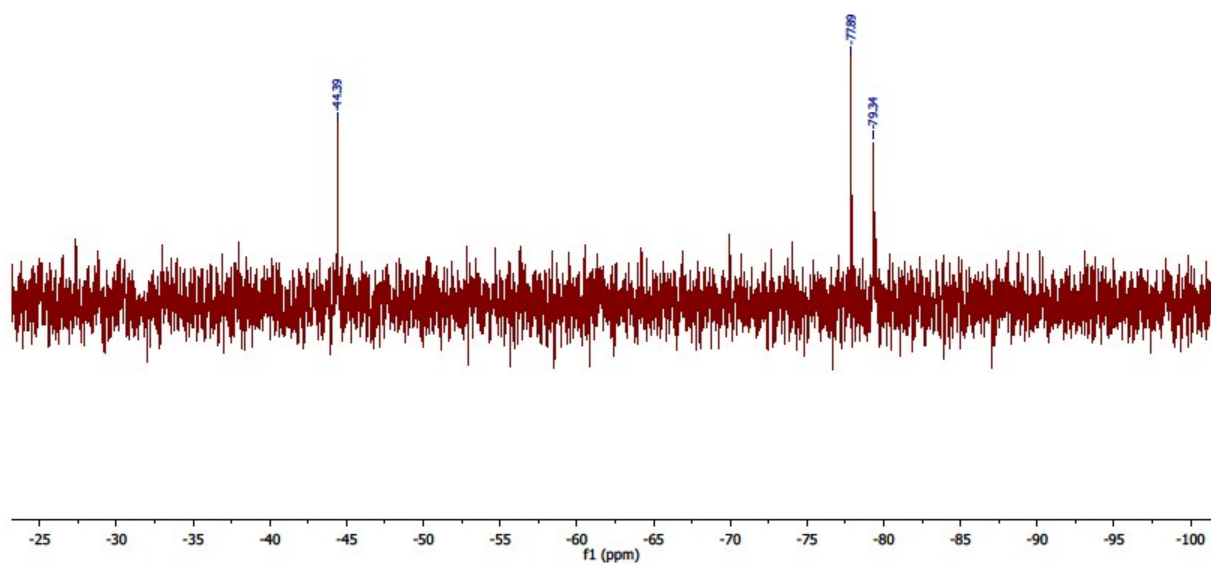
^1H NMR (CD_2Cl_2 , δ , ppm):



^{13}C NMR (CD_2Cl_2 , δ , ppm):

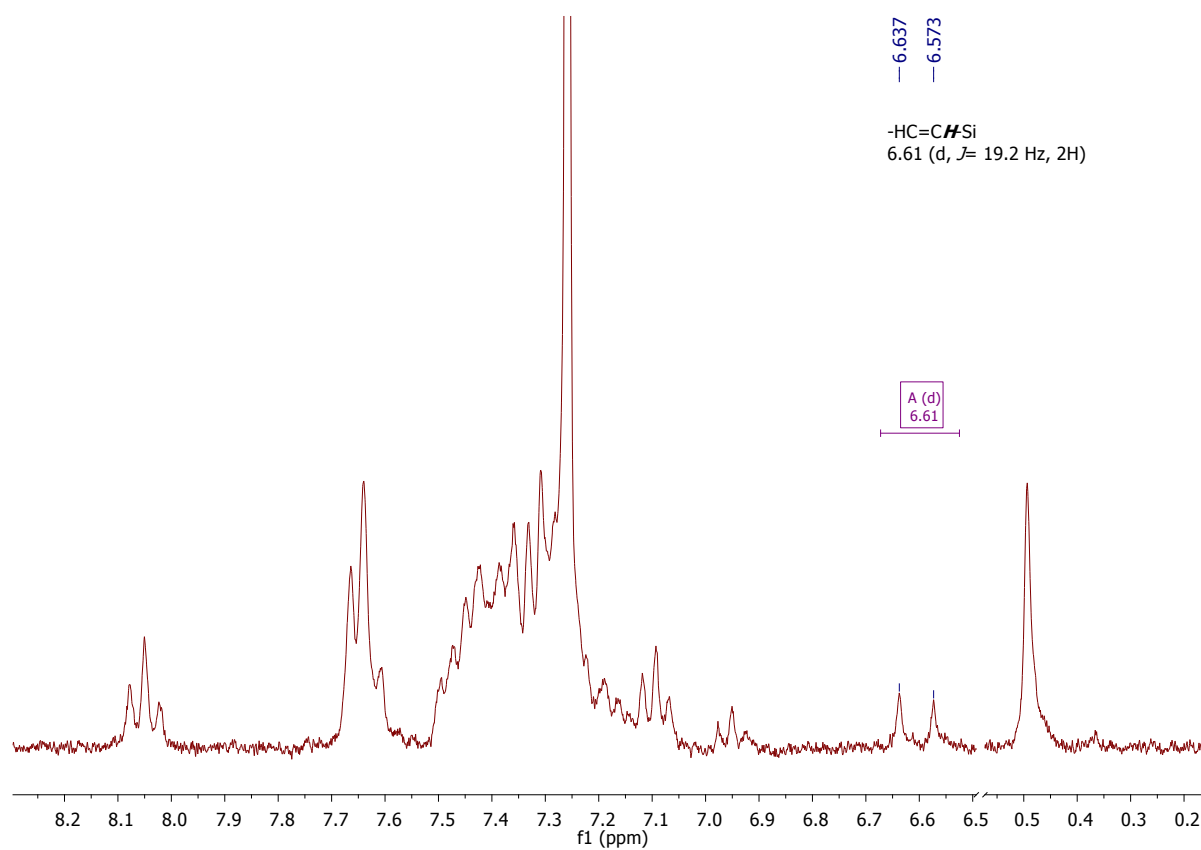


^{29}Si NMR (CD_2Cl_2 , δ , ppm):

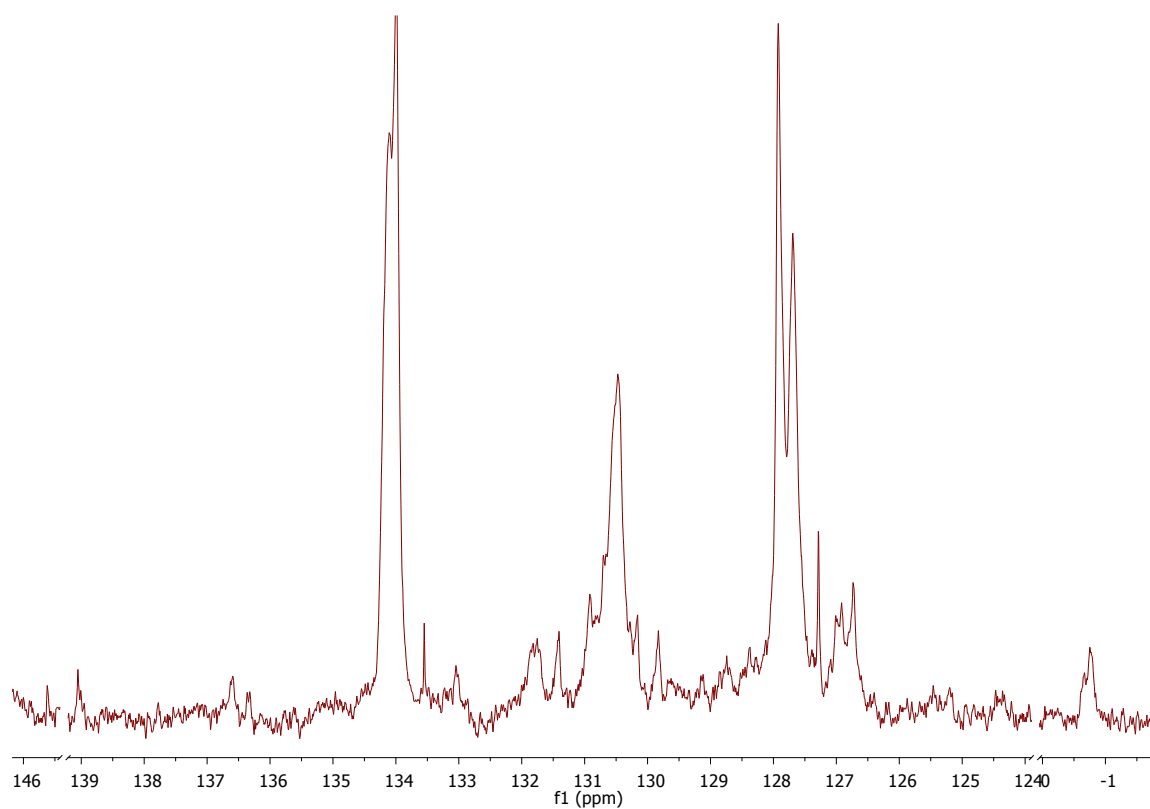


5: *cis* and *trans* mixture of di[9,19-(*E*)-4-(9-anthracenyl)styryl-methyl]-1,3,5,7,11,13,15,17-octaphenylpentacyclo [11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane

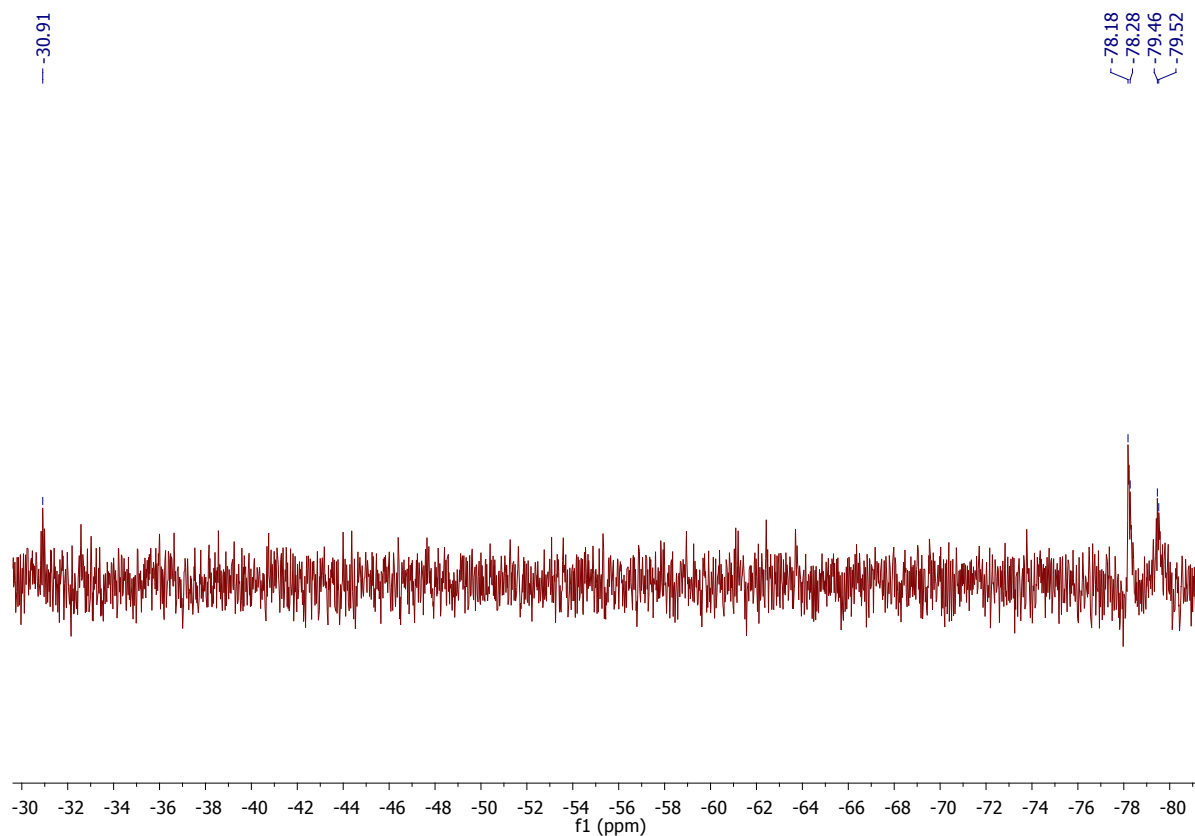
¹H NMR (CDCl₃, δ, ppm):



¹³C NMR (CDCl₃, δ, ppm):

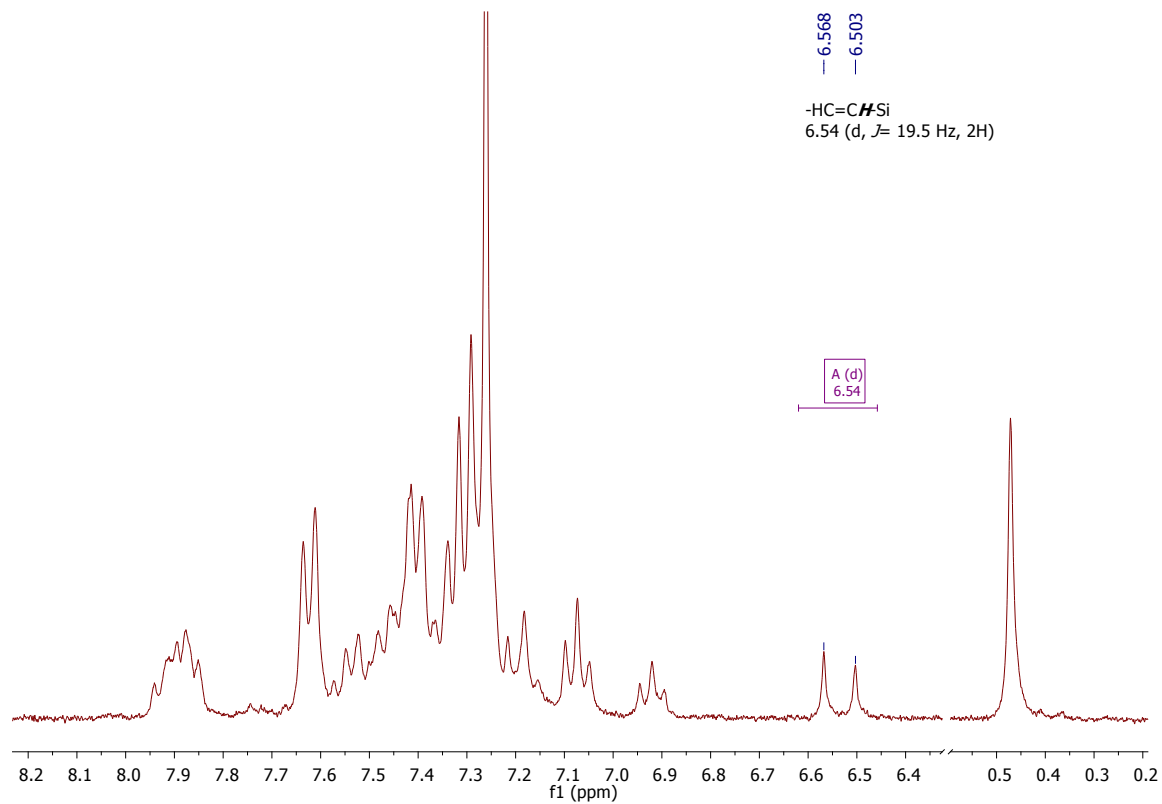


^{29}Si NMR (CDCl_3 , δ , ppm):

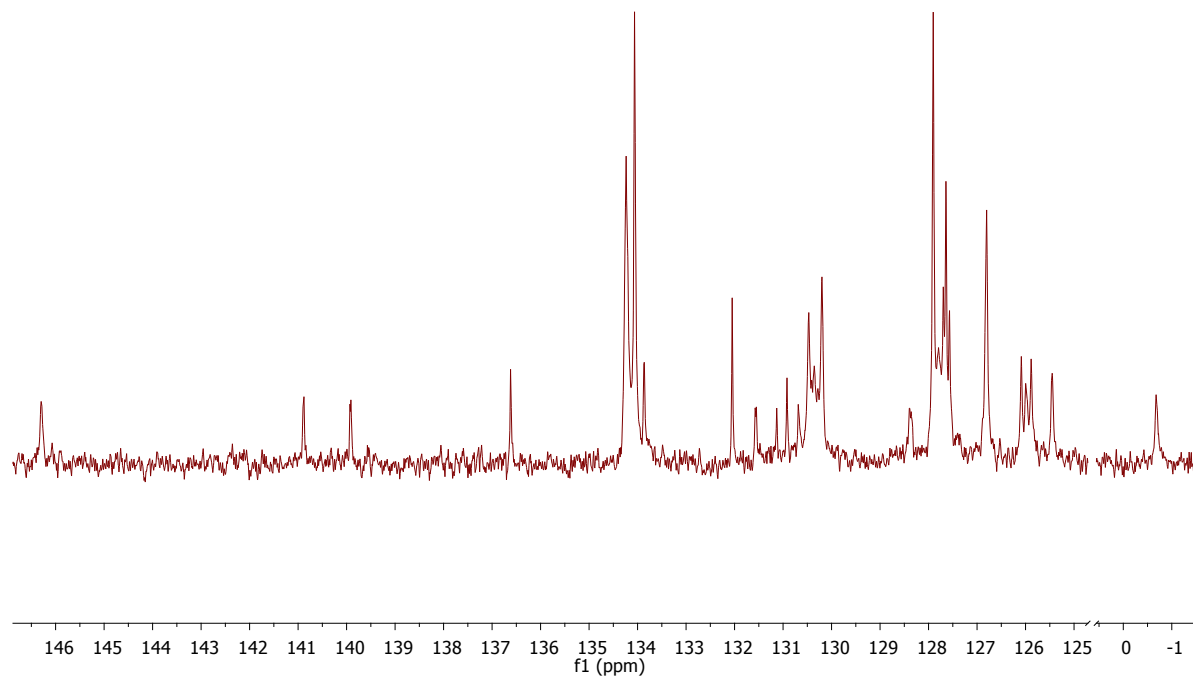


6: *cis* and *trans* mixture of di[9,19-(*E*)-4-(1-naphthyl)styrylmethyl]-1,3,5,7,11,13,15,17-octaphenylpentacyclo[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane

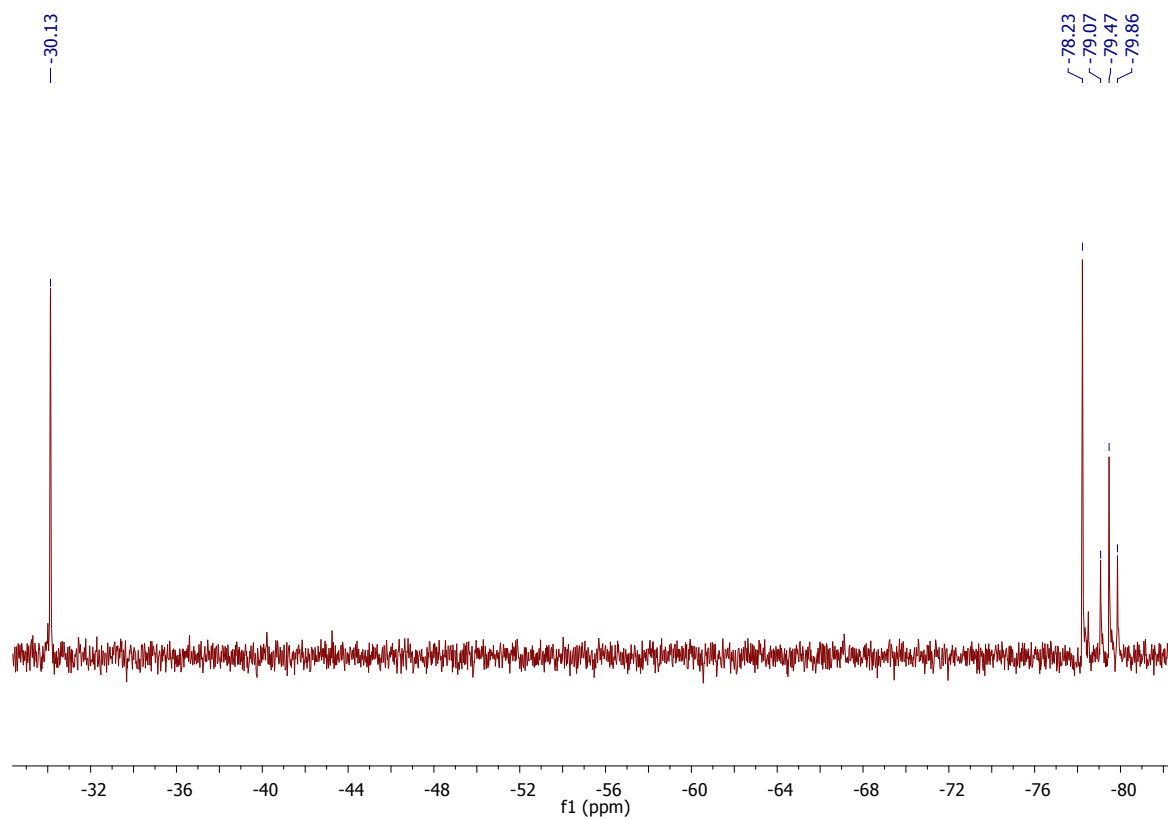
^1H NMR (CDCl_3 , δ , ppm):



^{13}C NMR (CDCl_3 , δ , ppm):

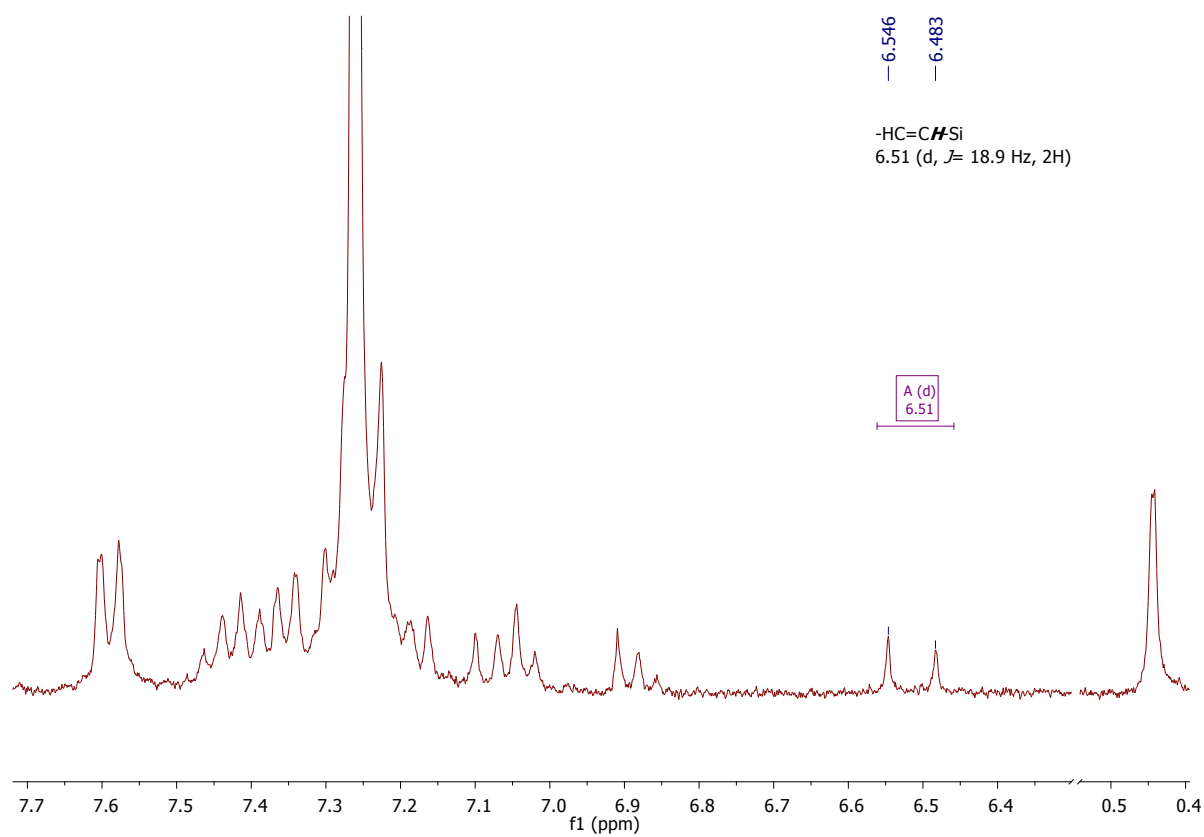


^{29}Si NMR (CDCl_3 , δ , ppm):

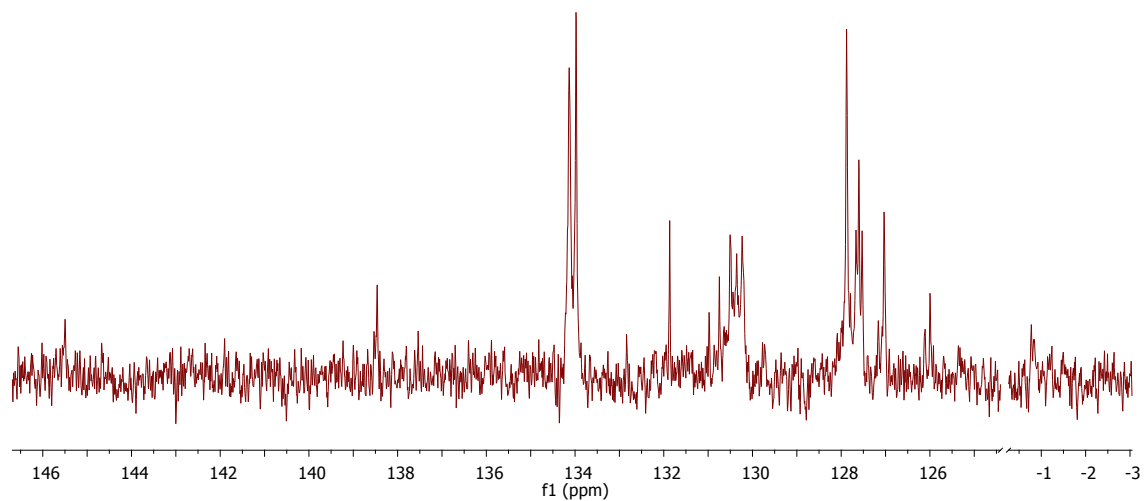


7: *cis* and *trans* mixture of di[9,19-(*E*)-4-(2,3,4,5,6-pentafluorophenyl)styrylmethyl]-1,3,5,7,11,13,15,17-octaphenylpentacyclo[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane

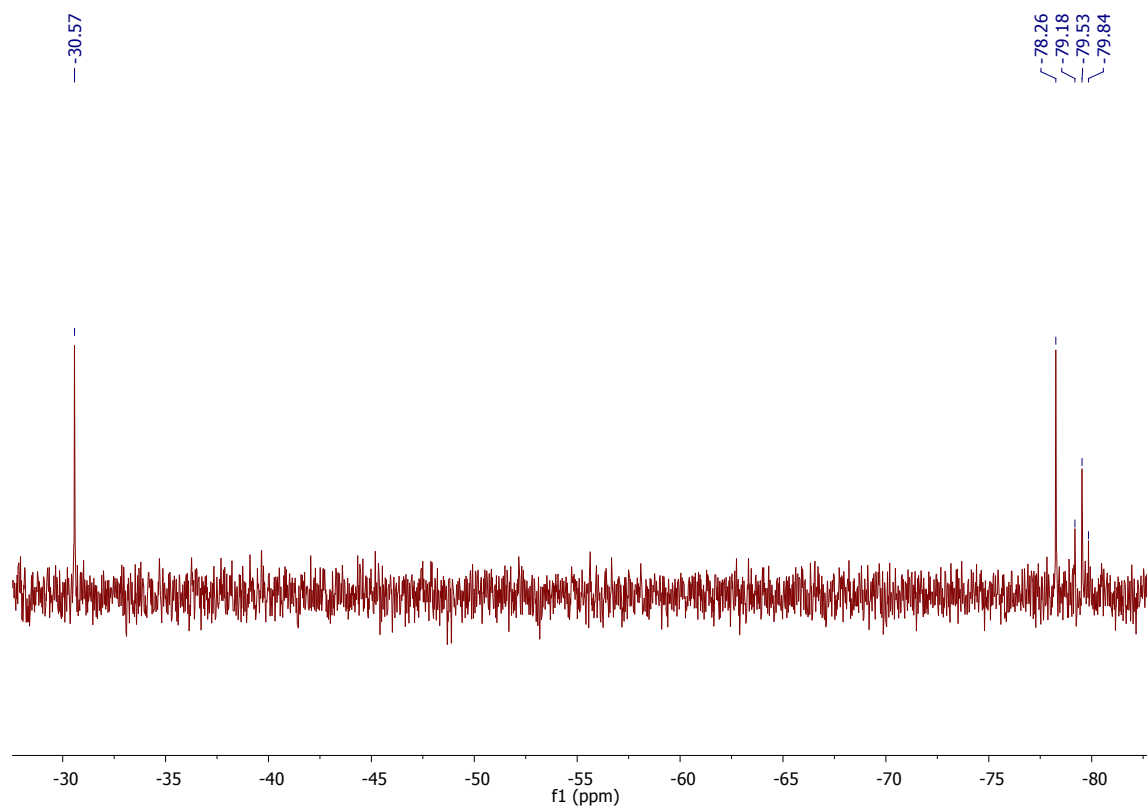
¹H NMR (CDCl₃, δ, ppm):



¹³C NMR (CDCl₃, δ, ppm):

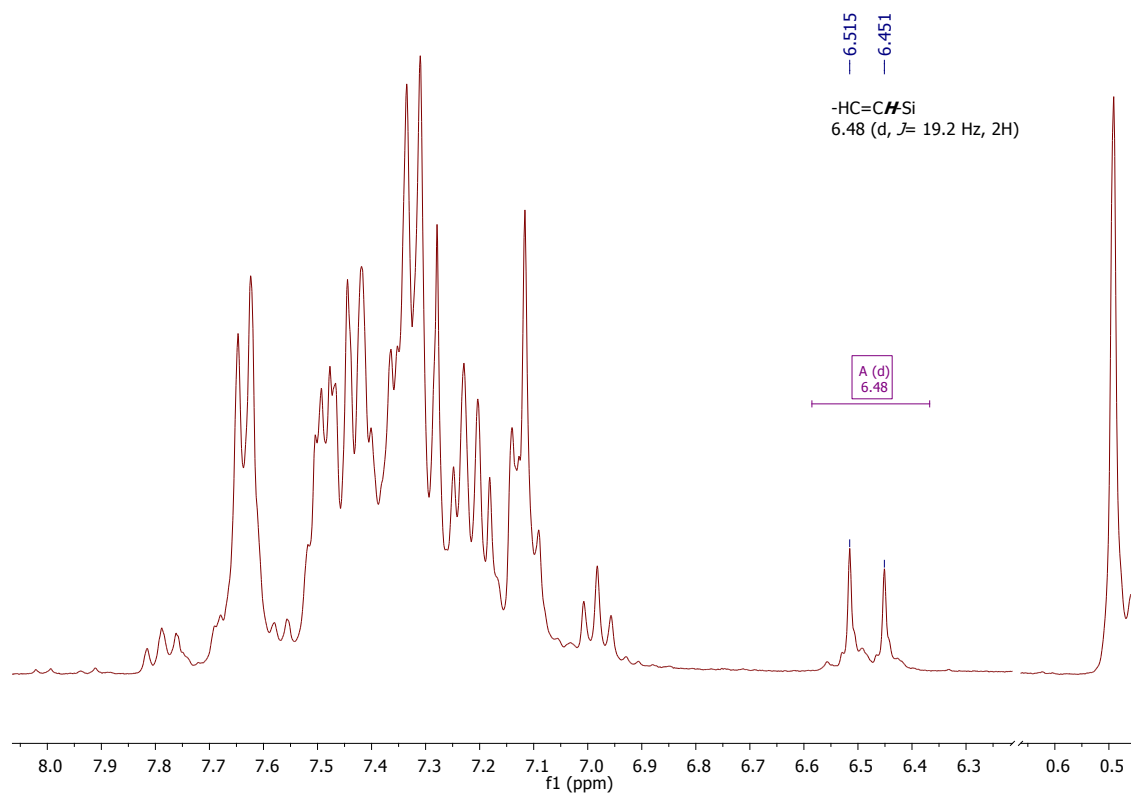


^{29}Si NMR (CDCl_3 , δ , ppm):

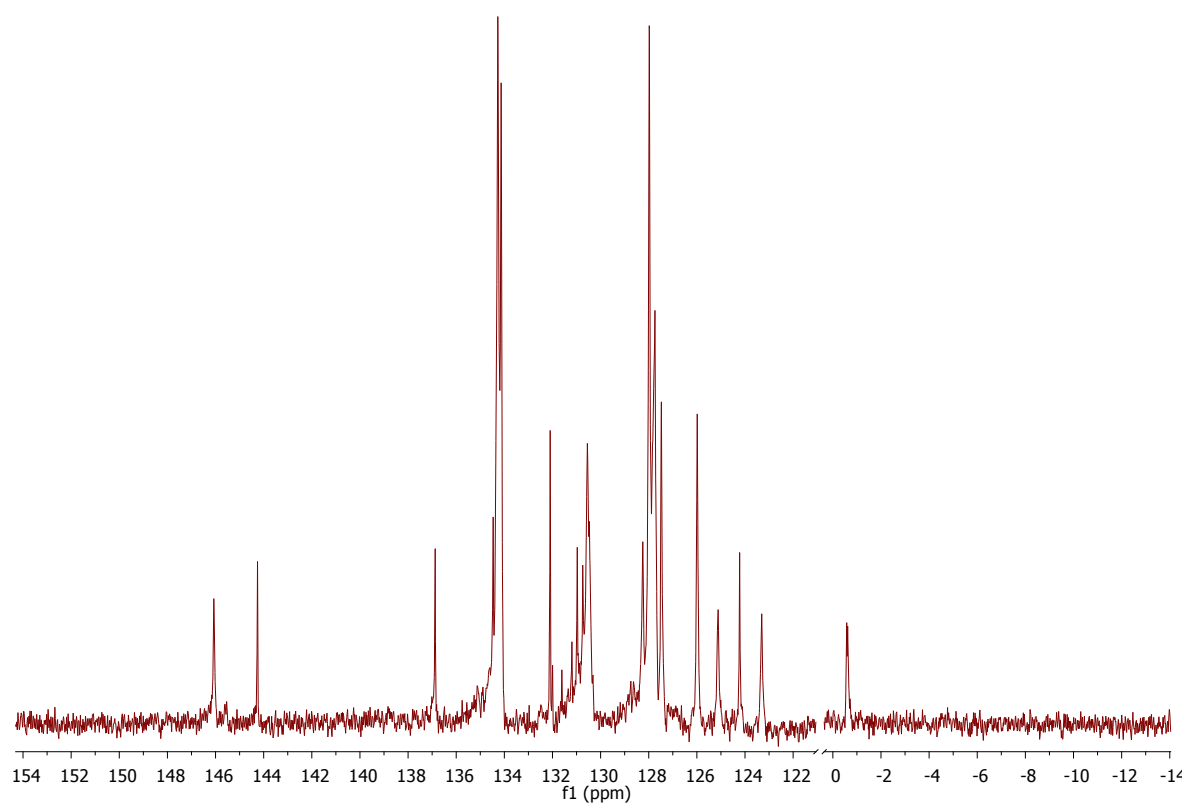


8: *cis* and *trans* mixture of di[9,19-(*E*)-4-(2-thienyl)styrylmethyl]-1,3,5,7,11,13,15,17-octaphenylpentacyclo[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane

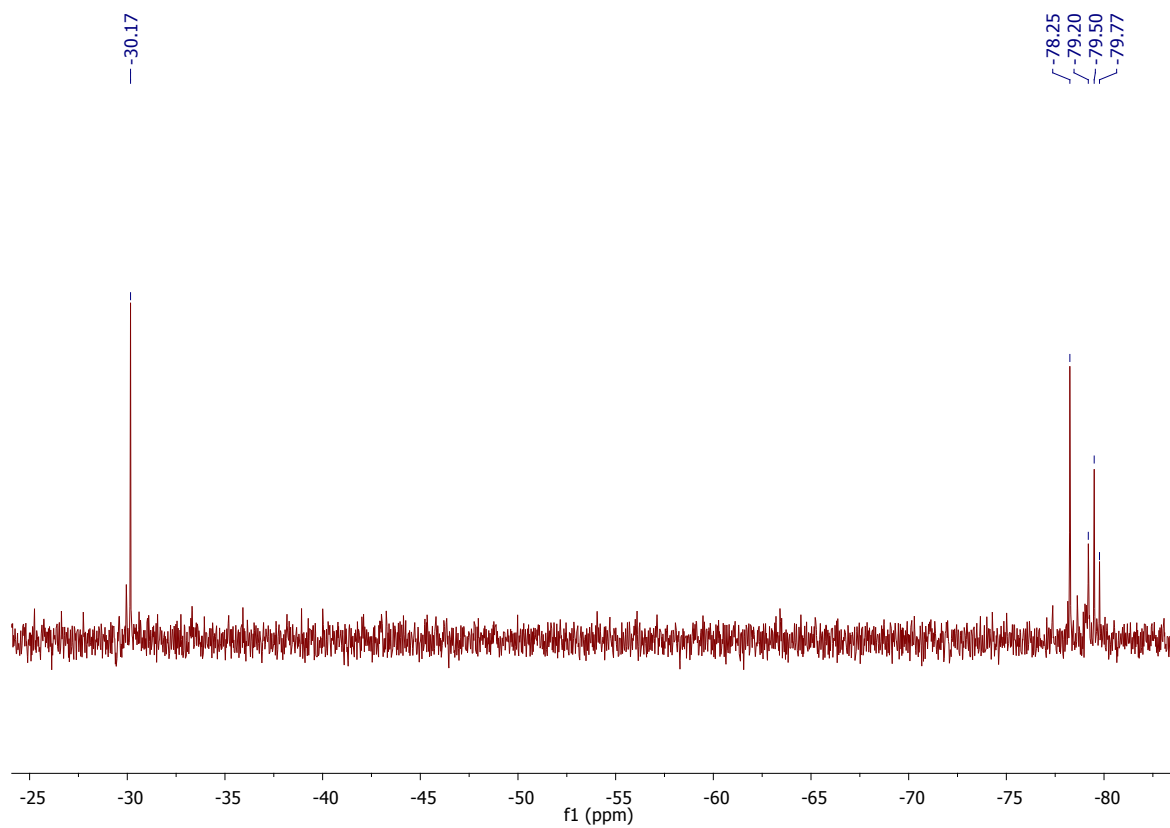
^1H NMR (CDCl_3 , δ , ppm):



^{13}C NMR (CDCl_3 , δ , ppm):



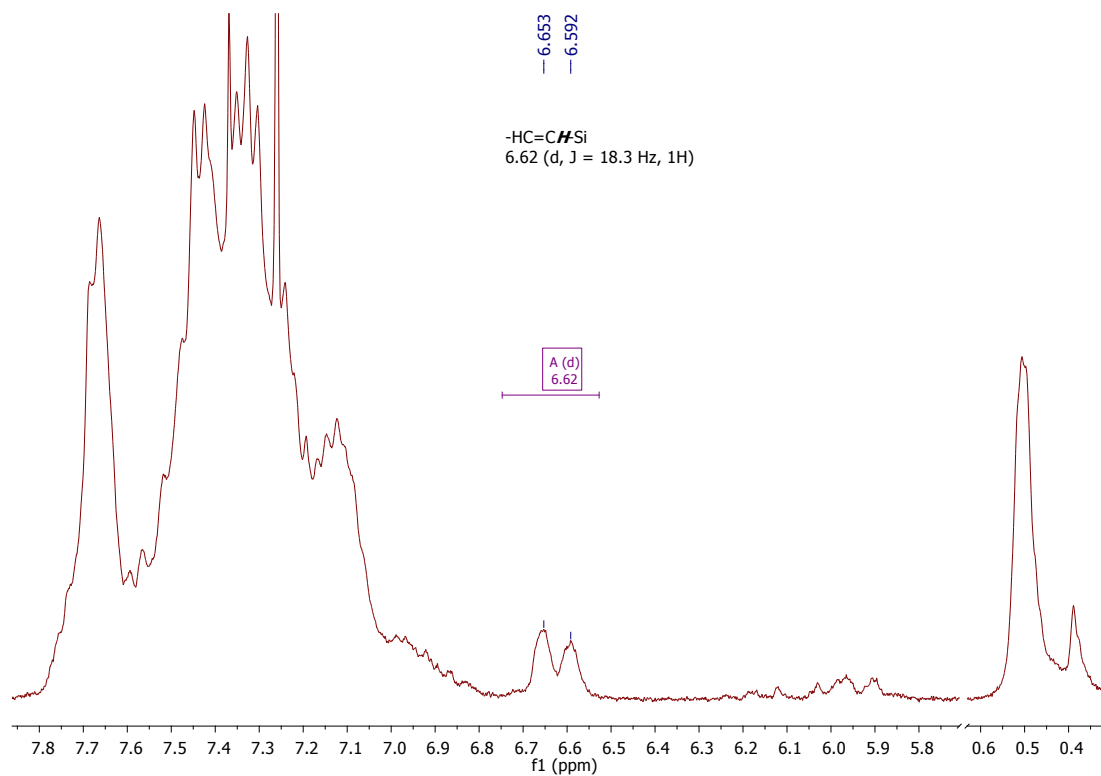
^{29}Si NMR (CDCl_3 , δ , ppm):



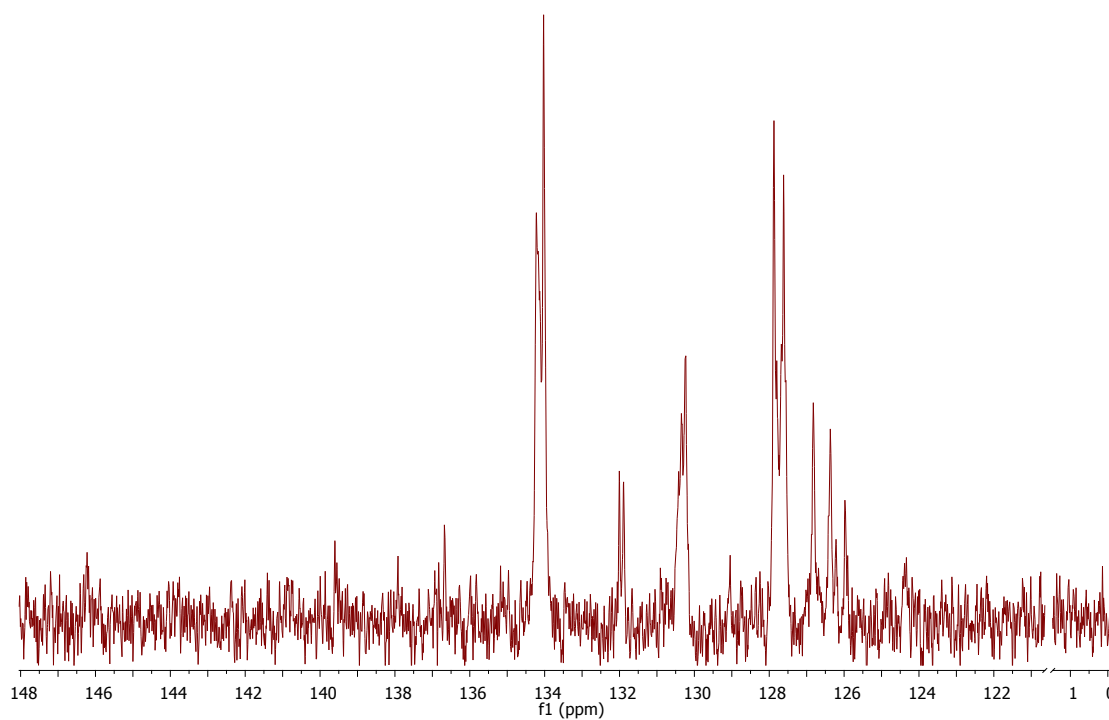
^1H , ^{13}C , ^{29}Si NMR spectra of the for isolated molecular compounds (9-15).

9: *poly{[9,19-(E)-4-(1-naphthyl)styrylmethyl]-1,3,5,7,11,13,15,17-octaphenylpentacyclo-[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane}s*

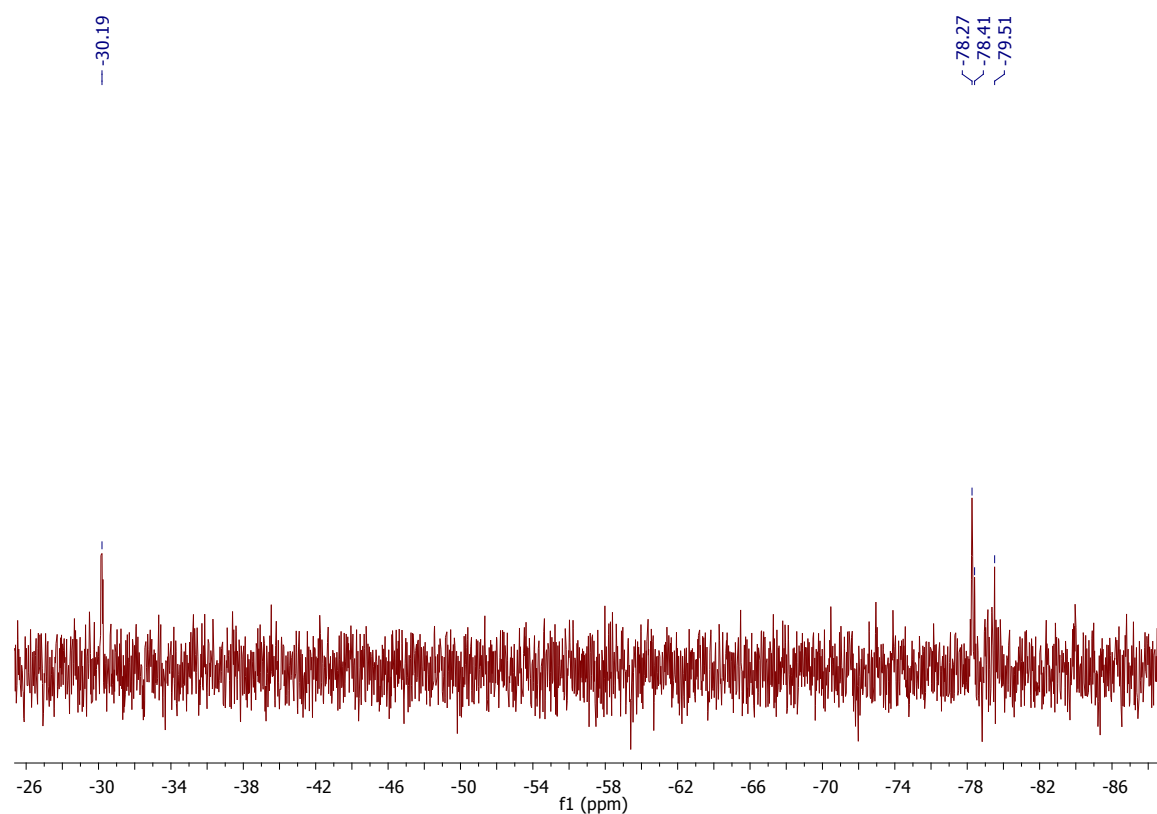
^1H NMR (CDCl_3 , δ , ppm):



^{13}C NMR (CDCl_3 , δ , ppm):

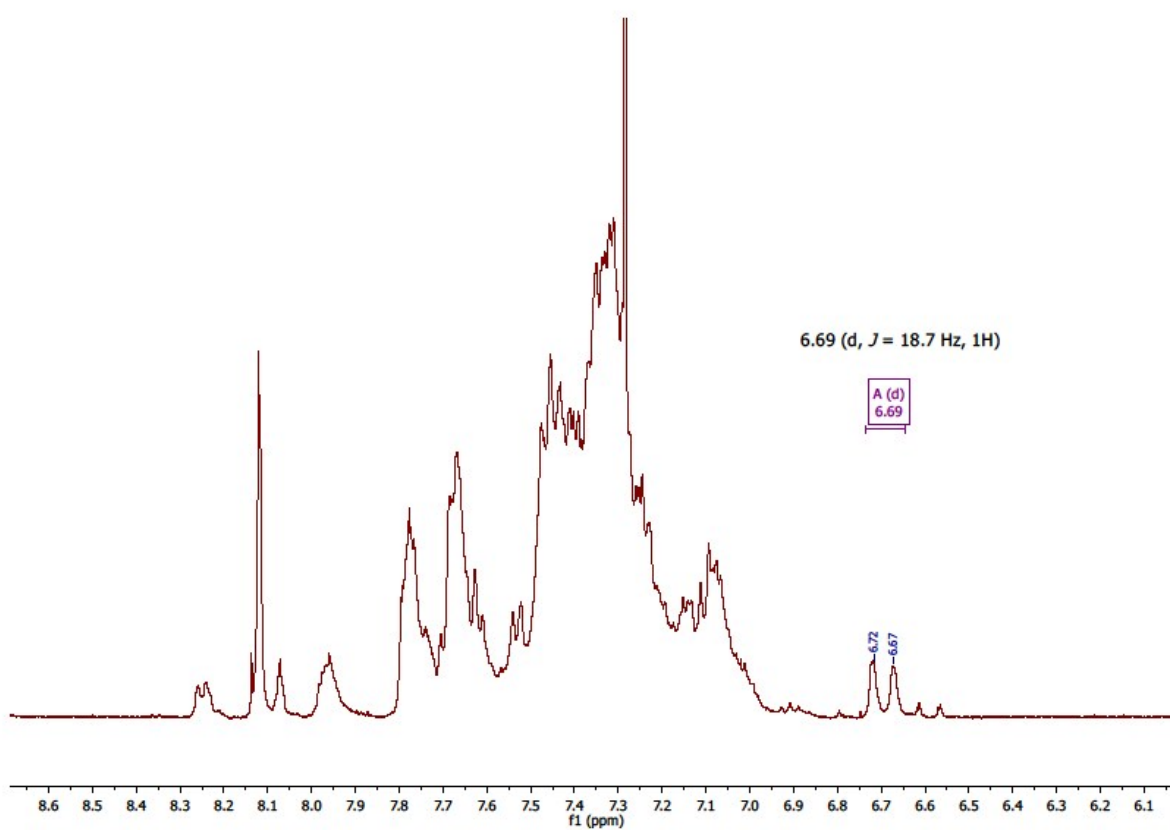


^{29}Si NMR (CDCl_3 , δ , ppm):

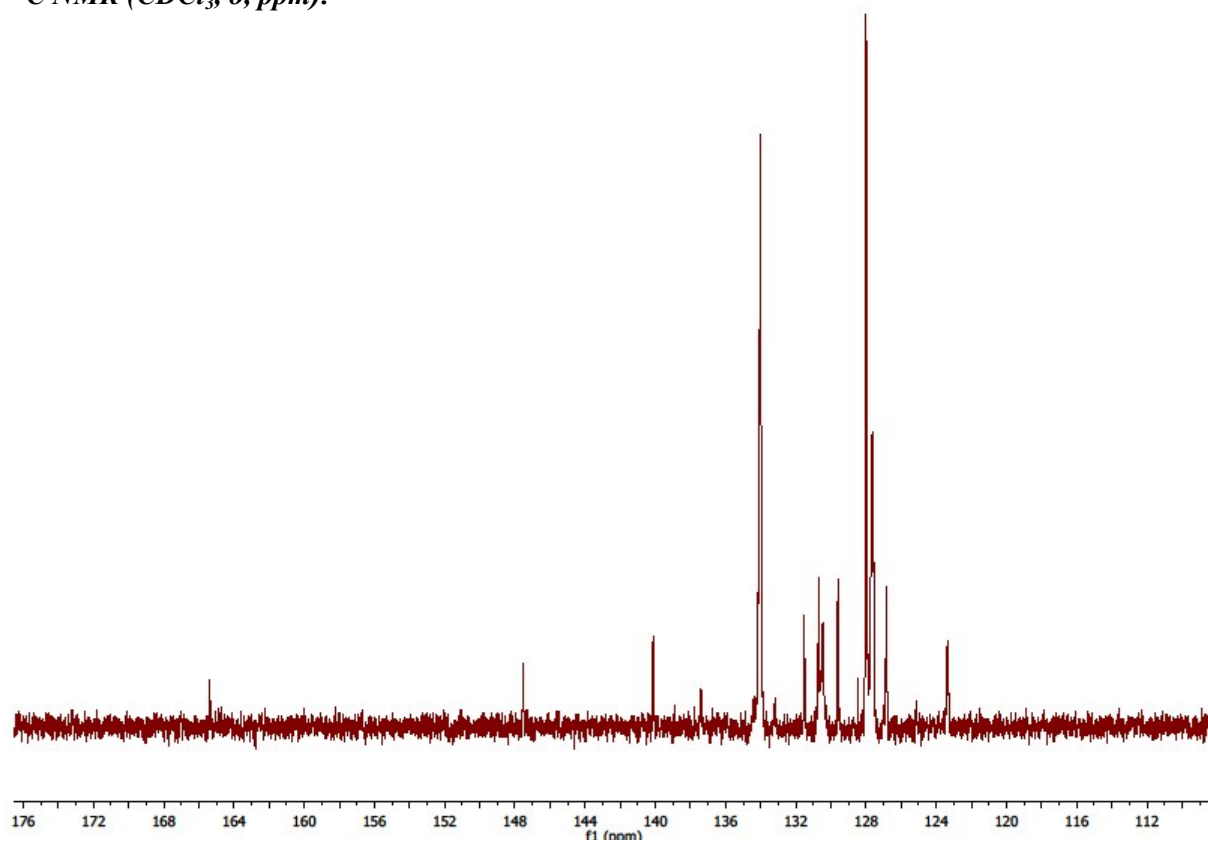


10: poly{[9,19-(E)-4-(1-naphthyl)styrylphenyl]-1,3,5,7,11,13,15, 17-octaphenylpentacyclo-[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane}s

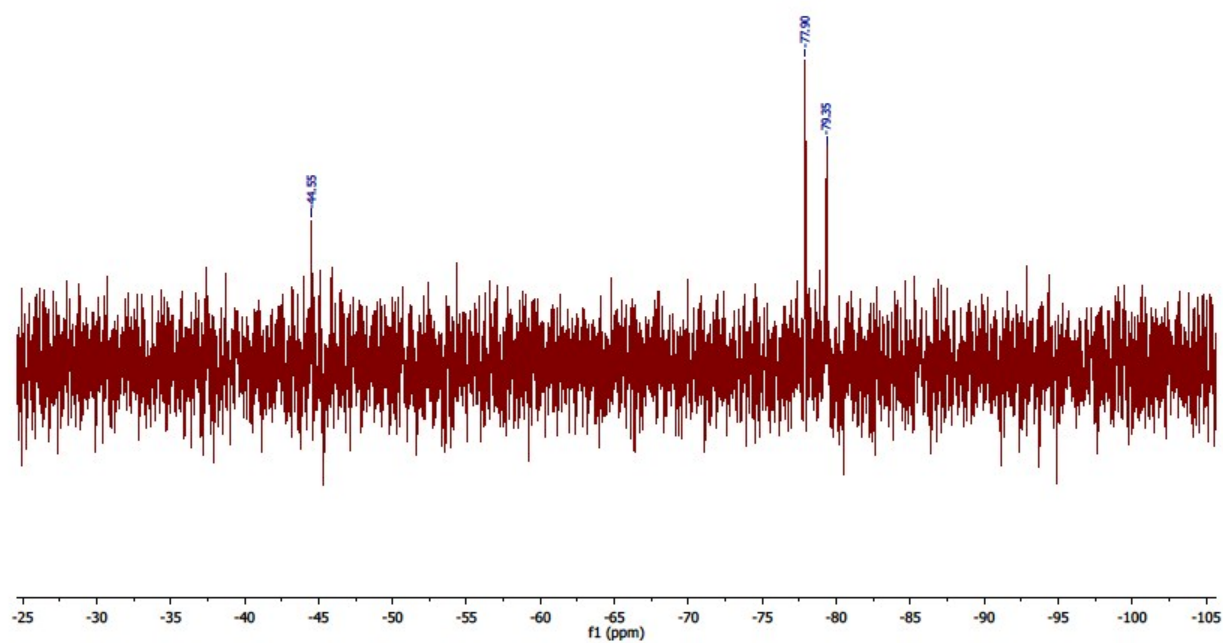
^1H NMR (CDCl_3 , δ , ppm):



^{13}C NMR (CDCl_3 , δ , ppm):

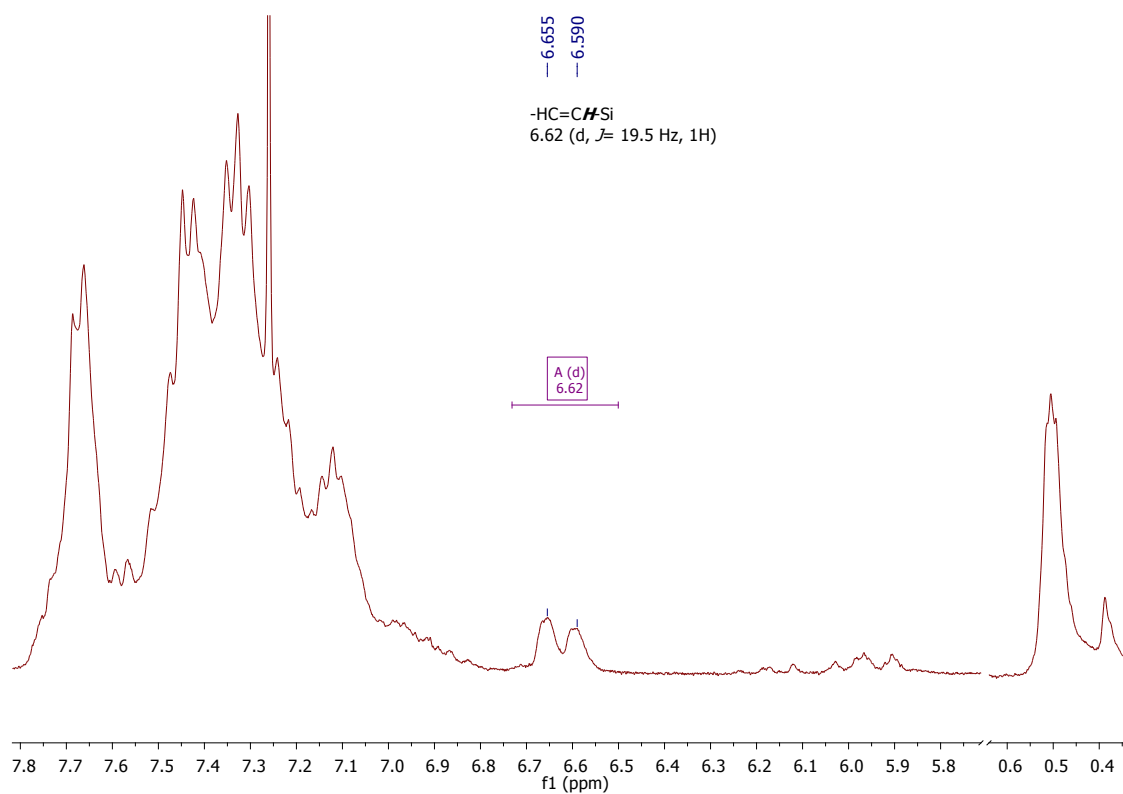


^{29}Si NMR (CDCl_3 , δ , ppm):

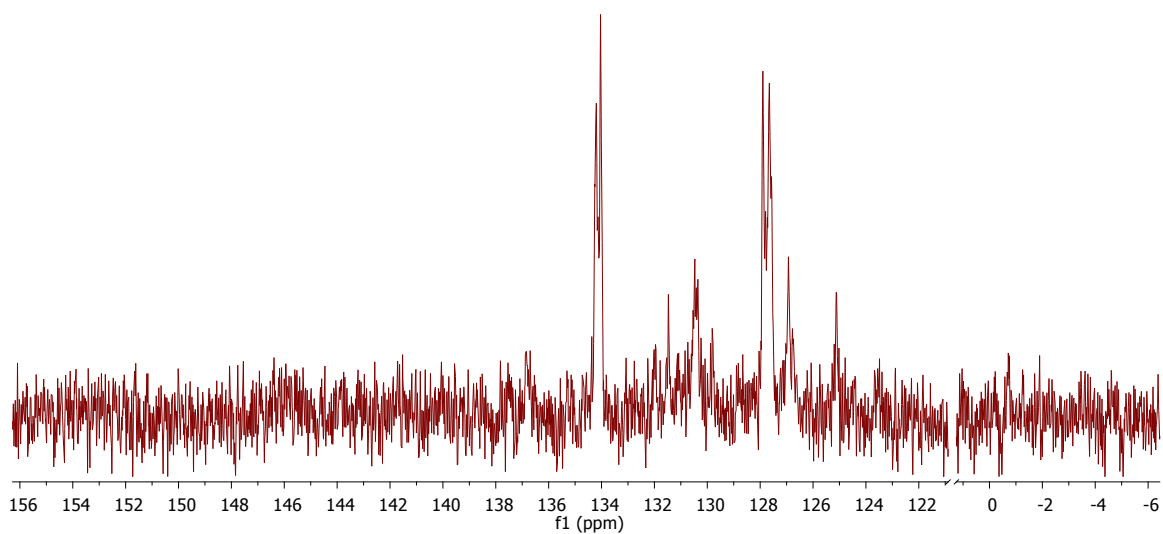


**11: poly{[9,19-(E)-4-(9-anthracenyl)styrylmethyl]-1,3,5,7,11,13,15,17-octaphenylpentacyclo
[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane}s**

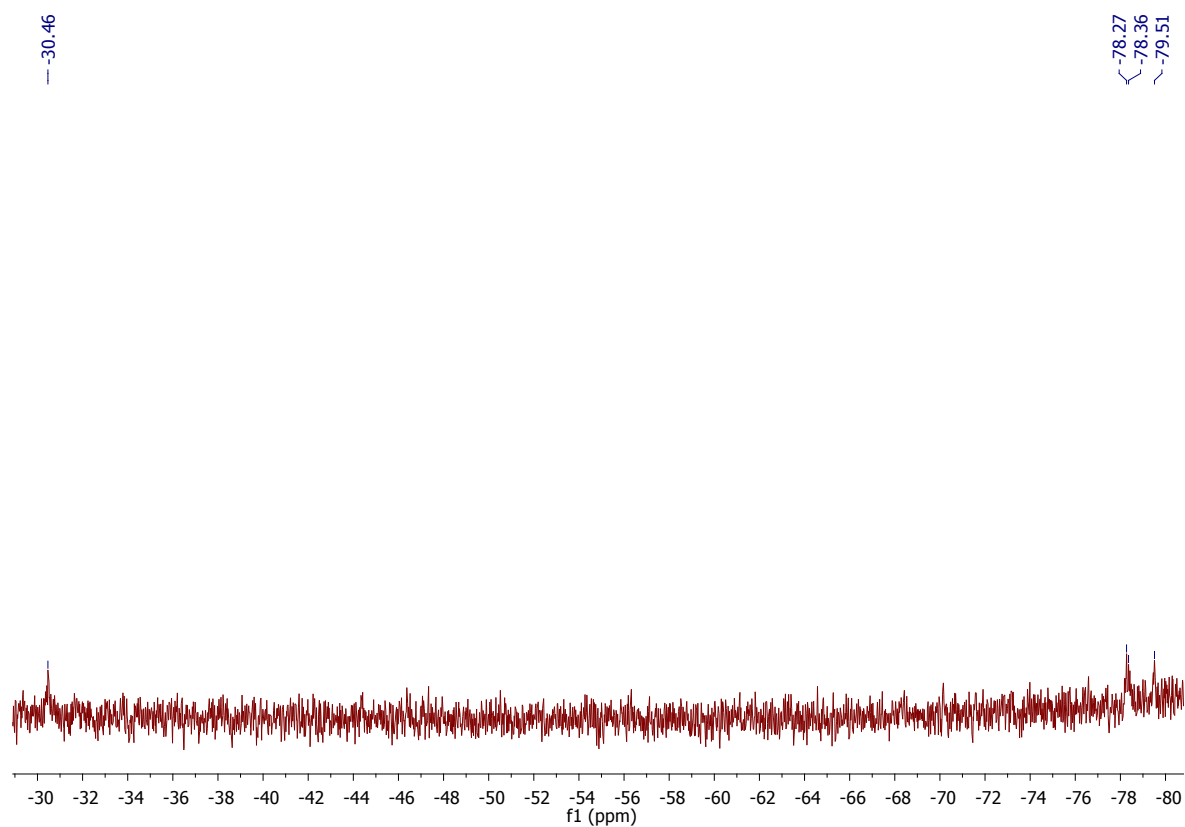
¹H NMR (CDCl₃, δ, ppm):



¹³C NMR (CDCl₃, δ, ppm):

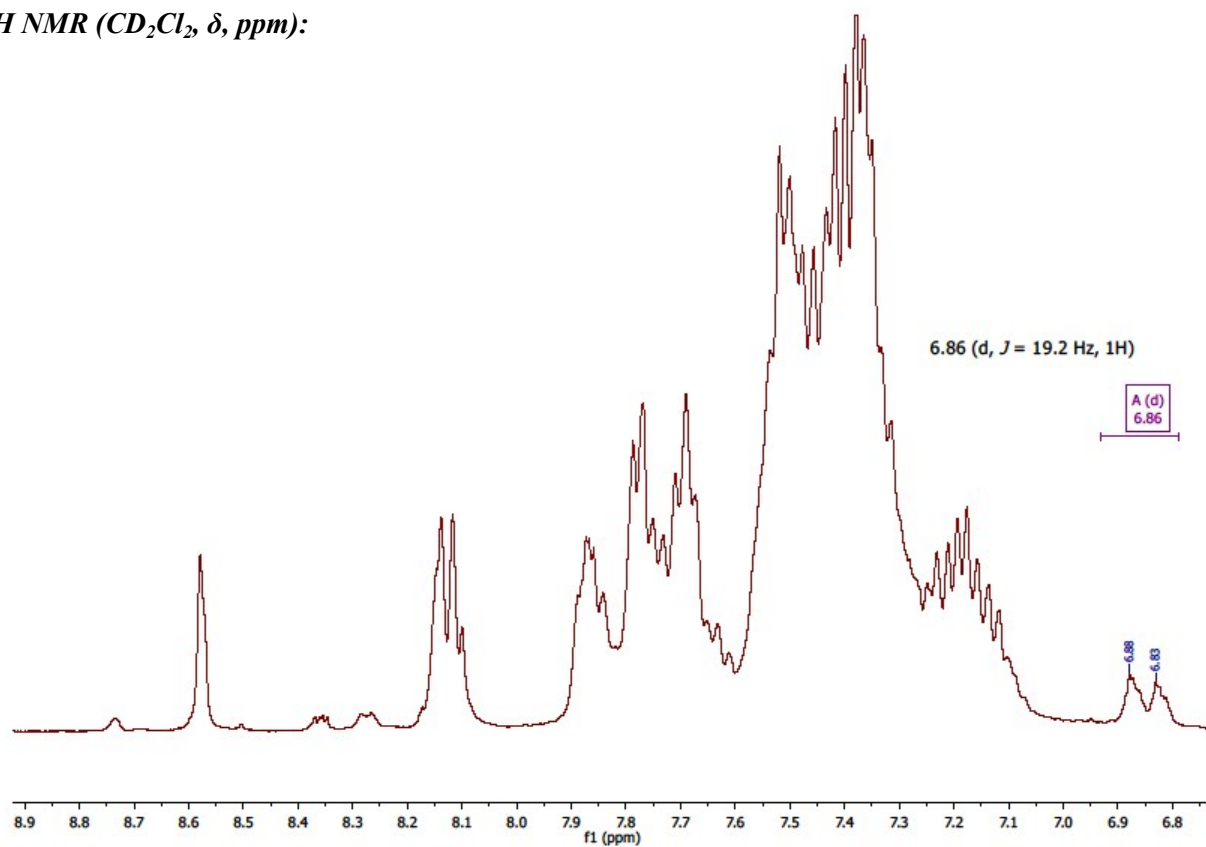


^{29}Si NMR (CDCl_3 , δ , ppm):

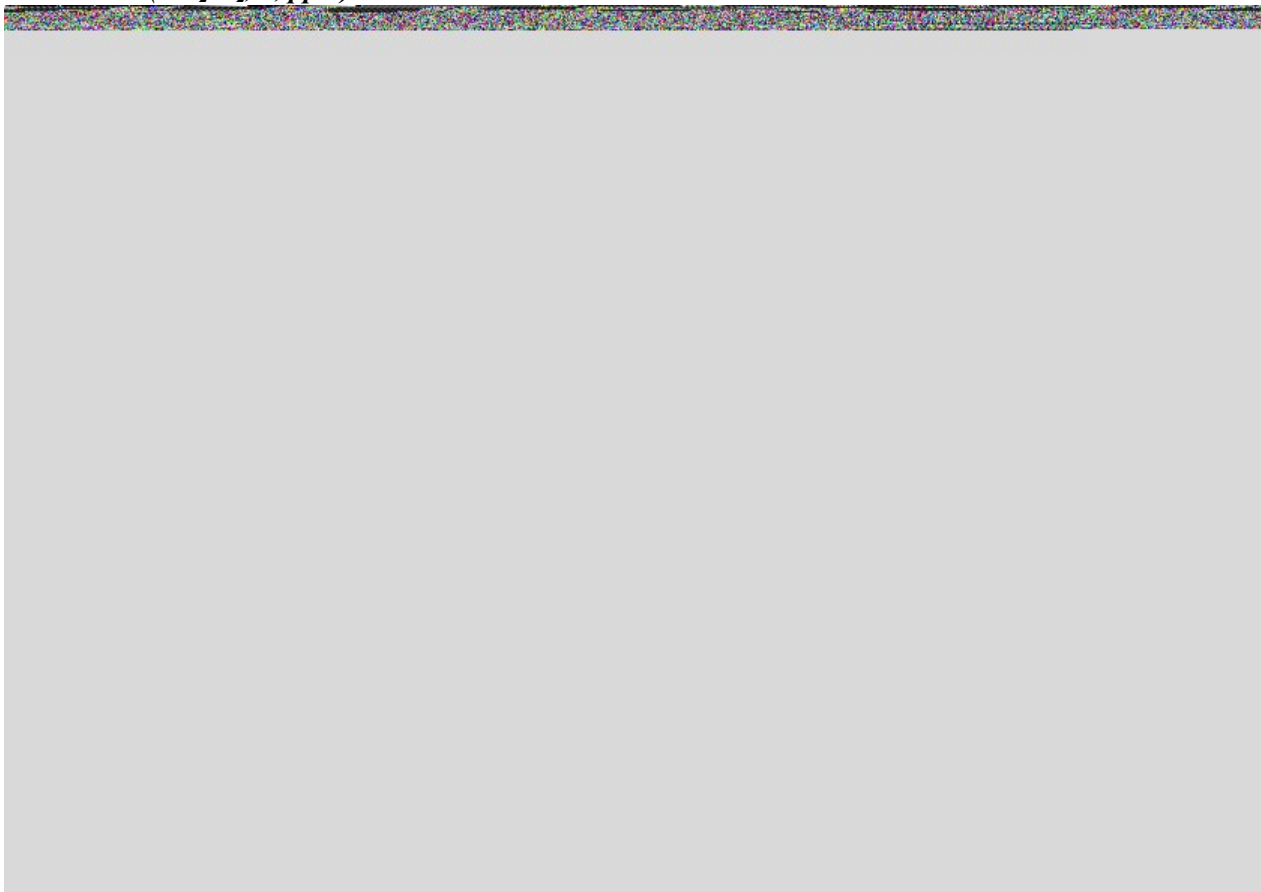


**12: poly{[9,19-(E)-4-(9-anthracenyl)styrylphenyl- 1,3,5,7,11,13,15,17-octaphenylpentacyclo-
[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane}s**

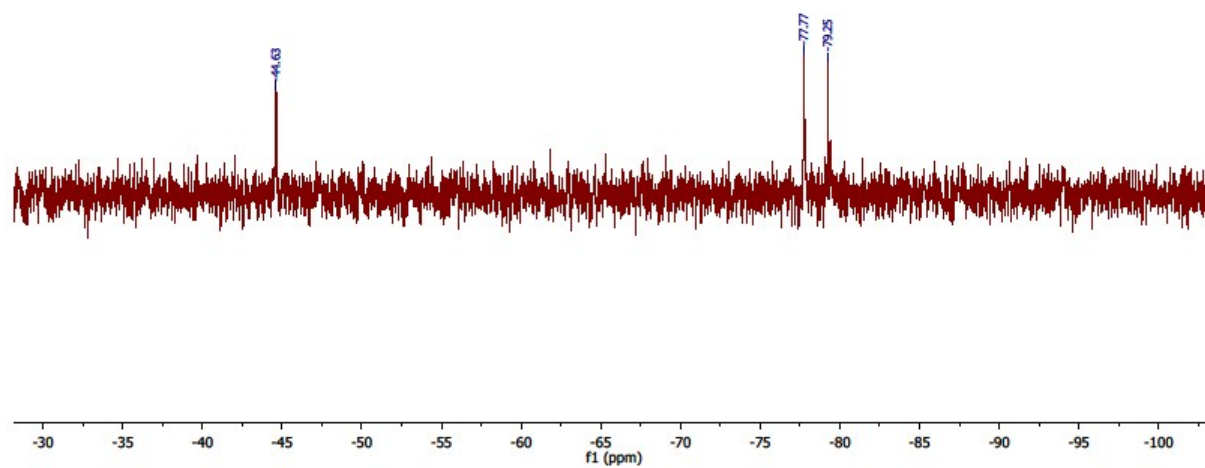
^1H NMR (CD_2Cl_2 , δ , ppm):



^{13}C NMR (CD_2Cl_2 , δ , ppm):

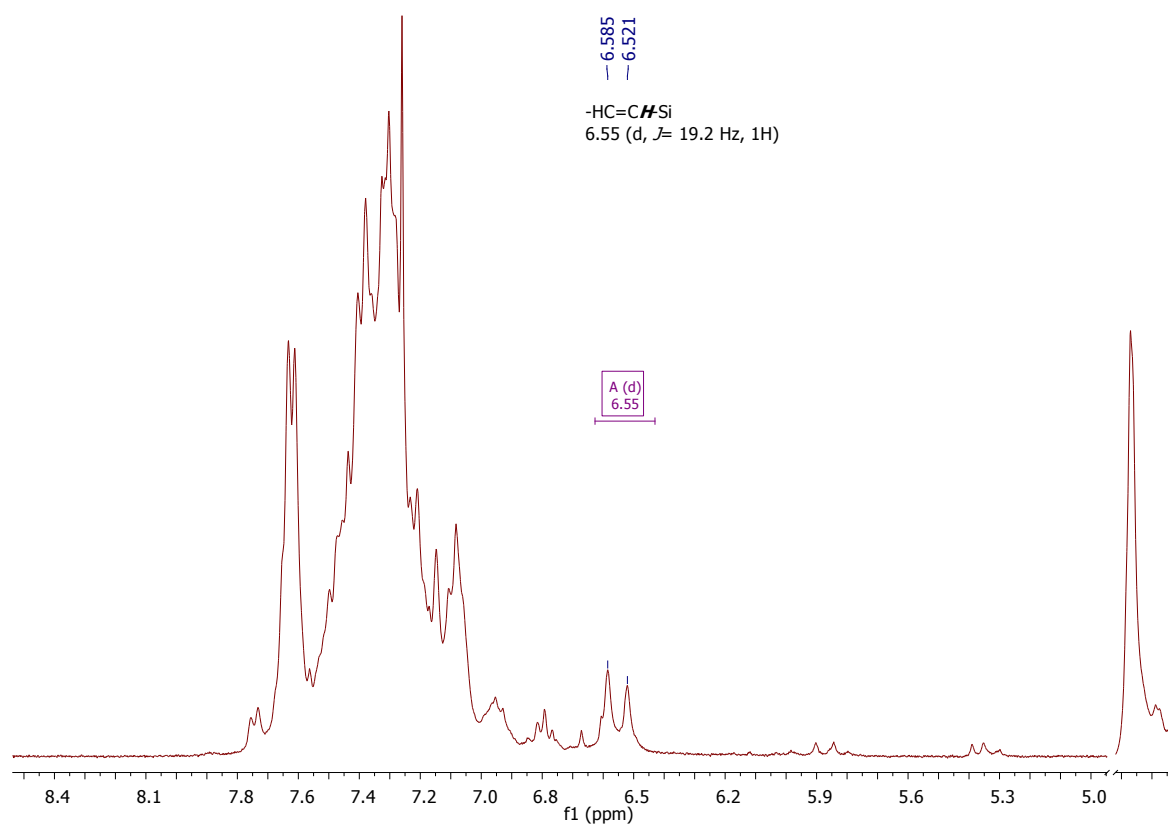


^{29}Si NMR (CD_2Cl_2 , δ , ppm):

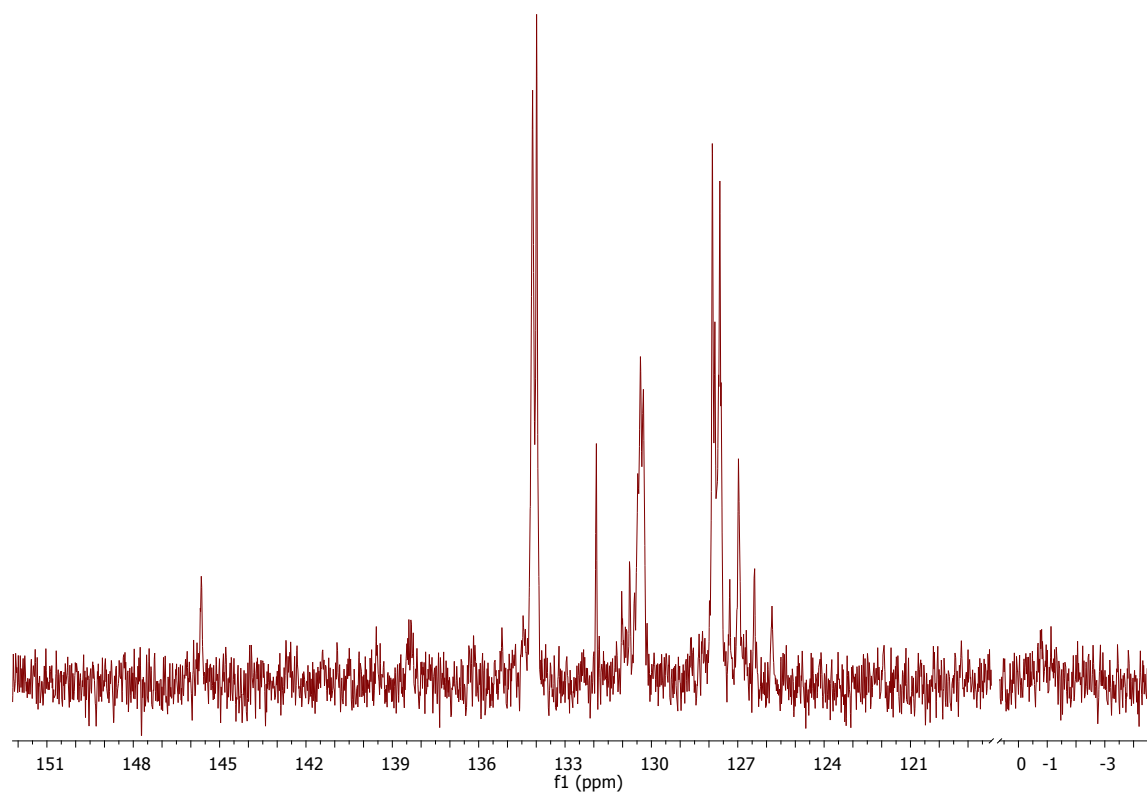


13: *poly{[9,19-(E)-4-(2,3,5,6-tetrafluorophenyl)styrylmethyl]-1,3,5,7,11,13,15,17-octaphenylpentacyclo[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane}s*

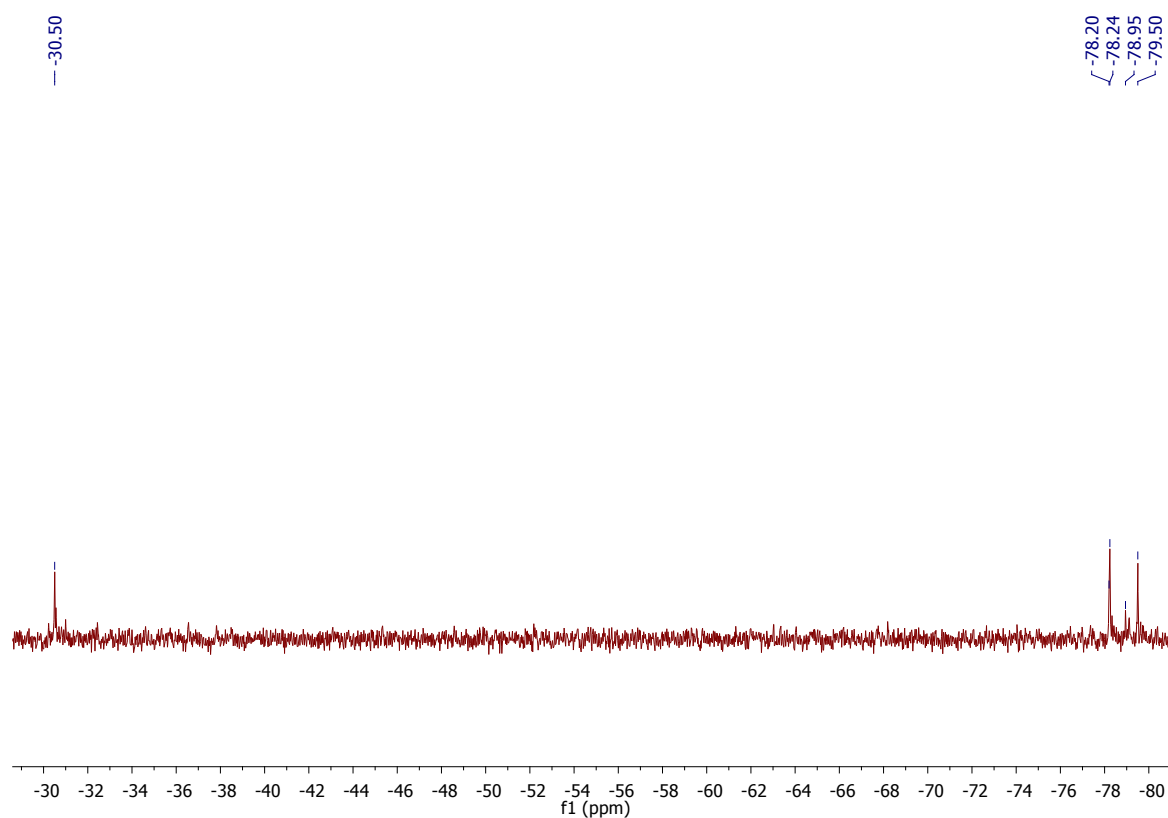
¹H NMR (CDCl₃, δ, ppm):



¹³C NMR (CDCl₃, δ, ppm):

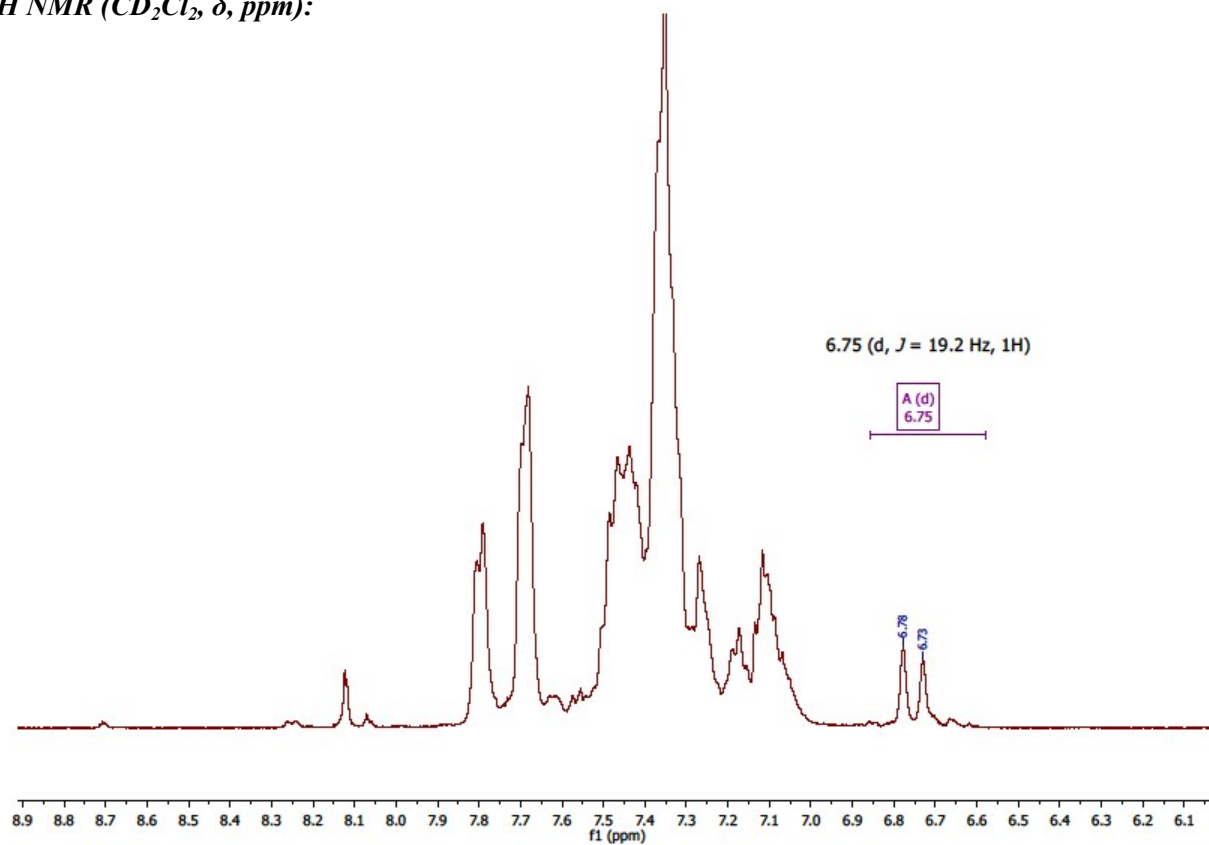


^{29}Si NMR (CDCl_3 , δ , ppm):

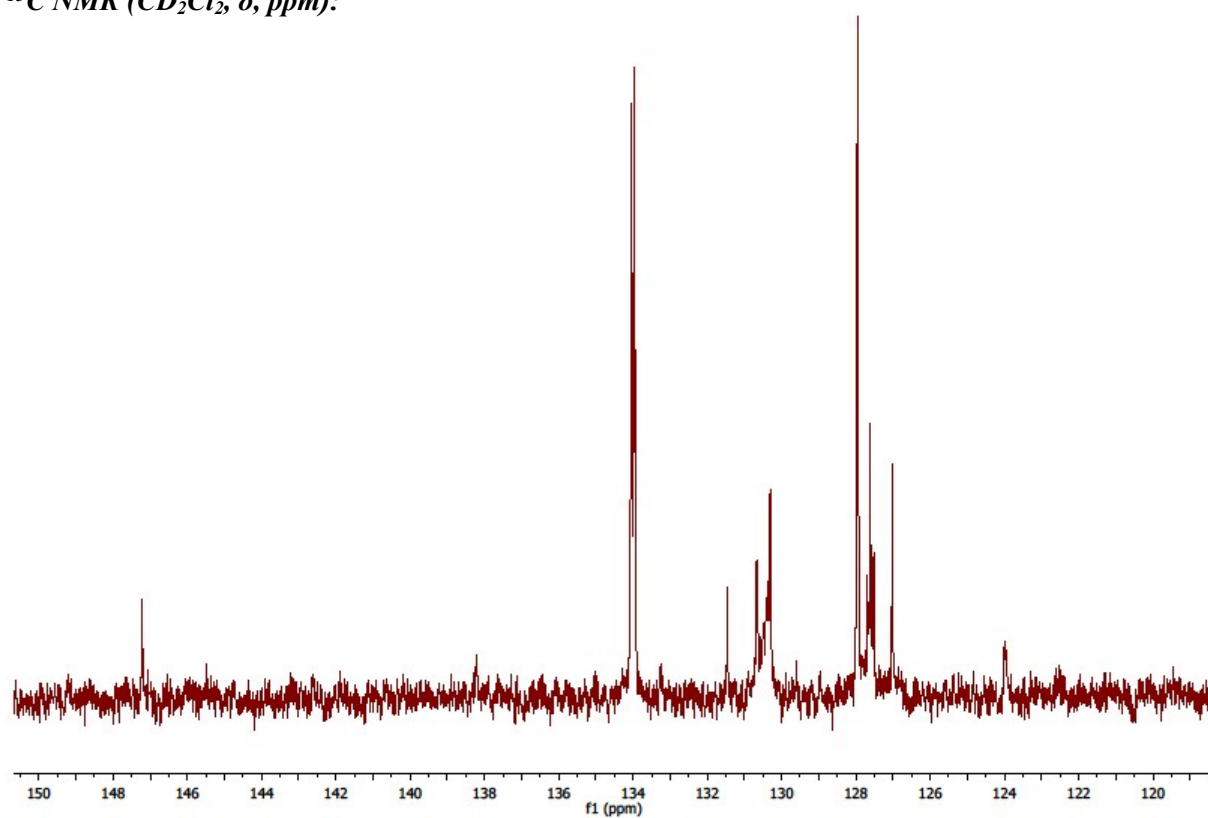


14: poly{[9,19-(*E*)-4-(2,3,5,6-tetrafluorophenyl)styrylphenyl]-1,3,5,7,11,13,15,17-octaphenylpenta-cyclo[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane}s

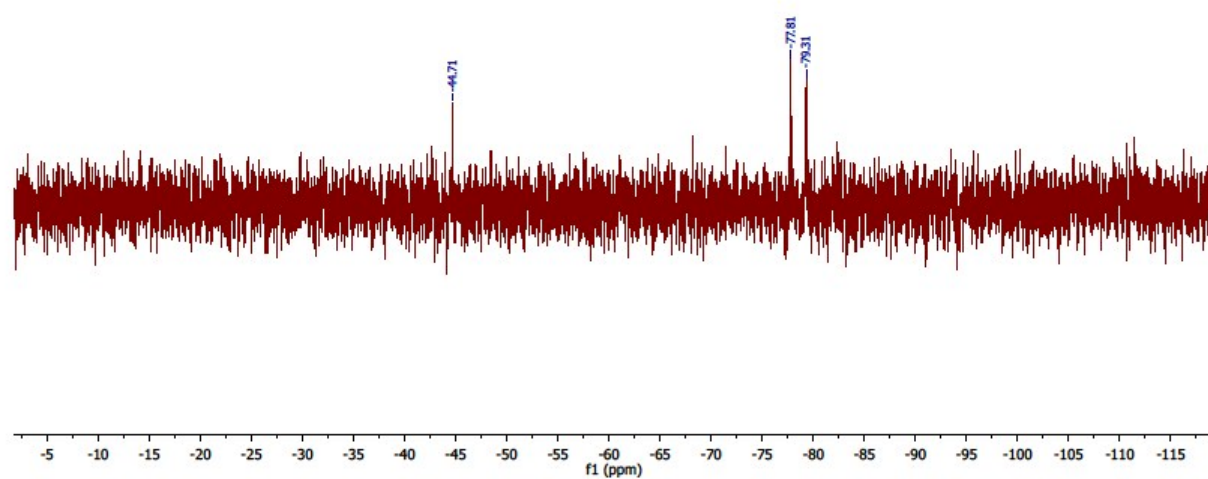
^1H NMR (CD_2Cl_2 , δ , ppm):



^{13}C NMR (CD_2Cl_2 , δ , ppm):

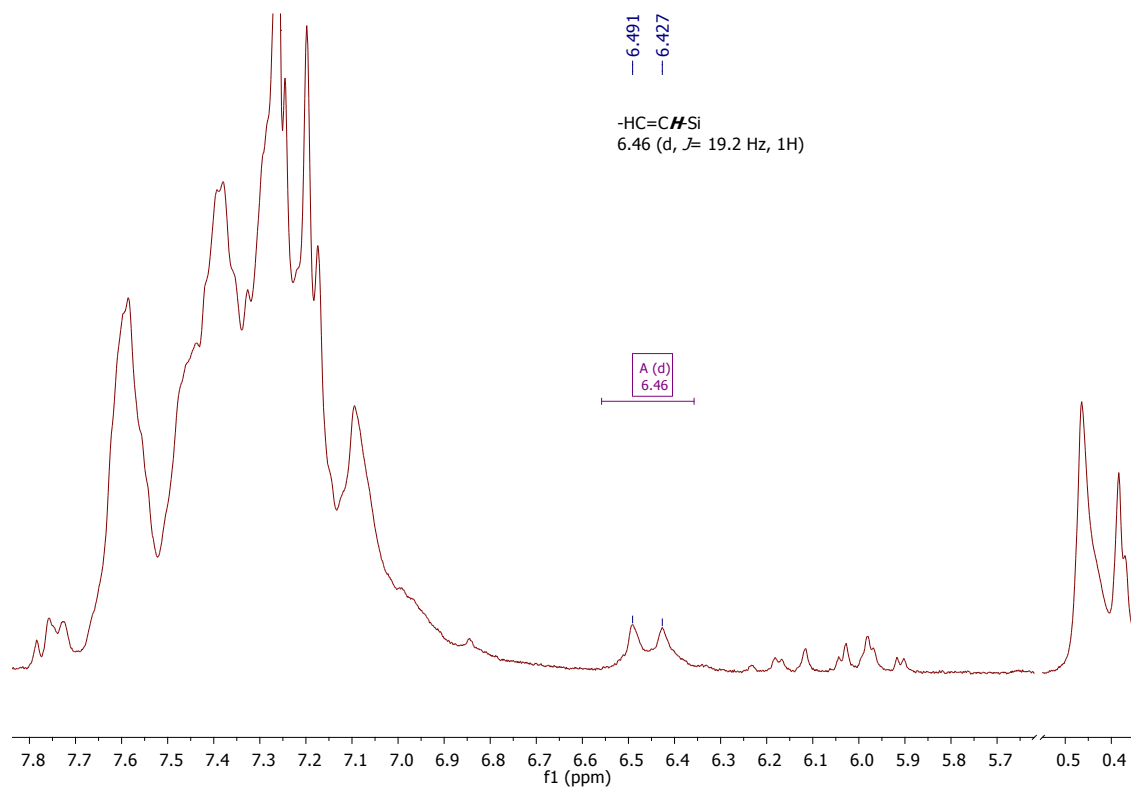


^{29}Si NMR (CD_2Cl_2 , δ , ppm):

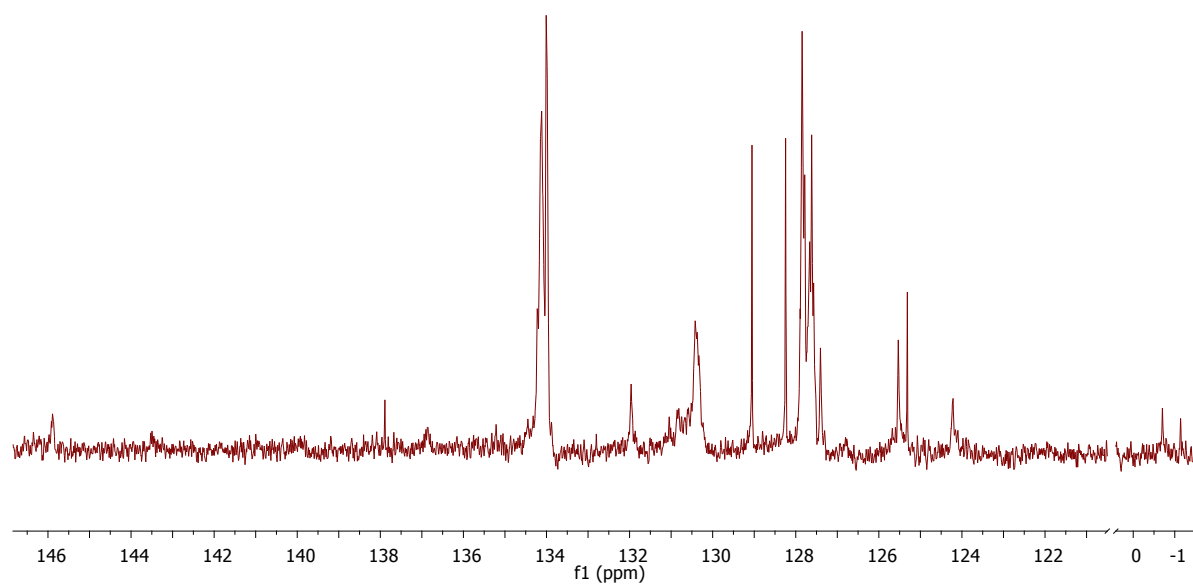


15: *poly{[9,19-(E)-4-(2-thienyl)styrylmethyl]-1,3,5,7,11,13,15,17-octaphenylpenta-*
cyclo[11.7.1.1^{3,11}.1^{5,17}.1^{7,15}]decasiloxane}s

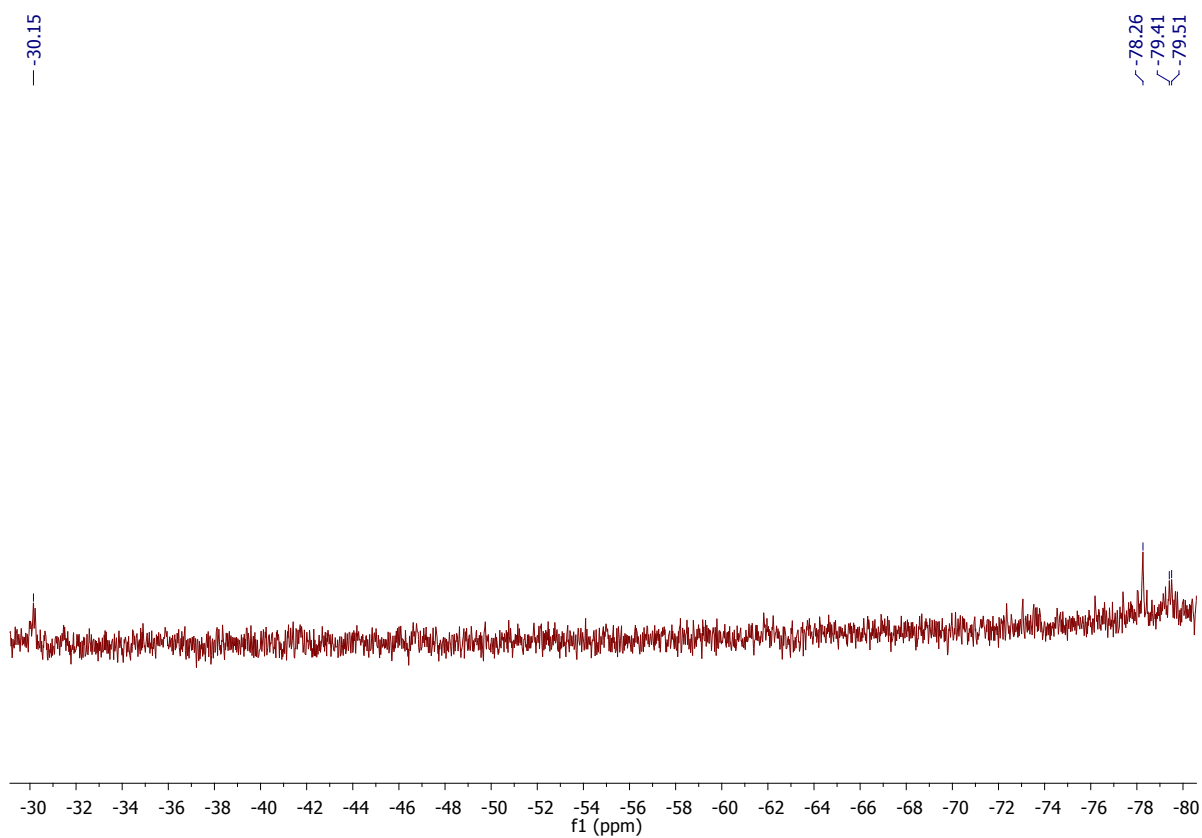
¹H NMR (CDCl₃, δ, ppm):



¹³C NMR (CDCl₃, δ, ppm):

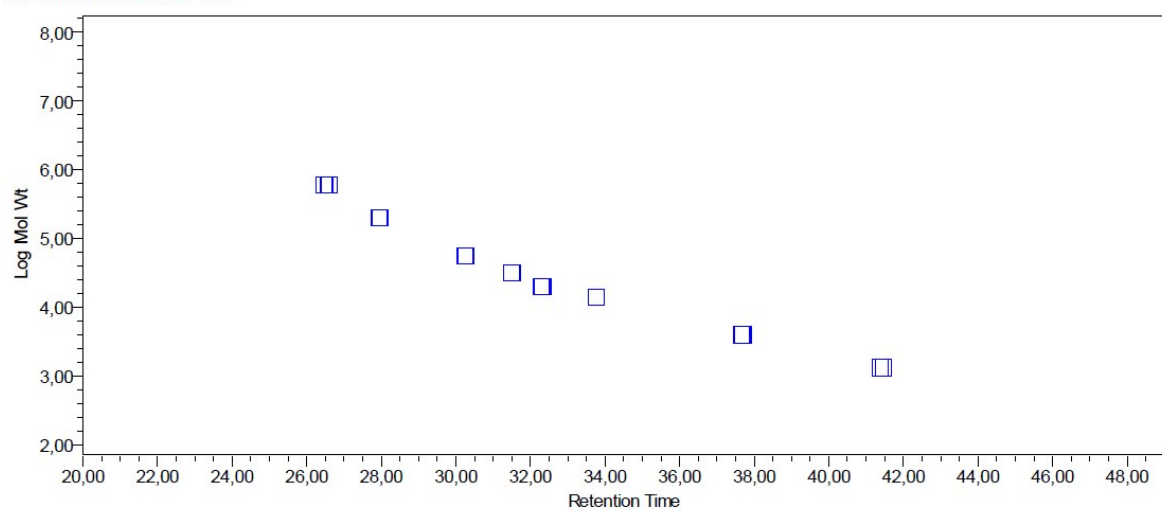


^{29}Si NMR (CDCl_3 , δ , ppm):



GPC Calibration Curve

GPC Calibration Plot



GPC Calibration Table

	Mol Wt (Daltons)	Retention Time (min)	Calculated Weight (Daltons)	% Residual
1	591000	26,483	610648	-3,218
2	591000	26,480	611933	-3,421
3	591000	26,609	554136	6,652
4	591000	26,590	562486	5,069
5	591000	26,465	619034	-4,529
6	197000	27,943	198660	-0,836
7	197000	27,959	196423	0,294
8	197000	27,954	196940	0,030
9	197000	27,947	198023	-0,517
10	197000	27,966	195711	0,659
11	55100	30,243	55250	-0,271
12	55100	30,242	55260	-0,290
13	55100	30,244	55200	-0,182
14	55100	30,246	55142	-0,075
15	55100	30,262	54738	0,662
16	31400	31,500	31441	-0,130
17	31400	31,507	31322	0,250
18	31400	31,495	31507	-0,338
19	31400	31,506	31352	0,153
20	31400	31,506	31343	0,182
21	20000	32,326	19920	0,403
22	20000	32,322	19941	0,296
23	20000	32,321	19944	0,279
24	20000	32,299	20124	-0,618
25	20000	32,282	20315	-1,548
26	13900	33,773	13862	0,273
27	13900	33,767	13890	0,075
28	13900	33,761	13914	-0,098
29	13900	33,762	13910	-0,074
30	13900	33,761	13914	-0,098
31	3950	37,668	3979	-0,734
32	3950	37,699	3940	0,248
33	3950	37,701	3939	0,283
34	3950	37,697	3943	0,167
35	3950	37,691	3951	-0,015
36	1310	41,376	1322	-0,920
37	1310	41,460	1289	1,593
38	1310	41,385	1319	-0,655