

*Supplementary:*

## ***N*<sup>1</sup>-[4-(1-methyl-1*H*-imidazol-2-yl)-[2,4'-bipyridin]-2'-yl]benzene-1,4-diamine**

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Figure S1: HPLC of title compound.

Figure S2: Mass spectra of title compound.

Figure S3: CNMR of title compound.

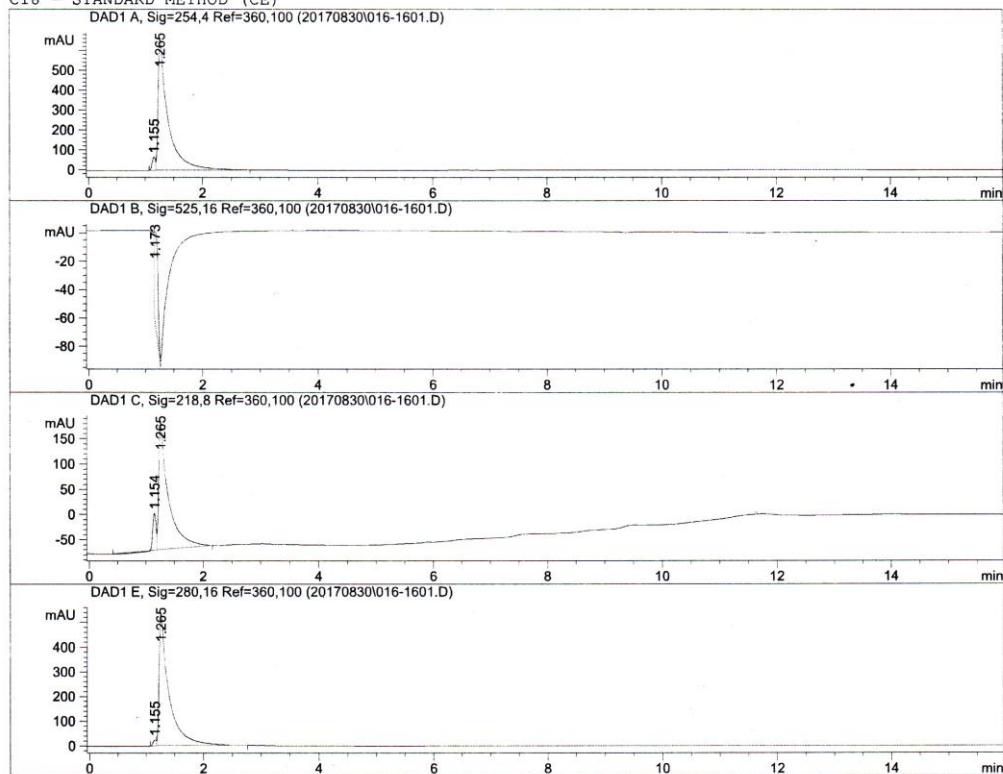
Figure S4: DEPT of title compound.

Figure S5: HNMR of title compound.

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Injection Date	: 30.08.2017 19:13:02	Seq. Line	: 16
Sample Name	: DZ-OX-7	Location	: Vial 16
Acq. Operator	: ME	Inj	: 1
		Inj Volume	: 10 µl

Acq. Method : C:\HPCHEM\1\METHODS\C18STD.M  
Last changed : 30.08.2017 19:11:57 by ME  
(modified after loading)  
Analysis Method : C:\HPCHEM\1\METHODS\C18STD.M  
Last changed : 01.02.2017 18:31:53 by CE  
C18 - STANDARD METHOD (CE)



=====  
Area Percent Report  
=====

Sorted By : Signal  
Multiplier : 1.0000  
Dilution : 1.0000  
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	1.155	PV	0.0572	247.31012	66.80379	3.0246
2	1.265	VB	0.1600	7929.44727	660.77563	96.9754

Totals : 8176.75739 727.57943

Results obtained with enhanced integrator!

Signal 2: DAD1 B, Sig=525,16 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	1.173	BP	0.0740	293.88531	63.44610	100.0000

Totals : 293.88531 63.44610

Results obtained with enhanced integrator!

Signal 3: DAD1 C, Sig=218,8 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	1.154	PV	0.0589	270.23532	73.54895	8.3000
2	1.265	VB	0.1573	2985.62646	253.74661	91.7000

Totals : 3255.86179 327.29556

Results obtained with enhanced integrator!

Signal 4: DAD1 E, Sig=280,16 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	1.155	PV	0.0532	89.88020	25.43465	1.3687
2	1.265	VB	0.1610	6477.16748	535.97656	98.6313

Totals : 6567.04768 561.41121

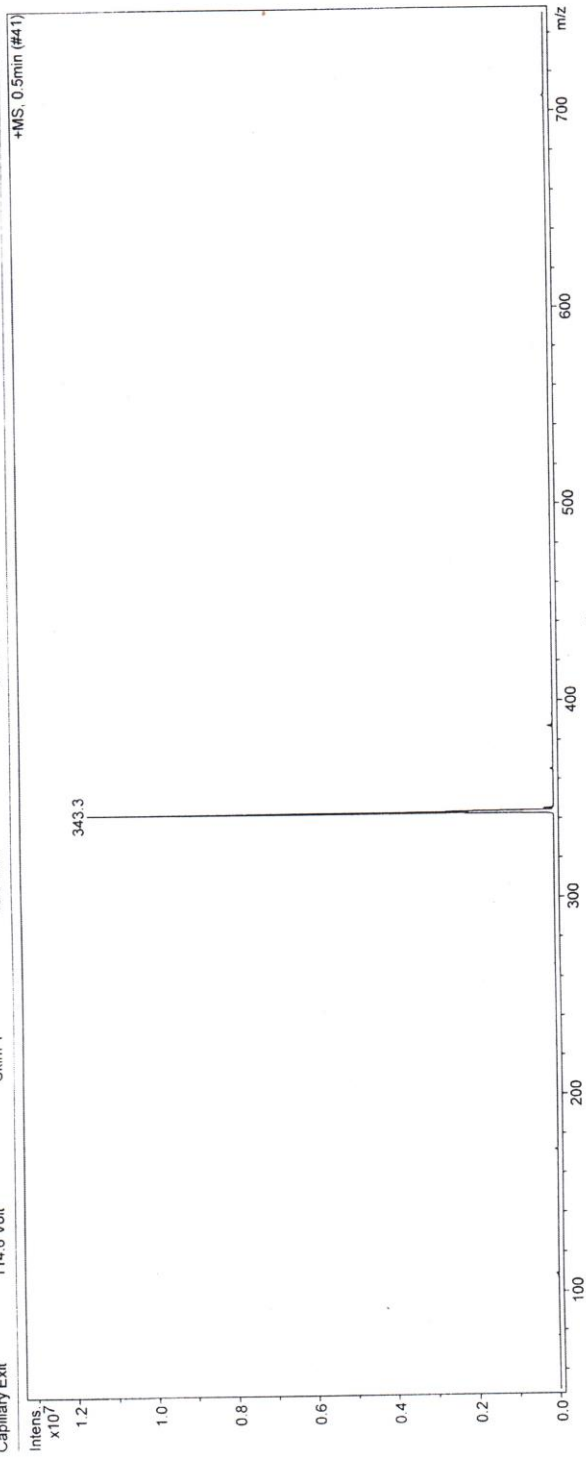
Results obtained with enhanced integrator!

\*\*\* End of Report \*\*\*

Figure S1: HPLC of title compound.

# Display Report

<b>Analysis Info</b>		ZINAD011.D		08/31/17 18:58:48		Operator		Default	
Analysis Name		DZ-OX-7		Kopfe von FIA M		Instrument		esquire3000plus_01038	
Sample Name		Dhafer Zinad, AG Koch		Method					
Comment									
<b>Acquisition Parameter</b>		Mass Range Mode		Std/Normal		Ion Polarity		Positive	
Ion Source Type		ESI		750 m/z		Averages		8 Spectra	
Scan Begin		50 m/z		40.0 Volt		Trap Drive		40.9	
Capillary Exit		114.5 Volt		Scan 1		Auto MS/MS		off	
						Alternating Ion Polarity		2528 $\mu$ s	
						Accumulation Time		off	



## Display Report

<b>Analysis Info</b>			
Analysis Name	ZINAD011 D	Acquisition Date	08/31/17 18:58:48
Sample Name	DZ-OX-7	Method	Kopie von FIA.M
Comment	Dhafer Zinad, AG Koch	Operator	Default
		Instrument	esquire3000plus_01038
<b>Acquisition Parameter</b>			
Ion Source Type	ESI	Mass Range Mode	Std/Normal
Scan Begin	50 m/z	Scan End	750 m/z
Capillary Exit	114.6 Volt	Skim 1	40.0 Volt
		Ion Polarity	Positive
		Averages	8 Spectra
		Trap Drive	40.9
		Alternating Ion Polarity	off
		Accumulation Time	2528 $\mu$ s
		Auto MS/MS	off

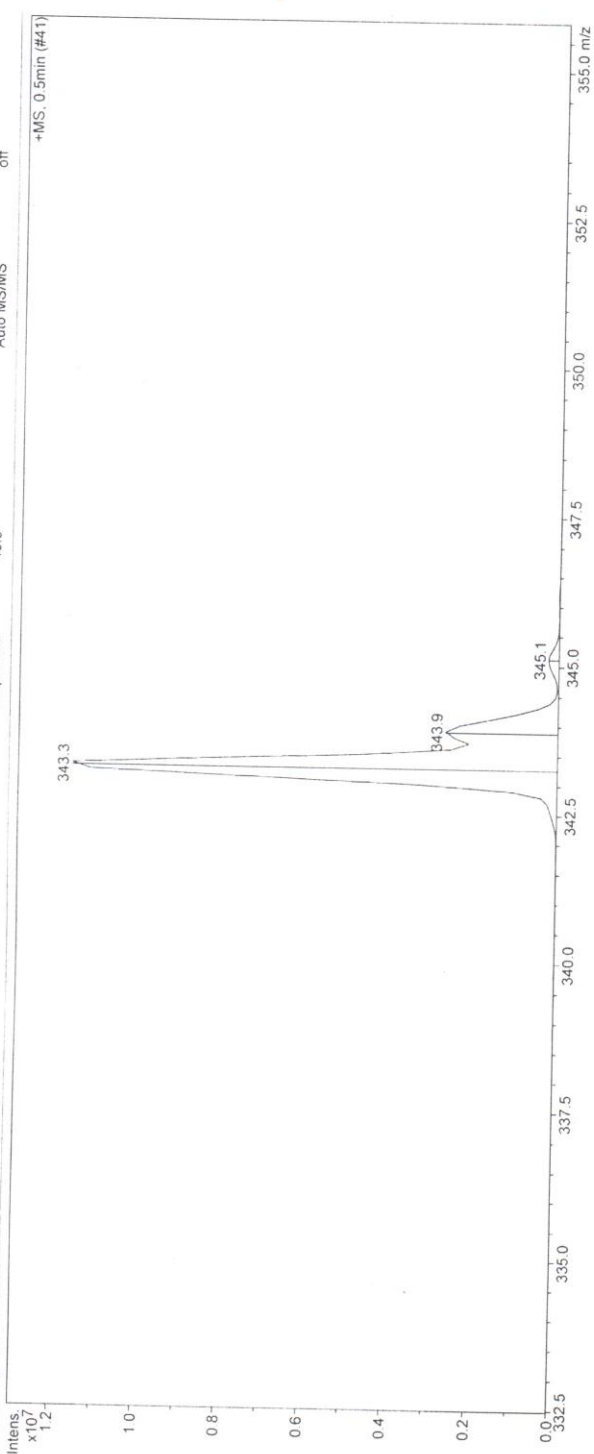
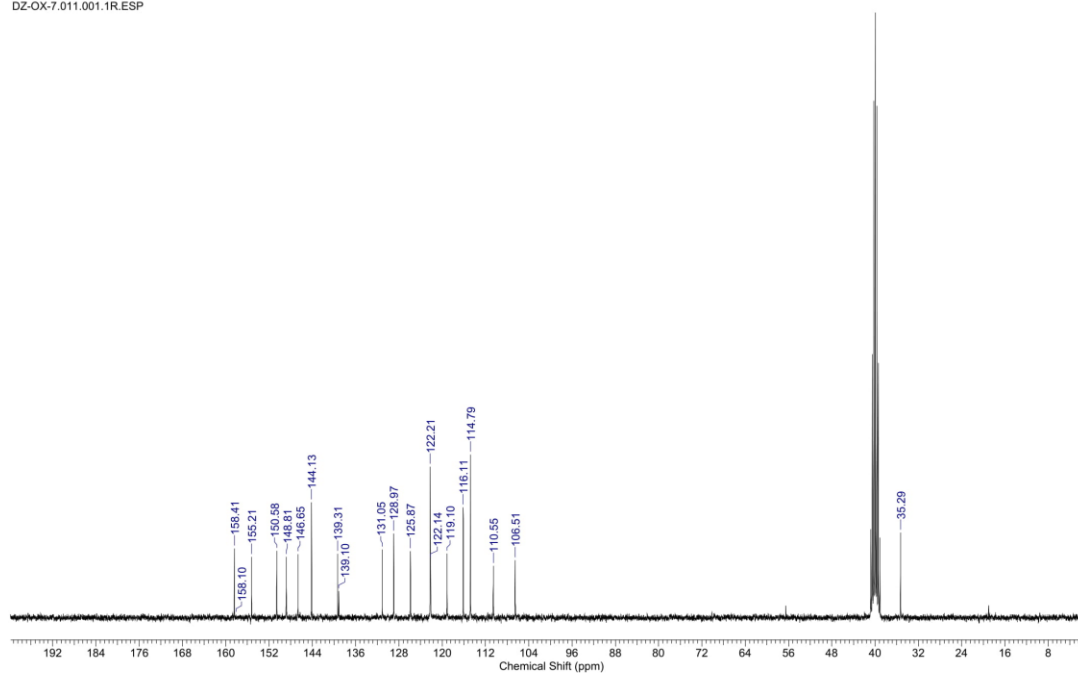


Figure S2: Mass spectra of title compound.

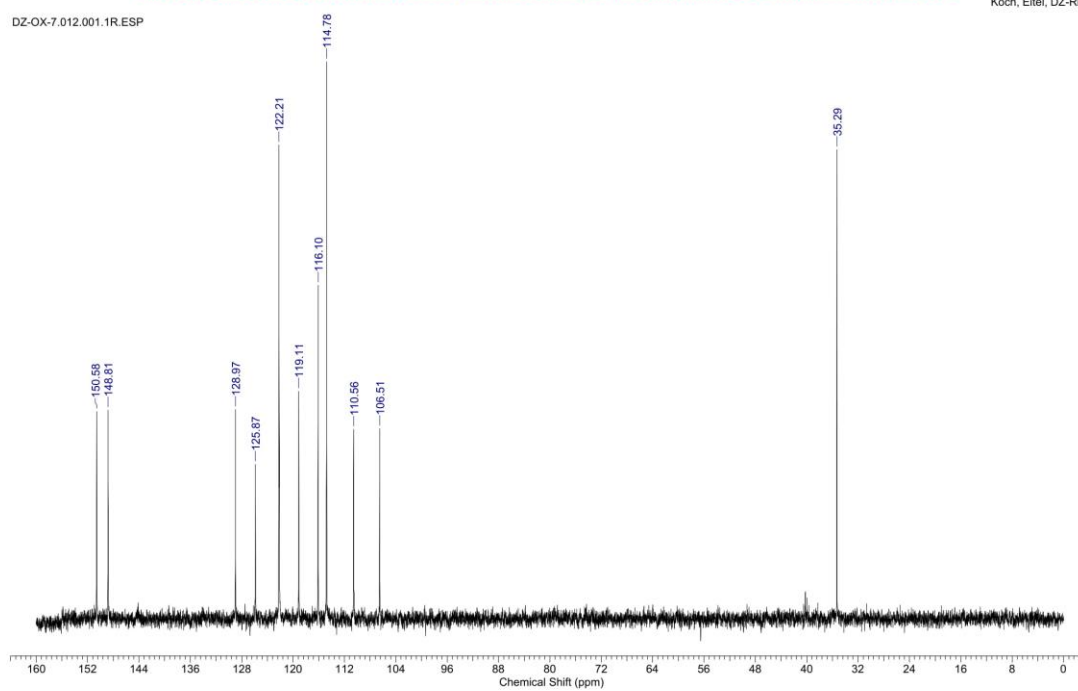
DZ-OX-7.011.001.1R.ESP



C:\USERS\SONY\DESKTOP\IDZ-OX-7\IDZ-OX-7.011.001.1R.ESP

Figure S3: CNMR of title compound.

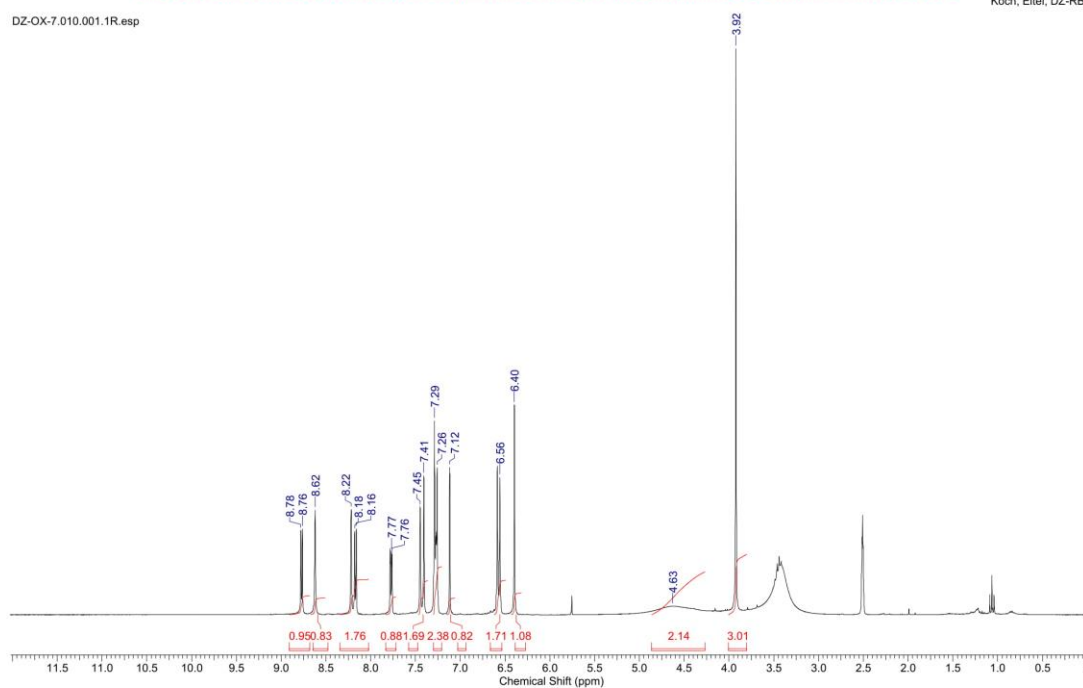
DZ-OX-7.012.001.1R.ESP



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Figure S4: DEPT of title compound.

DZ-OX-7.010.001.1R.esp



C:\USERS\SONY\DESKTOP\SPECTRA\IDZ-OX-7\IDZ-OX-7.010.001.1R.esp

Figure S5: <sup>1</sup>H NMR of title compound.