

Electronic Supplementary Information

Synthetic, Spectral and Structural Studies of a Schiff Base and its Anticorrosive Activity on Mild Steel in H₂SO₄

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Table S1. Selected bond lengths (Å), and bond angles (°) for **L**.

Bond length (Å)		Bond Angle (°)	
O(1)- C(3)	1.235(3)	C(7)-C(8)-C(9)	120.5(3)
N(1)- C(1)	1.356(3)	C(6)-C(7)-C(8)	120.3(3)
N(1)- N(2)	1.409(3)	C(18)-C(17)-C(16)	121.9(3)
N(1) C(10)	1.463(3)	C(1)-N(1)-N(2)	106.9(2)
N(3)- C(12)	1.277(3)	C(1)-N(1)-C(10)	123.9(2)
N(3)- C(2)	1.396(3)	N(2)-N(1)-C(10)	117.4(2)
N(2)- C(3)	1.405(3)	C(12)-N(3)-C(2)	121.0(2)
N(2)- C(4)	1.423(3)	C(3)-N(2)-N(1)	108.8(2)
		C(3)-N(2)-C(4)	122.0(2)
		N(1)-N(2)-C(4)	119.3(2)
		N(1)-C(1)-C(2)	110.5(2)
		N(1)-C(1)-C(11)	121.7(2)
		C(2)-C(1)-C(11)	127.8(3)
		C(1)-C(2)-N(3)	122.2(2)
		C(1)-C(2)-C(3)	108.0(2)
		N(3)-C(2)-C(3)	129.5(2)
		O(1)-C(3)-N(2)	122.9(2)
		O(1)-C(3)-C(2)	132.0(3)
		N(2)-C(3)-C(2)	105.0(2)
		C(9)-C(4)-C(5)	120.1(3)
		C(9)-C(4)-N(2)	121.5(2)
		C(5)-C(4)-N(2)	118.3(3)
		N(3)-C(12)-C(13)	121.5(3)
		C(6)-C(5)-C(4)	119.4(3)
		C(18)-C(13)-C(14)	117.6(3)
		C(18)-C(13)-C(12)	122.1(3)
		C(14)-C(13)-C(12)	120.3(3)
		C(15)-C(14)-C(13)	121.7(3)
		C(4)-C(9)-C(8)	119.4(3)
		C(16)-C(15)-C(14)	120.4(3)
		C(15)-C(16)-C(17)	118.0(3)
		C(15)-C(16)-C(19)	121.1(3)
		C(17)-C(16)-C(19)	120.9(3)
		C(7)-C(6)-C(5)	120.3(3)
		C(17)-C(18)-C(13)	120.5(3)

Table S2. Hydrogen bond parameters for **L**.

D-H···A-X	<i>d</i> H···A Å	<i>D</i> D···A Å	θ D-H···A°
1			
C(12)-H(12)··· O(1)	2.37(3)	3.047(4)	125(2)
C(15)-H(15)··· O(1) ^a	2.48	3.313(4)	149
symmetry equivalents: (a)2-X,-Y,-Z.			

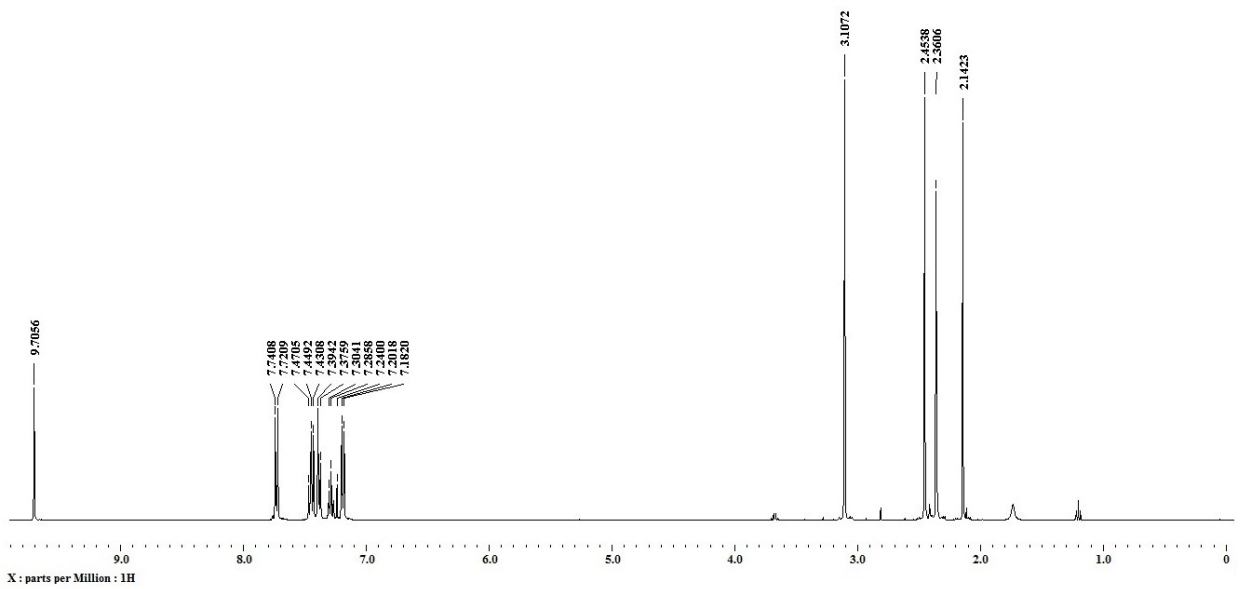


Figure F1. ^1H NMR spectrum of L.

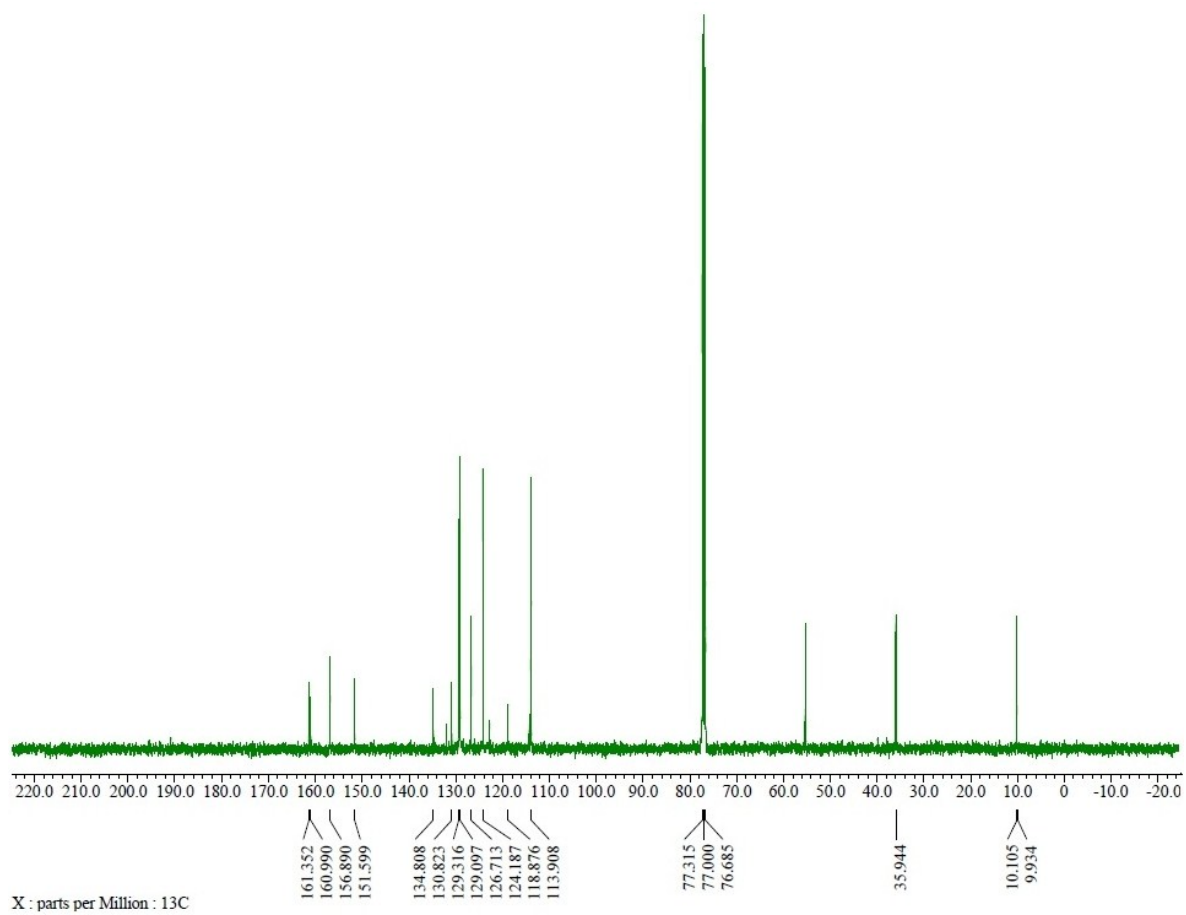


Figure F2. ^{13}C NMR spectrum of L.

Qualitative Compound Report

Data File	AK-schiff.d	Sample Name	AK-schiff
Sample Type	Sample	Position	P1-B4
Instrument Name	Instrument 1	User Name	
Acq Method	29.10.2014.m	Acquired Time	11-05-2017 13:25:16
IRM Calibration Status	Success	DA Method	Default.m
Comment			

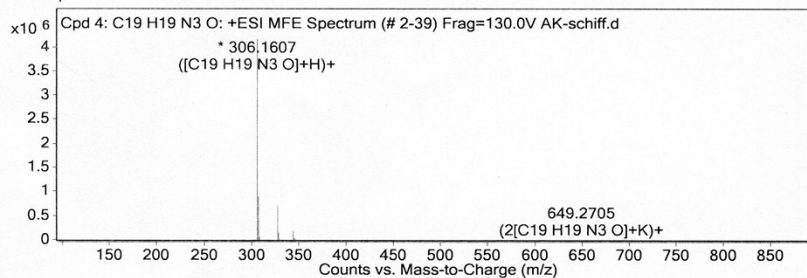
Sample Group	Info.
Acquisition SW	6200 series TOF/6500 series
Version	Q-TOF B.05.01 (B5125)

Compound Table

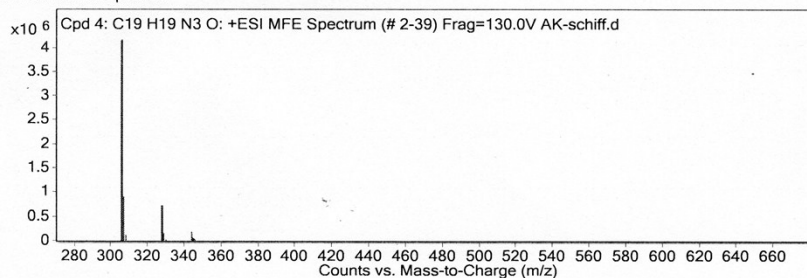
Compound Label	RT	Mass	Formula	MFG Formula	MFG Diff (ppm)	DB Formula
Cpd 4: C19 H19 N3 O	11	305.1534	C19 H19 N3 O	C19 H19 N3 O	-2.09	C19 H19 N3 O

Compound Label	m/z	RT	Algorithm	Mass
Cpd 4: C19 H19 N3 O	306.1607	11	Find by Molecular Feature	305.1534

MFE MS Spectrum



MFE MS Zoomed Spectrum



MS Spectrum Peak List

m/z	z	Abund	Formula	Ion
306.1607	1	4162072	C19 H19 N3 O	(M+H)+
307.1642	1	900961.33	C19 H19 N3 O	(M+H)+
308.1668	1	96179.13	C19 H19 N3 O	(M+H)+
309.1688	1	7836.04	C19 H19 N3 O	(M+H)+
328.1429	1	718592.69	C19 H19 N3 O	(M+Na)+
329.1457	1	145851.92	C19 H19 N3 O	(M+Na)+
330.1487	1	16339.05	C19 H19 N3 O	(M+Na)+
344.1166	1	185627.09	C19 H19 N3 O	(M+K)+
345.1195	1	40311.35	C19 H19 N3 O	(M+K)+
346.1169	1	16867.37	C19 H19 N3 O	(M+K)+

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