

อนุพันธ์แนพทริดลพิรดีนและสมบัติภัยภาคเชิงแสง



นางสาวกอบกุล แซ่ปัง



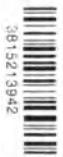
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ลิขสิทธิ์ของจุฬาลงกรณ์มหาวิทยาลัย



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NAPHTHYRIDYL PYRIDINE DERIVATIVES AND THEIR PHOTOPHYSICAL PROPERTIES

Miss Kobkun Sae Pang



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By	Miss Kobkun Sae Pang
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Thesis Advisor	Assistant Professor Rojrit Rojanathanes, Ph.D.
Thesis Co-Advisor	Assistant Professor Patchanita Thamyongkit, Dr.rer.nat.

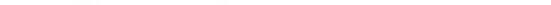
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 Dean of the Faculty of Science  
(Professor Supot Hannongbua, Dr.rer.nat.)

### THESIS COMMITTEE

Warinthon Chavasiri Chairman  
(Assistant Professor Warinthon Chavasiri, Ph.D.)

Rajit Bajaj Thesis Advisor

(Assistant Professor Rojrit Rojanathanes, Ph.D.)  


(Austin, LP, from P. L. Austin, The Great Lakes Region, 1974)

Nuanplum Chintarayini Examiner

(Associate Professor Nuanphun Chantarasiri, Ph.D.)

Vachiraporn Ativakom External Examiner

(Vachiraporn Ajavakom, Ph.D.)

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ได้สังเคราะห์ลิแกนด์ bis-naphthyridyl pyridines ชนิดใหม่ และสารเชิงซ้อนกับ Ru ขึ้นในงานวิจัยนี้โดยลิแกนด์มี 2 หน่วยฟังก์ชันหลักคือ หมู่คาร์บอชิลิก สำหรับยึดต่อ กับผิว ออกไซด์ในเซลล์สุริยะ และ ของที่มีในโตรเจนอยู่มากเพื่อสร้างสารเชิงซ้อนกับไอออนของโลหะ ทั้ง ลิแกนด์และสารเชิงซ้อนของ Ru(III) ได้รับการพิสูจน์เอกลักษณ์โดยสมบูรณ์ การทดสอบทาง กายภาพเชิงแสงชี้ให้เห็นว่า การดูดกลืนและการคายแสงสูงสุดของสารเชิงซ้อนเคลื่อนไปทางด้านสี แดง เนื่องจากความสามารถในการละลายในตัวทำละลายทั่วไปต่ำ จึงจำเป็นต้องปรับปรุง โครงสร้างของทั้งลิแกนด์และสารเชิงซ้อนนี้ต่อไป



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ปีการศึกษา 2556

ลายมือชื่อ อ.ที่ปรึกษาวิทยานิพนธ์หลัก Rit B.N.

ลายมือชื่อ อ.ที่ปรึกษาวิทยานิพนธ์ร่วม Patchanita T.

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KOBKUN SAE PANG: NAPHTHYRIDYL PYRIDINE DERIVATIVES AND THEIR PHOTOPHYSICAL PROPERTIES. ADVISOR: ASST. PROF. ROJIT ROJANATHANES, Ph.D., CO-ADVISOR: ASST. PROF. PATCHANITA THAMYONGKIT, Dr.rer.nat., 71 pp.

In this research a novel bis-naphthyridine ligand and its Ruthenium complex were synthesized. The ligand contains two major functionalities; carboxylic group for attaching with an oxide surface in solar cell and nitrogen – rich pocket for complexing with metal ion. Both ligand and its Ruthenium complex were fully characterized. The investigation of the photophysical properties indicated that their absorption and emission maxima of Ru(III) complex red shifted with its ligand. Due to their poor solubility in common solvents, further structural modification of both ligand and its complex is required.



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Student's Signature Kobkun Saengsang  
Advisor's Signature Rojit Rojanathanes  
Co-Advisor's Signature Patchanita T.

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## LIST OF ABBREVIATIONS

$^{\circ}\text{C}$	: degree Celcius
calcd	: calculated
$\text{CDCl}_3$	: deuterated chloroform
$^{13}\text{C}$ -NMR	: carbon-13 nuclear magnetic resonance spectroscopy
d	: day(s)
d	: doublet (NMR)
dd	: doublet of doublet (NMR)
DSSCs	: dye sensitized solar cells
$\text{DMSO}-d_6$	: hexadeuterated dimethylsulfoxide
equiv	: equivalent(s)
g	: grams(s)
h	: hours(s)
$^1\text{H}$ -NMR	: proton nuclear magnetic resonance spectroscopy
HOMO	: highest occupied molecule orbital
IR	: Infrared spectroscopy
$J$	: coupling constant
LMCT	: ligand to metal charge transfer
LUMO	: lowest unoccupied molecular orbital
m	: multiplet (NMR)
mmol	: millimole(s)
mg	: milligram(s)

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MHz	: megahertz (million Hertz)
ml	: millilitre(s)
MLCT	: metal to ligand charge transfer
m/z	: mass per charge ratio
obsd	: observed
rt	: room temperature
s	: singlet (NMR)
t	: triplet (NMR)
TCO	: transparent conductive oxide
$\delta$	: chemical shift
$\lambda$	: wavelength
$\lambda_{\text{ex}}$	: excitation wavelength
$\lambda_{\text{abs}}$	: absorption wavelength
$\lambda_{\text{em}}$	: emission wavelength
$\epsilon$	: molar absorptivity

