## **Electronic Supporting Information**

# Title: *p*-TSA·H<sub>2</sub>O catalyzed metal-free and environmentally benign synthesis of 4-aryl quinolines from Arylamine, Arylacetylene, and Dimethyl sulfoxide

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<sup>1</sup>H NMR Spectrum of 6-Methoxy-4-phenylquinoline(3a)





#### HRMS Spectrum of 6-Methoxy-4-phenylquinoline(3a)





<sup>13</sup>C NMR Spectrum of 6-Methoxy-4-(m-*tolyl*)quinoline (3b)

## HRMS Spectrum of 6-Methoxy-4-(m-tolyl)quinoline (3b)



 <sup>1</sup>H NMR Spectrum of 6-Methoxy-4-(p-tolyl)quinoline(3c)

 SF-PA-PE-DMSO-1H.1.fid – 1H

 6
 8
 8
 9
 1
 1
 1

 6
 8
 8
 9
 1
 1
 1
 1

 8
 8
 8
 9
 1
 1
 1
 1
 1

 8
 8
 9
 1
 1
 1
 1
 1
 1

 9
 8
 9
 1
 1
 1
 1
 1
 1

 9
 8
 9
 1
 1
 1
 1
 1
 1







MeO

# <sup>13</sup>C NMR Spectrum of 6-Methoxy-4-(*p*-tolyl)quinoline(3c)

SF-PA-PE-DMSO-13C.3.fid — 13C

157.95 147.59 147.59 144.83 138.39 135.51 131.25 123.35 129.33 127.95 121.90 121.80 103.88	77.37 77.16 76.95	55.57	21.44
	$\checkmark$		





# HRMS Spectrum of 6-Methoxy-4-(p-tolyl)quinoline(3c)



<sup>1</sup>H NMR Spectrum of 4-(4-Ethylphenyl)-6-methoxyquinoline(3d) SF-PA-4EE-T-1H.1.fid — SF-PA-4EE-T-1H 8.08 8.06 7.45 7.45 7.39 7.39 7.35 7.35 7.35 7.35 7.25 7.25 7.25 8.78 8.77 2.79 2.77 2.76 2.74 1.35 1.33 1.32 3.79 14 2 7.45 7.43 7.39 7.38 7.35 7.35 7.27 7.27 7.25 7.25 7.25 8.78 8.77 8.08 8.06 MeO 8.9 8.8 8.7 8.6 8.5 8.4 8.3 8.2 8.1 8.0 7.9 7.8 7.7 7.6 7.5 7.4 7.3 7.2 7.1 7.0 f1 (ppm) **1.00**-1.00H 1.99 3.01 1.00 1.00 3.06₌ 2.03<sub>1</sub> **3.05**∃ 5.0 f1 (ppm) 9.5 9.0 8.0 7.5 7.0 6.0 5.5 4.5 4.0 3.5 3.0 2.5 2.0 1.5 L0.0 8.5 6.5 1.0 0.5 0.0







 $\mathsf{SF}\text{-}\mathsf{PT}\text{-}\mathsf{3}\mathsf{D}\mathsf{M}\mathsf{S}\text{-}\mathsf{1}\mathsf{H}\text{.}\mathsf{1}\text{.}\mathsf{fid} - \mathsf{1}\mathsf{H}$ 8.81 8.80 8.12 8.10 7.57 7.57 7.57 7.57 7.49 7.49 7.40 7.40 7.40 7.40 7.40 7.31 7.32 7.33 3.83 1.44 $\sqrt{-}$ 1 8.12 8.10 7.57 7.57 7.50 7.50 7.42 7.42 7.42 7.40 7.40 7.40 7.31 7.31 7.33 7.33 トイ MeO 7.7 7.6 7.5 7.4 7.3 7.2 f1 (ppm) 8.2 8.1 8.0 7.9 7.8  $1.00_{\pm}$ 0.98₌ 2.03 2.03 0.99<sup>₹</sup> 2.03 3.01₌ 9.09₌ 5.0 4.5 f1 (ppm) 10.0 9.5 6.5 6.0 5.5 4.0 3.5 1.5 9.0 8.5 8.0 7.5 7.0 3.0 2.5 2.0 1.0 0.5 0.0

<sup>1</sup>H NMR Spectrum of 4-(4-(*tert*-Butyl)phenyl)-6-methoxyquinoline (3e)

34.81 31.45 f1 (ppm) 



#### HRMS Spectrum of 4-(4-(*tert*-Butyl)phenyl)-6-methoxyquinoline(3e)

<sup>1</sup>H NMR Spectrum of 6-Methoxy-4-(4-methoxyphenyl)quinoline(3f)

SF-PA-4FE-ST-1H.1.fid —	SF-PA-4FE-ST-1H	J
8.77 8.76	8.05 8.05 7.45 7.45 7.39 7.38 7.37 7.37 7.37 7.37 7.37 7.37 7.25 7.25 7.25 7.07 7.05	3.90
$\checkmark$		$\mathbf{X}$



# <sup>13</sup>C NMR Spectrum of 6-Methoxy-4-(4-methoxyphenyl)quinoline(3f)

011				
SF-PA-4FE-ST-13C.3.fid — SF-PA-4FE-	ST-13C	-		
159.92 157.97	147.67 147.01 144.98	131.37 130.67 128.06 121.80 121.76 121.76 116.57 114.29 103.94	77.41 77.16 76.91 55.58	55.51
$\sim$	$\searrow$ ///		$\checkmark \checkmark \checkmark$	/



#### HRMS Spectrum of 6-Methoxy-4-(4-methoxyphenyl)quinoline(3f)









<sup>19</sup>F NMR Spectrum of 4-(4-Fluorophenyl)-6-methoxyquinoline(3g)



#### HRMS Spectrum of 4-(4-Fluorophenyl)-6-methoxyquinoline(3g)

<sup>1</sup>H NMR Spectrum of 4-(4-Bromophenyl)-6-methoxyquinoline(3h)





## HRMS Spectrum of 4-(4-Bromophenyl)-6-methoxyquinoline(3h)







<sup>19</sup> FNMR Spectrum of 6-Methoxy-4-(2-(trifluoromethyl)phenyl)quinoline(3i).



10 0 -10 -20 -30 -40 -50 -60 -70 -80 -90 -100 -110 -120 -130 -140 -150 -160 -170 -180 -190 -200 -210 f1 (ppm)

#### HRMS Spectrum of 6-Methoxy-4-(2-(trifluoromethyl)phenyl)quinoline(3i)





<sup>1</sup>H NMR Spectrum of 4-(3-Ethynylphenyl)-6-methoxyquinoline (3j)







<sup>1</sup>H NMR Spectrum of 6-Methoxy-4-(pyridin-2-yl)quinoline (3k)

MeO

SF-PA-PYR-DMS-1H.1.fid - 1H





8.8 8.7 8.6 8.5 8.4 8.3 8.2 8.1 8.0 7.9 7.8 7.7 7.6 7.5 7.4 7.3 7.2 7.1 f1 (ppm)



<sup>13</sup>C NMR Spectrum of 6-Methoxy-4-(pyridin-2-yl)quinoline (3k)



## HRMS Spectrum of 6-Methoxy-4-(pyridin-2-yl)quinoline (3k)








# HRMS Spectrum of 6-Methoxy-4-(thiophen-3-yl)quinoline(3l)





## HRMS Spectrum of 6-Methoxy-3,4-diphenylquinoline(3m)







## HRMS Spectrum of 4-(4-(*tert*-Butyl)phenyl)-6-(methylthio)quinoline (3n)





<sup>1</sup>H NMR Spectrum of 4-Phenylquinoline(30)



## HRMS Spectrum of 4-Phenylquinoline(30)









<sup>1</sup>H–<sup>1</sup>H- COSY Spectrum of 6-Methyl-4-phenylquinoline (3p) SF-PT-DM-4-2\_COSY.17.ser – SF-PT-DM-4-2\_COSY

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<sup>1</sup>H–<sup>13</sup>C- COSY Spectrum of 6-Methyl-4-phenylquinoline (3p)

SF-PT-DM-4-2\_HSQC.15.ser — SF-PT-DM-4-2\_HSQC



# HRMS Spectrum of 6-Methyl-4-phenylquinoline (3p)





<sup>1</sup>H NMR Spectrum of 7-Methoxy-4-phenylquinoline (3q)





# HRMS Spectrum of 7-Methoxy-4-phenylquinoline (3q)







## HRMS Spectrum of 6-(Methylthio)-4-phenylquinoline (3r)



<sup>1</sup>H NMR Spectrum of 6,7-Dimethyl-4-phenylquinoline (3s)



<sup>13</sup>C NMR Spectrum of 6,7-Dimethyl-4-phenylquinoline (3s)

# HRMS Spectrum of 6,7-Dimethyl-4-phenylquinoline (3s)





<sup>1</sup>H NMR Spectrum of 4-Phenylbenzo[*h*]quinoline(3t)



<sup>13</sup>C NMR Spectrum of 4-Phenylbenzo[*h*]quinoline(3t)

# HRMS Spectrum of 4-Phenylbenzo[*h*]quinoline(3t)





<sup>1</sup>H NMR Spectrum of 4-Phenyl-6*H*-indeno[2,1-*g*]quinoline (3u)



## HRMS Spectrum of 4-Phenyl-6*H*-indeno[2,1-*g*]quinoline (3u)



<sup>1</sup>H NMR Spectrum of 4-Phenyl-1,10-phenanthroline(3v)





<sup>13</sup>C NMR Spectrum of 4-Phenyl-1,10-phenanthroline(3v)

## HRMS Spectrum of 4-Phenyl-1,10-phenanthroline(3v)

