

Supporting information for 2-{9-(10-Bromoanthracenyl)}-1,3-dihydro-1*H*-[*d*]-1,3,2-benzodiazaborole

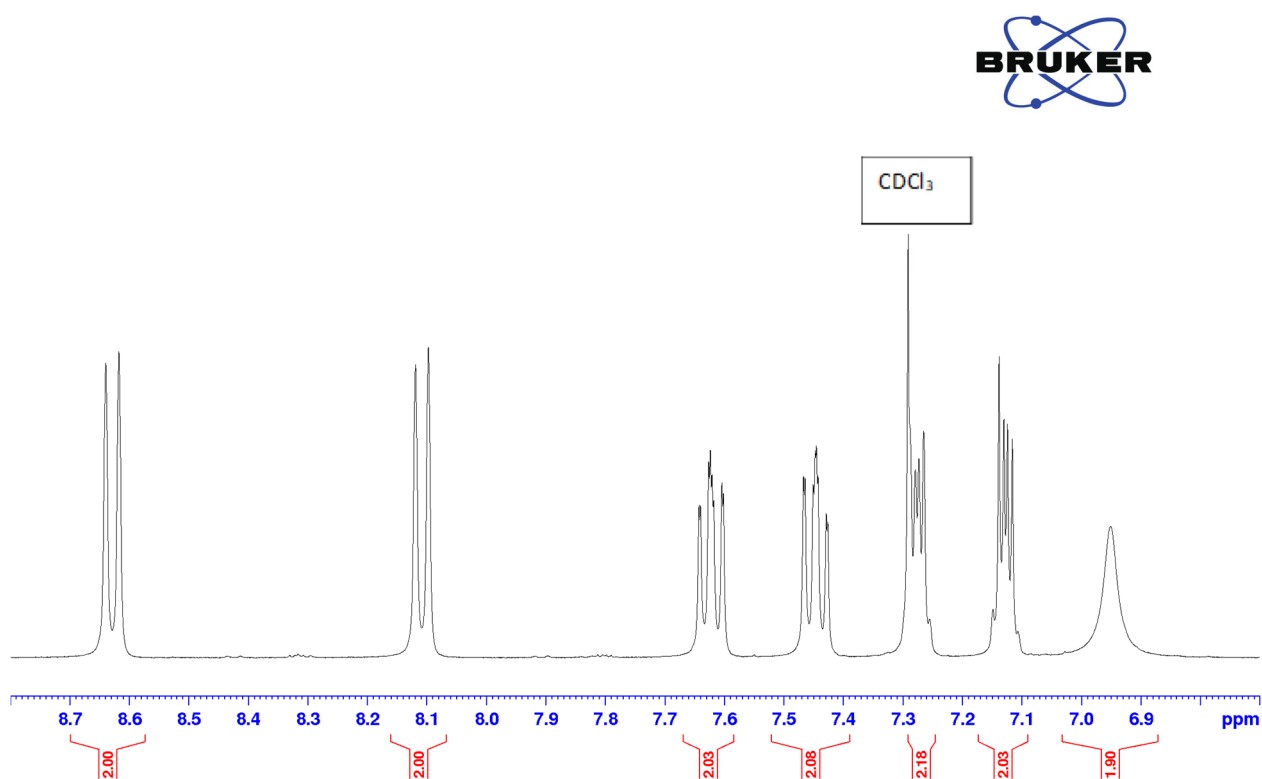
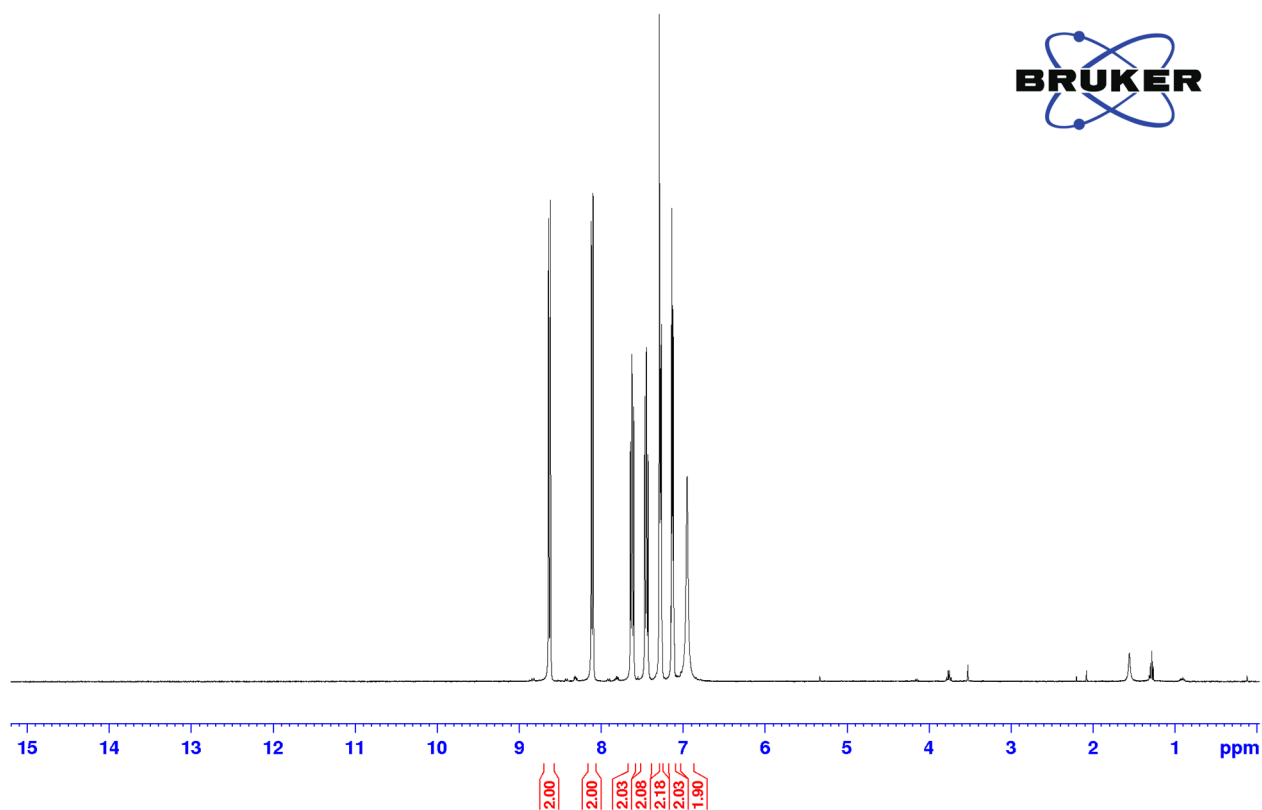


Figure S1. ¹H NMR spectrum of 2-{9-(10-Bromoanthracenyl)}-1,3-dihydro-1*H*-[*d*]-1,3,2-benzodiazaborole.

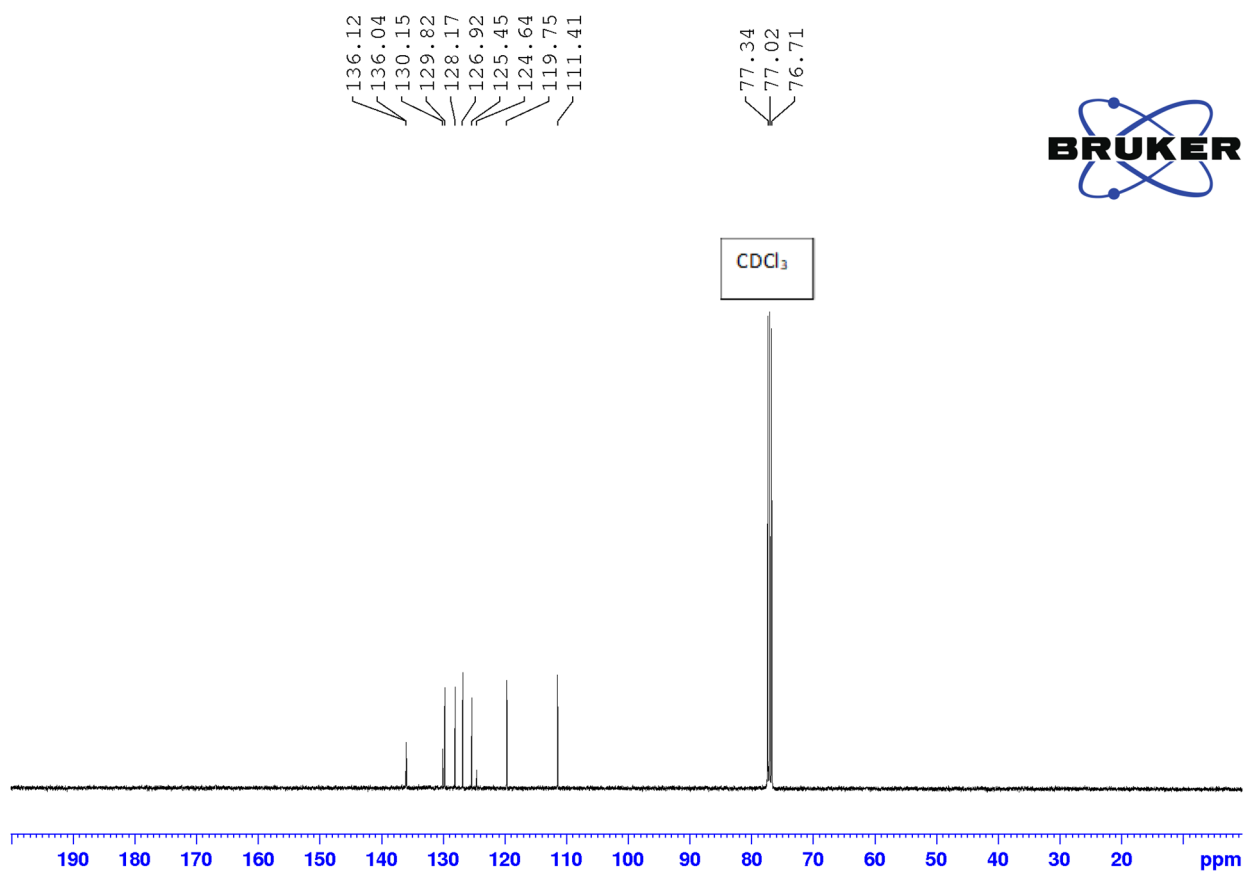


Figure S2. ¹³C NMR of 2-[9-(10-Bromoanthracenyl)]-1,3-dihydro-1H-[d]-1,3,2-benzodiazaborole.

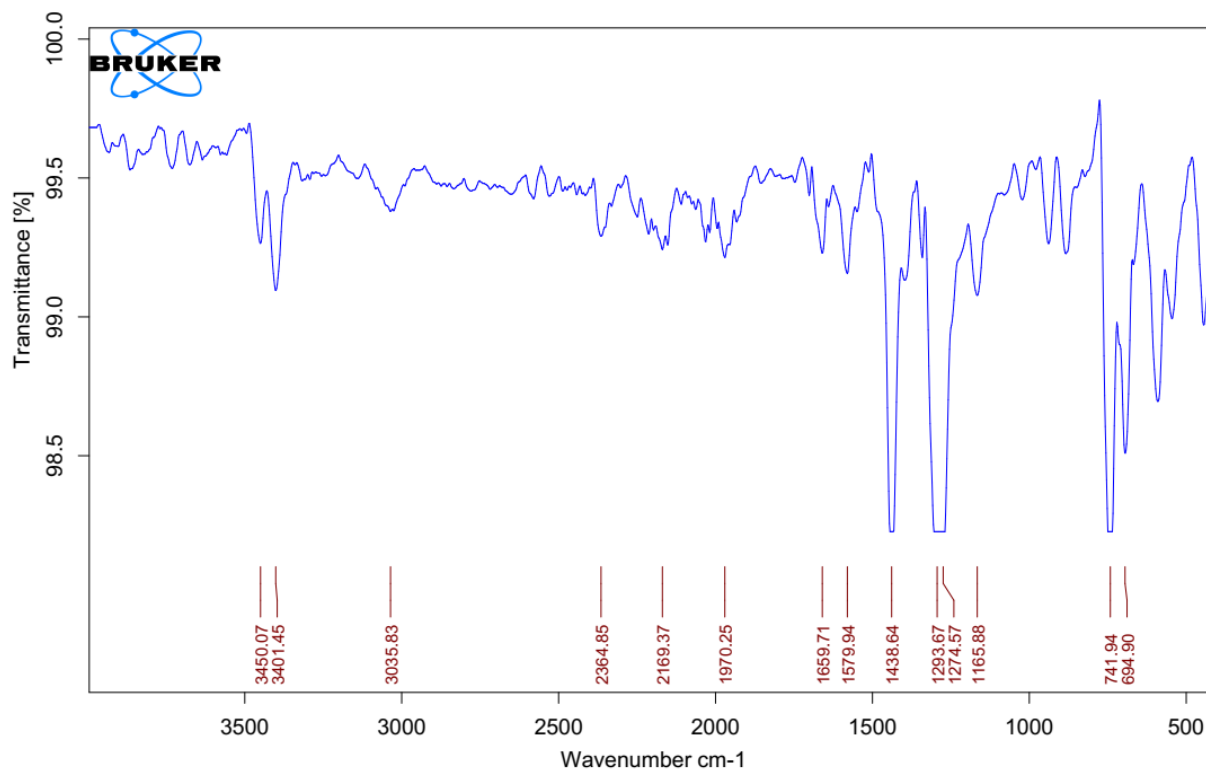


Figure S3. FT-IR of 2-[9-(10-Bromoanthracenyl)]-1,3-dihydro-1H-[d]-1,3,2-benzodiazaborole.

Single Mass Analysis

Tolerance = 5.0 PPM / DBE: min = -1.5, max = 100.0

Element prediction: Off

Number of isotope peaks used for i-FIT = 3

Monoisotopic Mass, Even Electron Ions

41 formula(e) evaluated with 1 results within limits (up to 20 closest results for each mass)

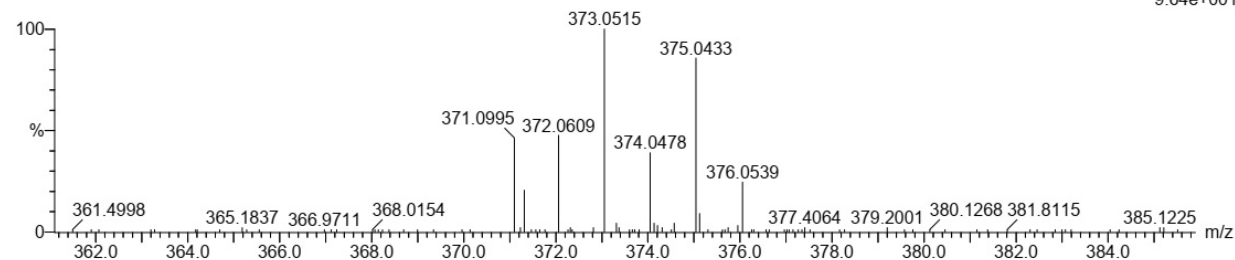
Elements Used:

C: 20-25 H: 10-15 N: 0-5 Br: 0-5 B: 0-1

diaza 1 29 (0.251)

TOF MS AP+

9.64e+001



Minimum:

Maximum:

-1.5

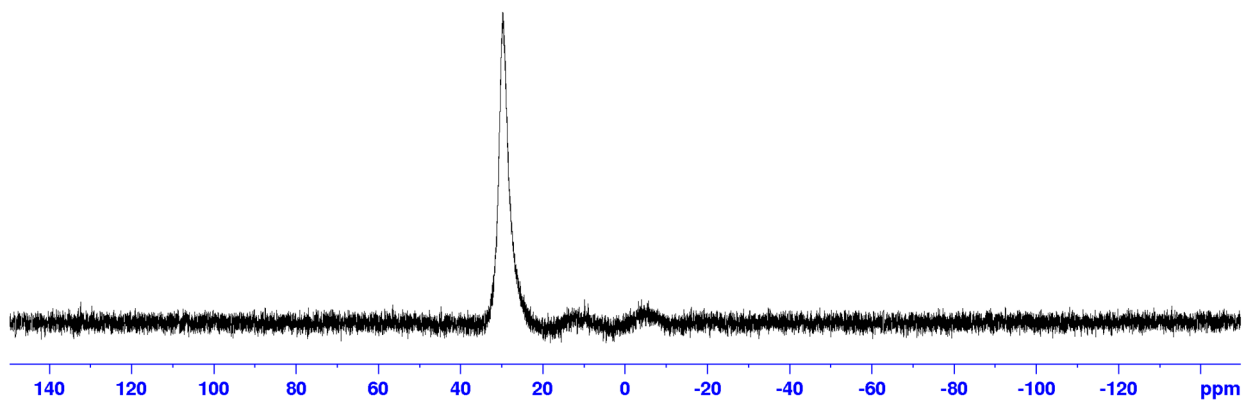
100.0

Mass	Calc. Mass	mDa	PPM	DBE	i-FIT	i-FIT (Norm)	Formula
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373.0515	373.0512	0.3	0.8	14.5	47.5	0.0	C20 H15 N2 Br B
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Figure S4. HRMS of 2-[9-(10-Bromoanthracenyl)]-1,3-dihydro-1H-[d]-1,3,2-benzodiazaborole.

— 29.64

Figure S5: ^{11}B NMR of 2-[9-(10-Bromoanthracenyl)]-1,3-dihydro-1H-[d]-1,3,2-benzodiazaborole.