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APPENDICES

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APPENDIX A

Calculation of the correction factor

The correction factor was calculated based upon the results obtained from gas chromatographic analysis (see also the experimental section). The substrate used was 5 mmol, whereas the internal standard, cyclohexanone, 0.02 mmol was employed.

Example

A : the exact amount of prepared desired product (mmol)

B : total volume of the reaction (mL)

C : peak area of the desired product

D : peak area of the internal standard

The calculation of the correction factor can be described as follows:

The amount of the desired product from the reaction mixture

$$= (0.02 \times C / D) = F \text{ mmol}$$

The amount of the desired product in B mL (total volume of the reaction)

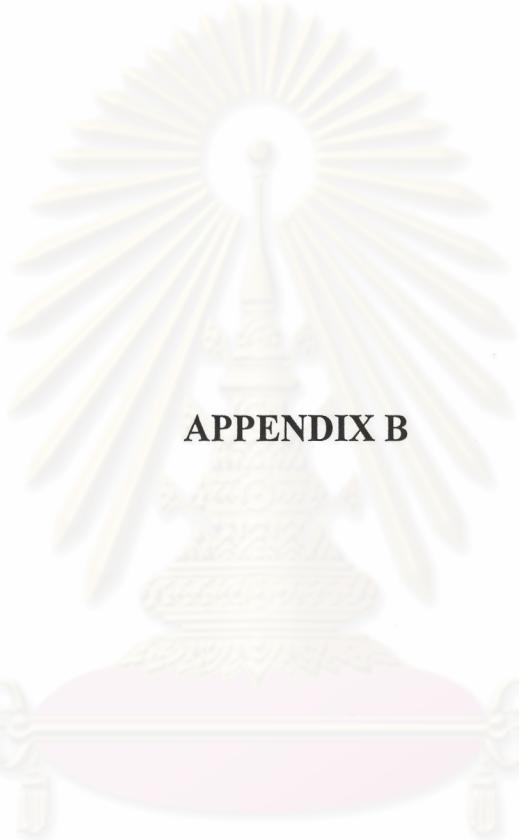
$$= (F) \times (B) = G \text{ mmol}$$

Thus, the correction factor of the desired product

$$= (A) / (G) = H$$

The calculation of % yield of the desired product can be calculated as:

$$\% \text{ Yield of product} = (G) \times (H) / 5 \times 100$$



APPENDIX B

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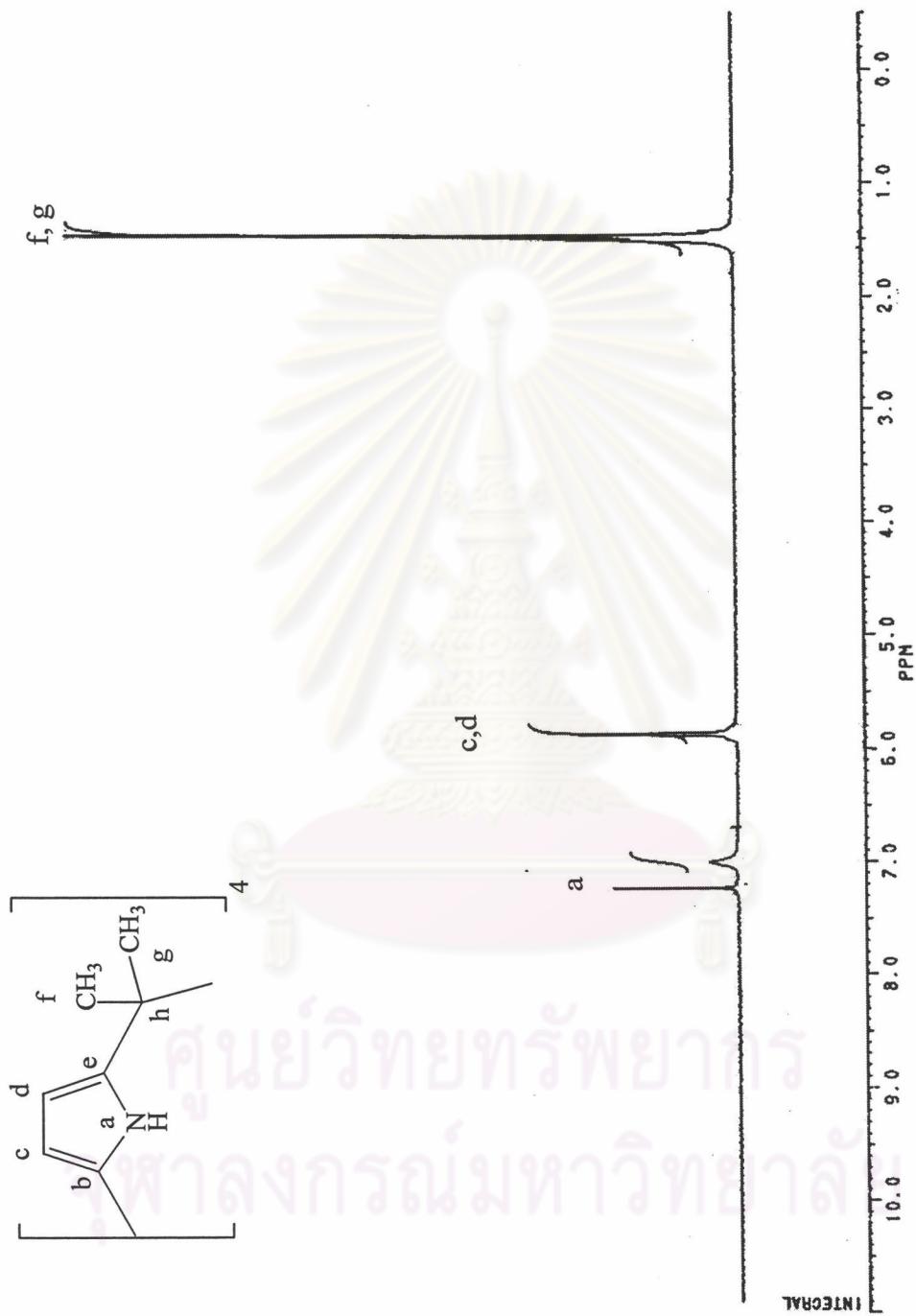


Figure 1 ^1H NMR spectrum of *meso*-octamethyl calix[4]pyrrole

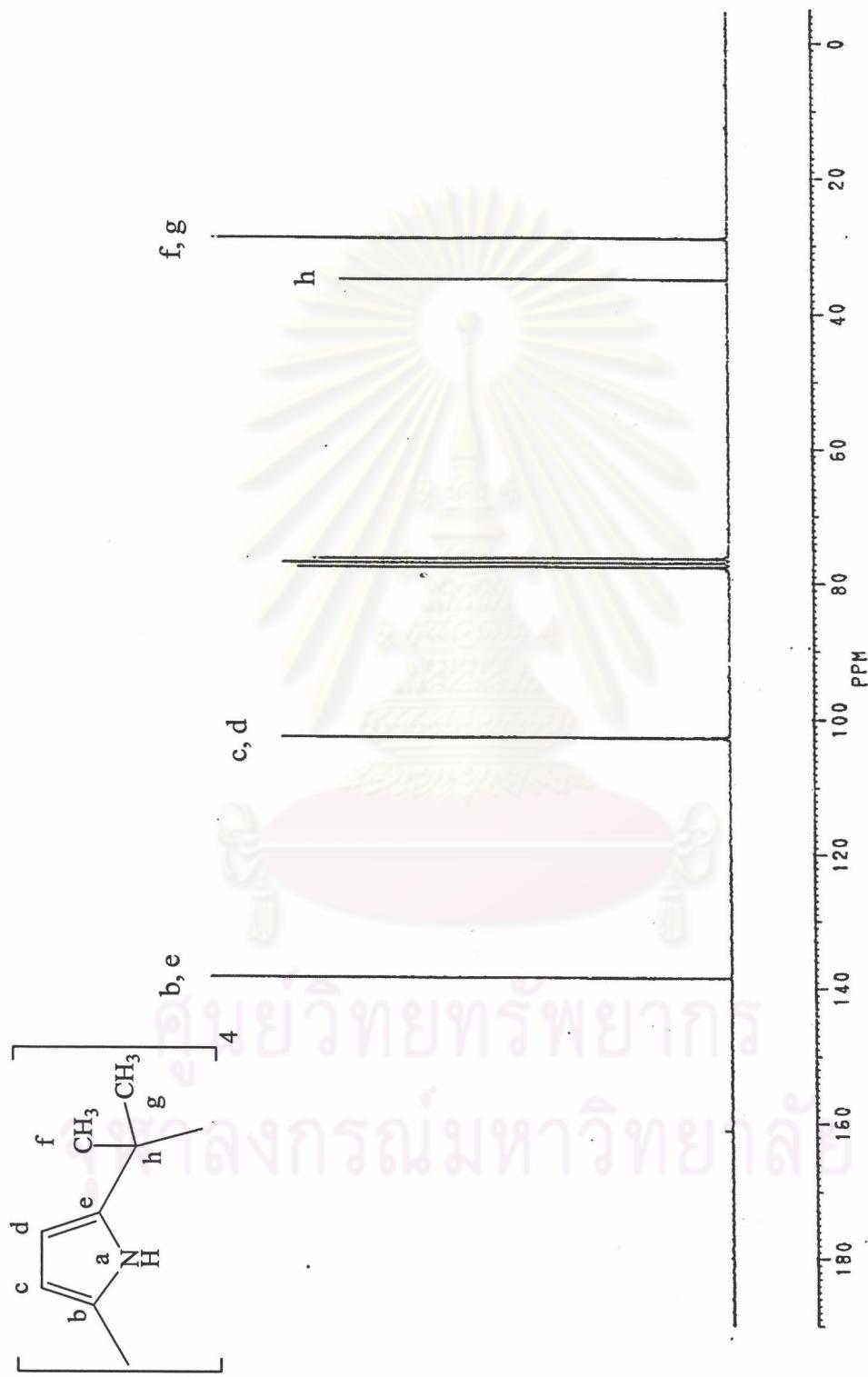


Figure 2 ^{13}C NMR spectrum of *meso*-octamethyl calix[4]pyrrole

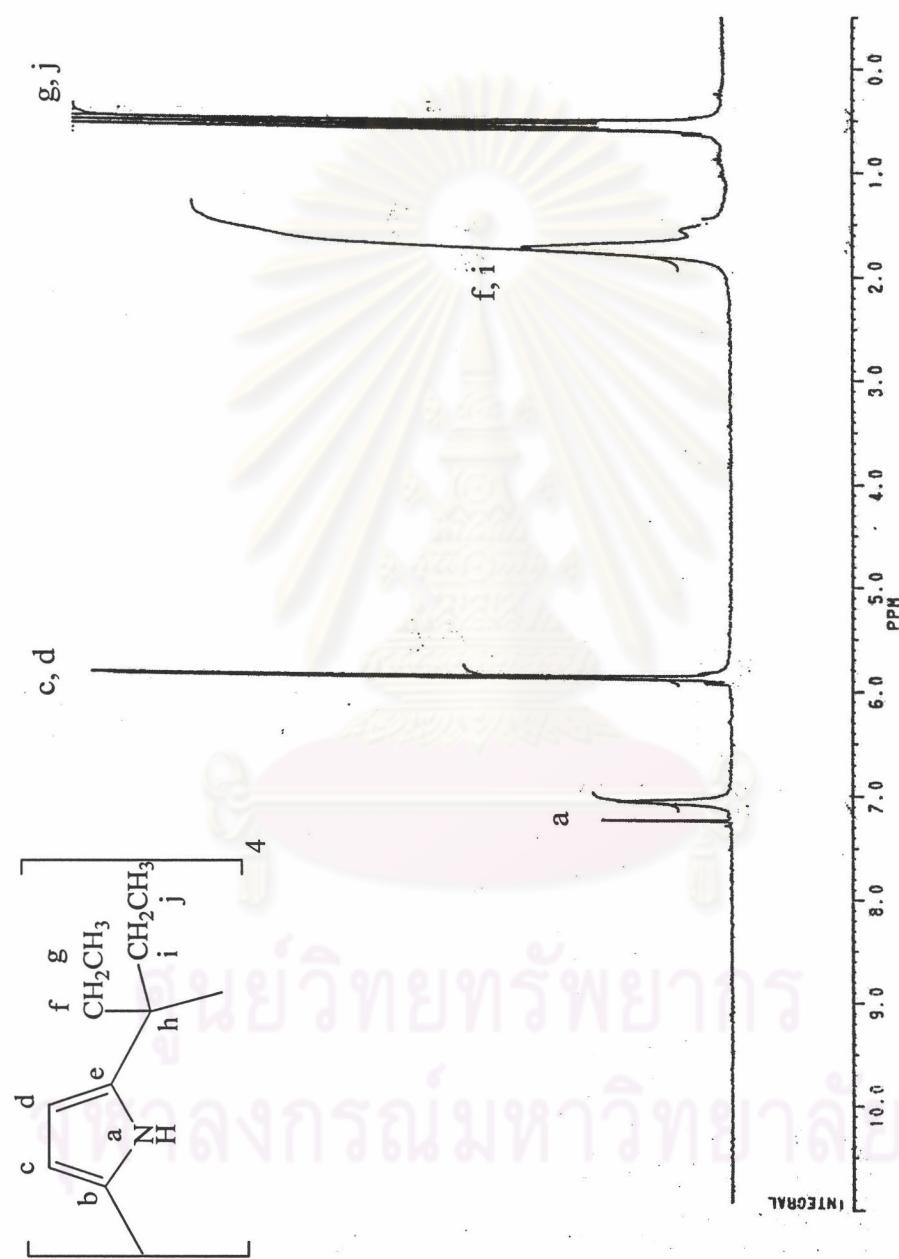


Figure 3 ^1H NMR spectrum of *meso*-octaethyl-calix[4]pyrrole

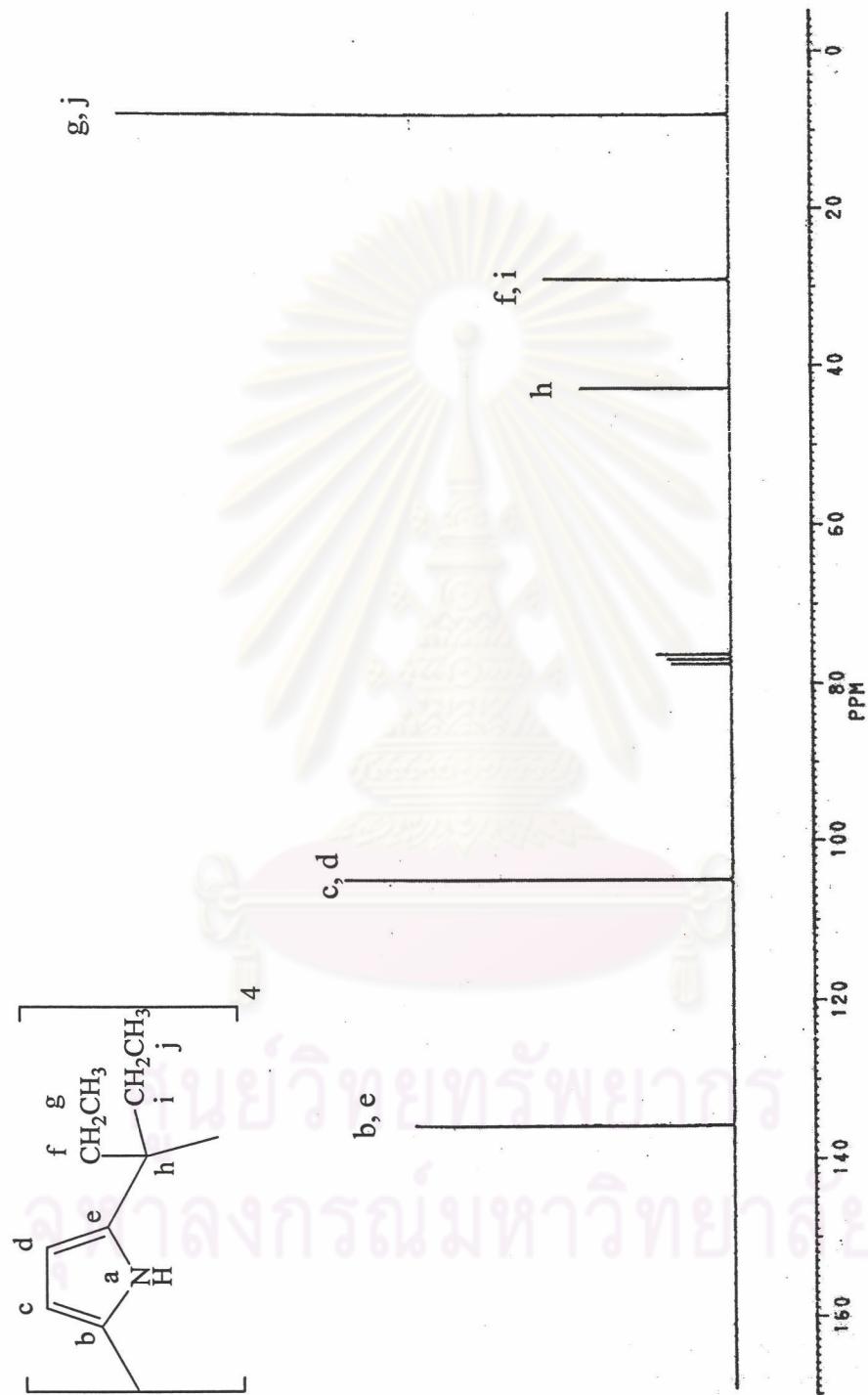


Figure 4 ^{13}C NMR spectrum of *meso*-octaethyl-calix[4]pyrrole

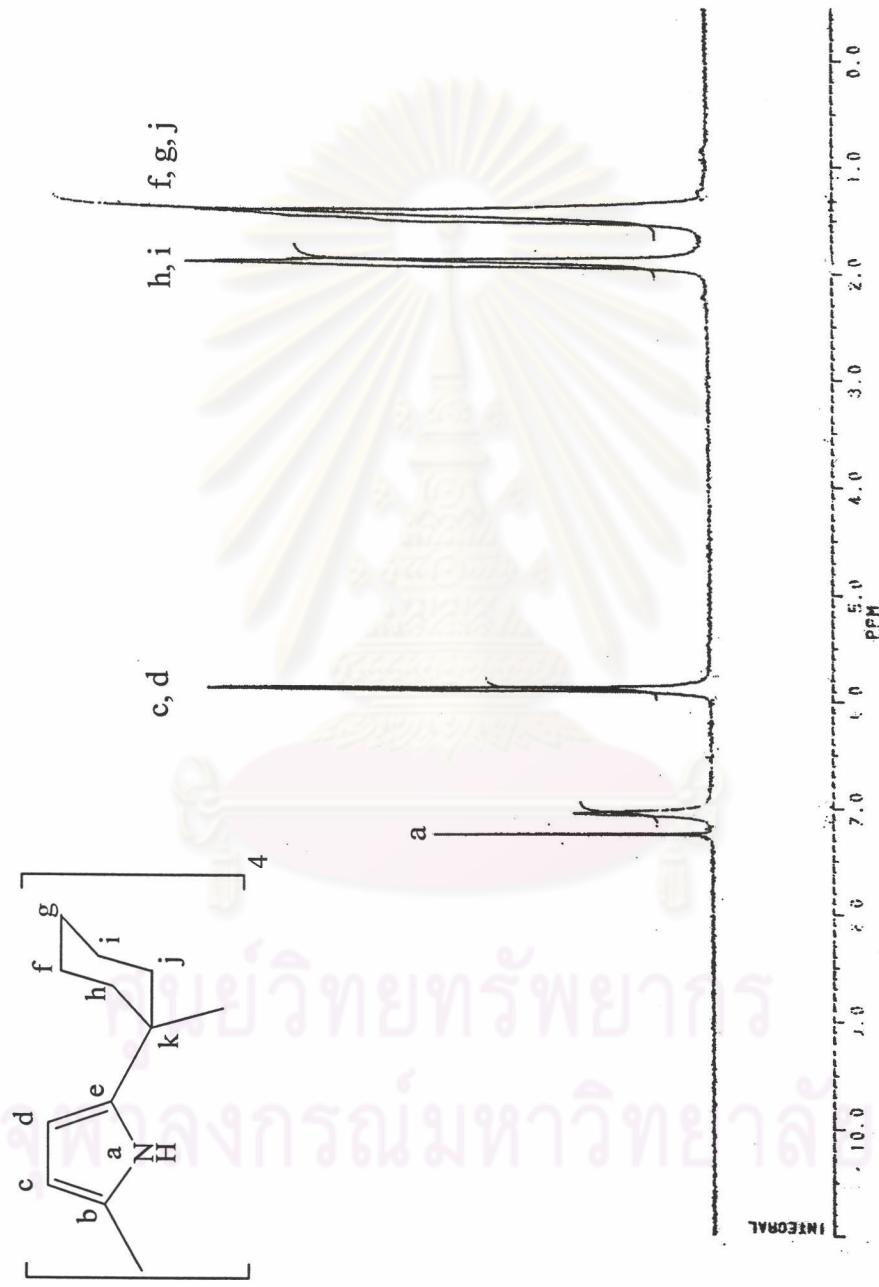


Figure 5 ^1H NMR spectrum of tetraspirohexyl-calix[4]pyrrole

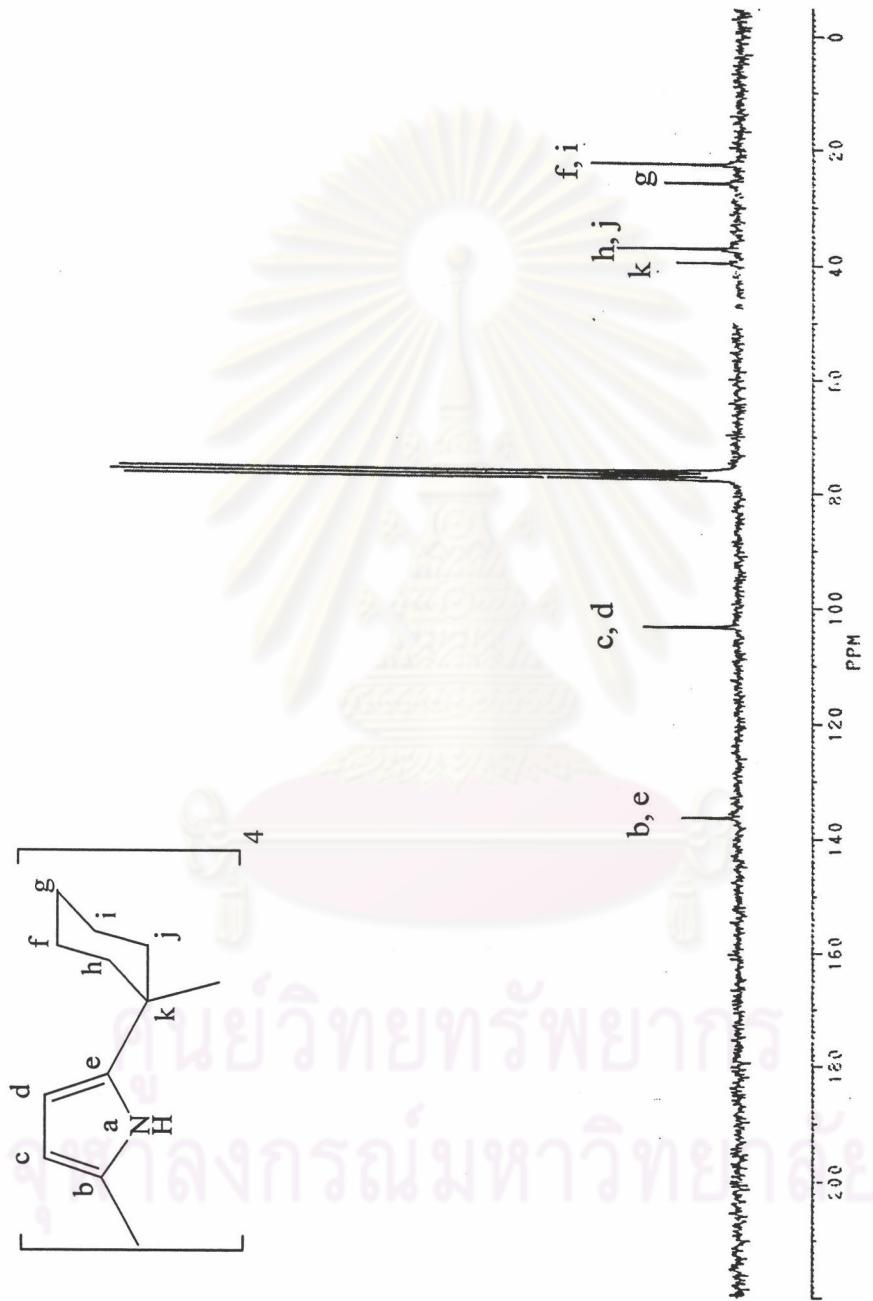


Figure 6 ^{13}C NMR spectrum of tetraspirohexyl-calix[4]pyrrole

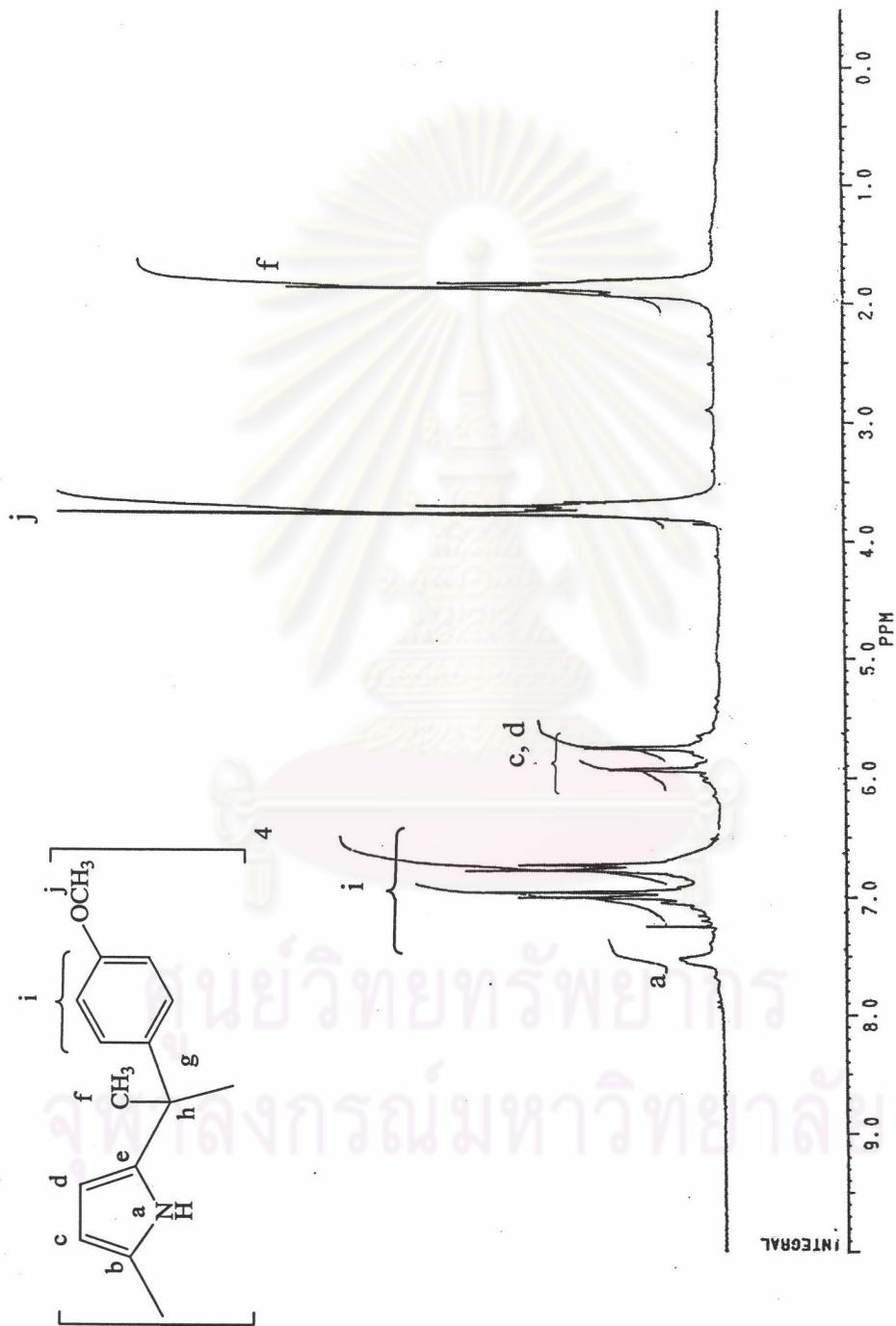


Figure 7 ^1H NMR spectrum of *meso*-tetrakis(4-methoxyphenyl)-tetramethyl-calix[4]pyrrole

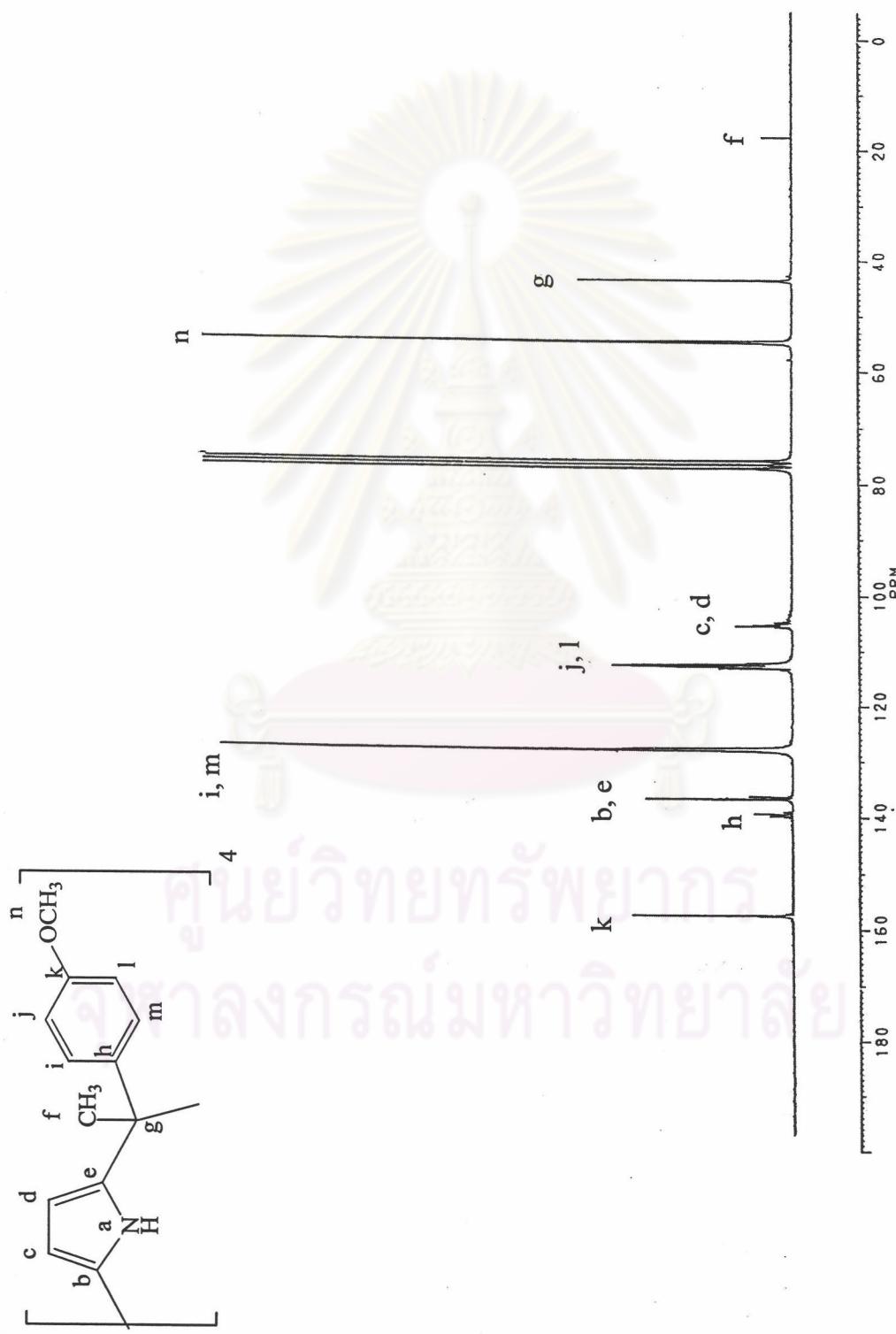


Figure 8 ^{13}C NMR spectrum of *meso*-tetrakis(4-methoxyphenyl)-tetramethyl-calix[4]pyrrole

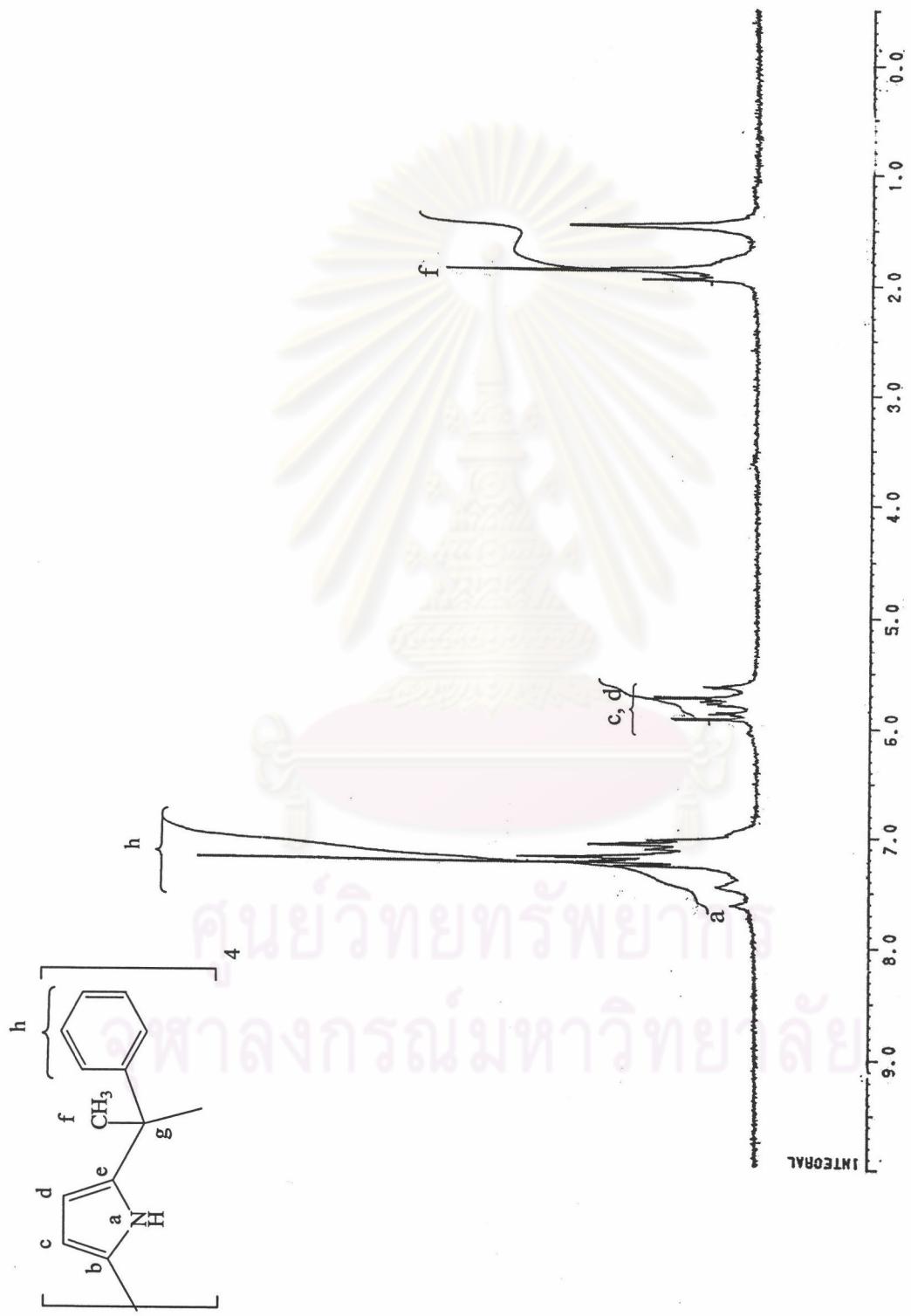


Figure 9 ^1H NMR spectrum of *meso*-tetramethyl-tetraphenyl-calix[4]pyrrole

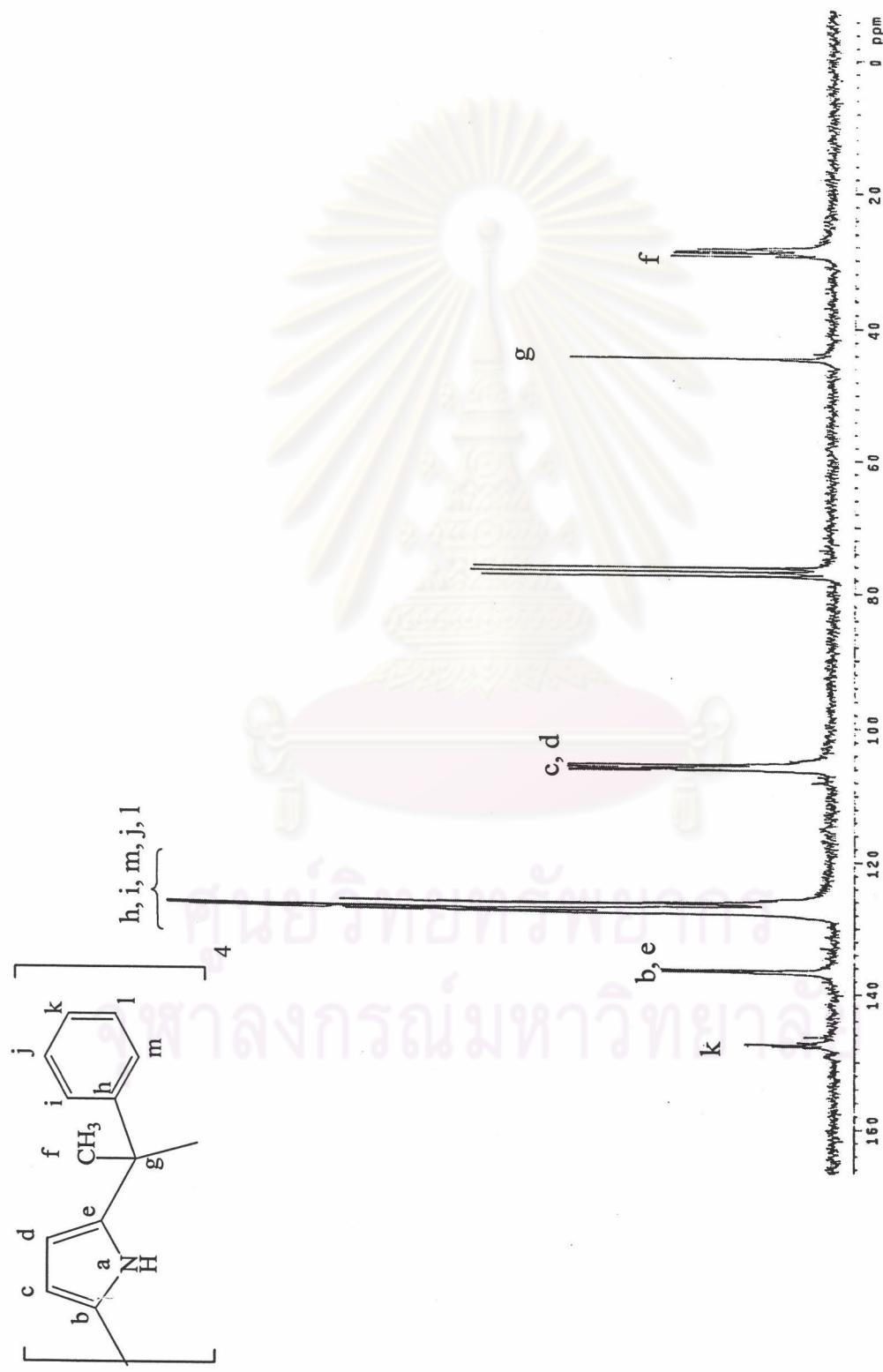


Figure 10 ^{13}C NMR spectrum of *meso*-tetramethyl-tetraphenyl-calix[4]pyrrole

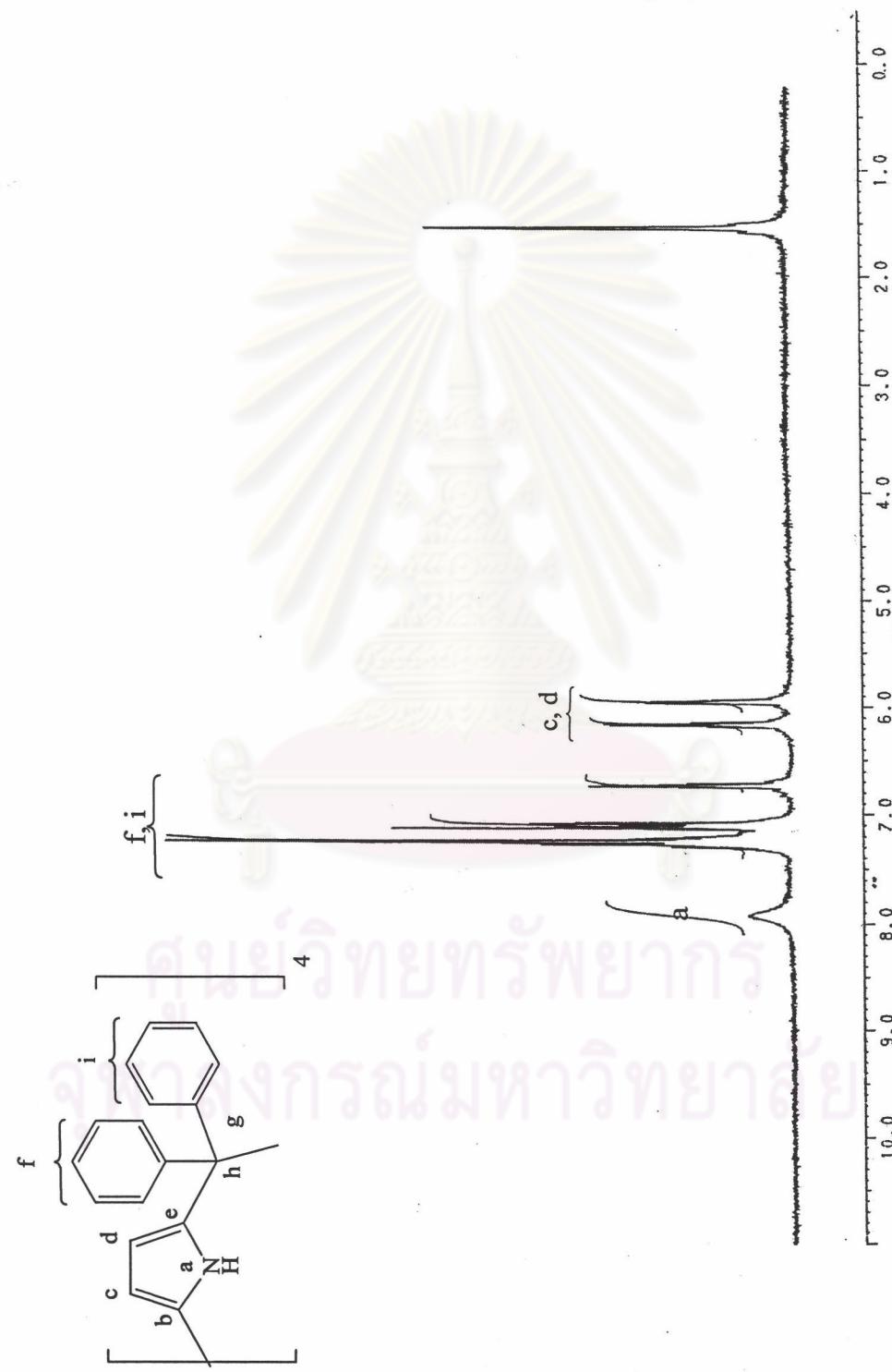


Figure 11 ^1H NMR spectrum of *meso*-octaphenyl-calix[4]pyrrole

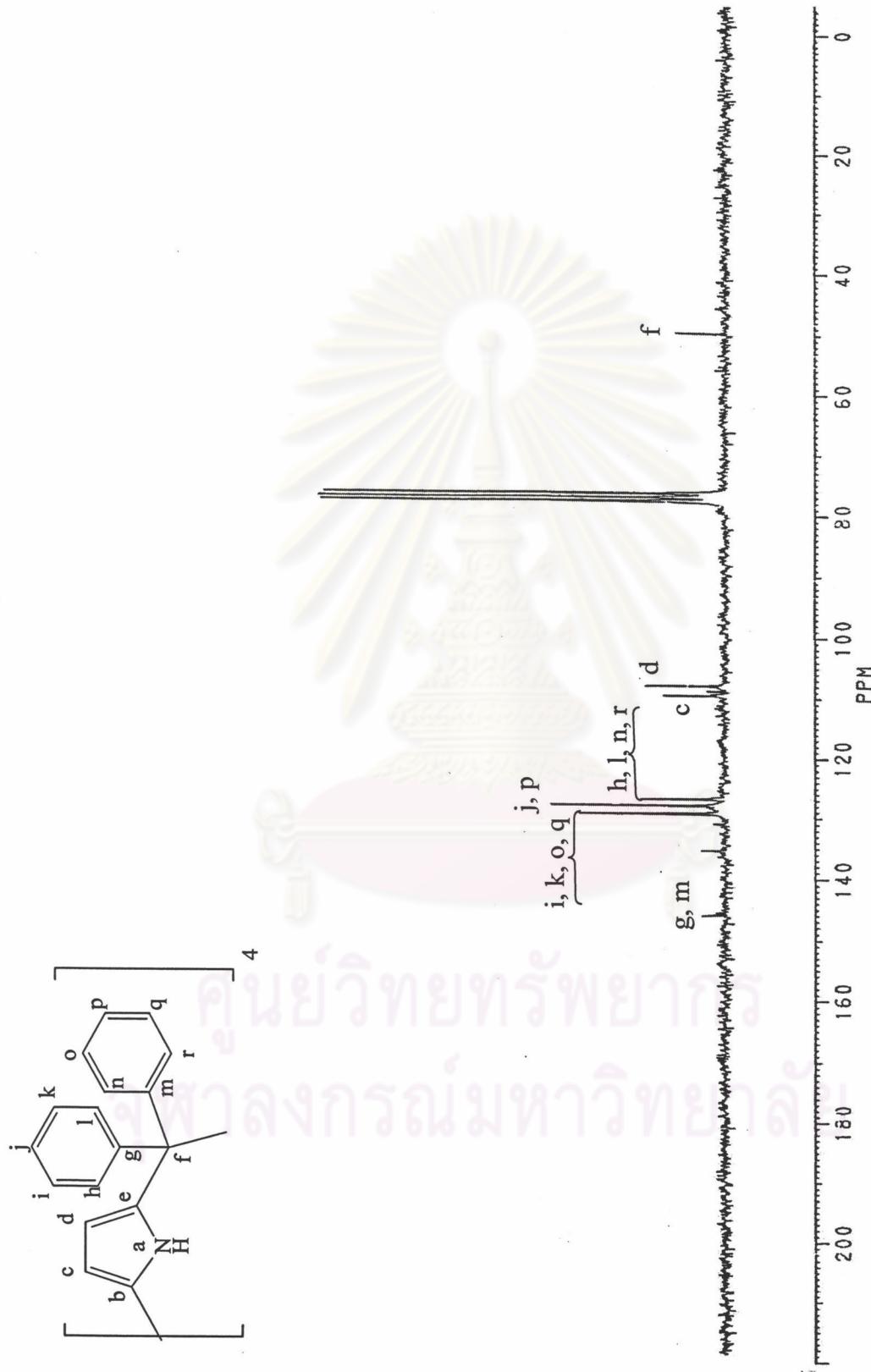


Figure 12 ^1H NMR spectrum of *meso*-octaphenyl-calix[4]pyrrole

VITA

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