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กับไอกอนโลหะแทรนซิชันและโลหะหนัก

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**COMPLEXING PROPERTIES OF PHENOLIC DIAZACROWN
ETHERS WITH TRANSITION AND HEAVY METAL IONS**

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ศูนย์วิทยบริการ
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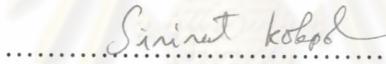
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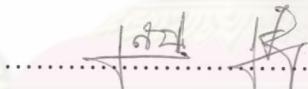
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ศึกษาการเกิดสารประกอบเชิงซ้อนระหว่างอนุพันธ์ของฟีโนลิกไดอะคราวน์อีเทอร์ (**L1-L6**) กับไอออนโลหะแทرنซิชันและโลหะหนักได้แก่ Co^{2+} Ni^{2+} Cu^{2+} Zn^{2+} Cd^{2+} Hg^{2+} และ Pb^{2+} ด้วยเทคนิคยูวีสเปกโตรโฟโตเมคริคไทเทรชัน พบสารประกอบเชิงซ้อนแบบ ML เป็นส่วนใหญ่ โดยสถิติรากของสารประกอบเชิงซ้อนของโลหะสูงขึ้น เมื่อหมู่แทนที่ (R) มีความยาวมากขึ้น ยกเว้นสารประกอบเชิงซ้อนของ Ni^{2+} และ Hg^{2+} นอกจากนั้นพบว่าลิเกนด์ทุกตัวเกิดสารประกอบเชิงซ้อนที่ถูกต้องได้ ($\log \beta_{\text{ML}} \geq 7$) ยกเว้น **L6** ($\log \beta_{\text{ML}} [\text{Pb}^{2+}\text{L6}] = 6.17$) ลำดับสถิติรากสำหรับสารประกอบเชิงซ้อนของโลหะแทرنซิชันแตกต่างกันคือ $\text{Zn}^{2+} < \text{Ni}^{2+} < \text{Co}^{2+}$ และสำหรับโลหะหมู่ IIB คือ $\text{Zn}^{2+} < \text{Hg}^{2+} < \text{Cd}^{2+}$ ยกเว้นระบบของ **L1** ซึ่งสารประกอบเชิงซ้อนของ Hg^{2+} จะมีสถิติรากสูงสุด เมื่อเปรียบเทียบลำดับของความเฉพาะเจาะจงต่อโลหะพบว่ามีแนวโน้มเดียวกับลำดับของความเสถียร

| | R | R_1 |
|-----------|---|----------------------------------|
| L1 | $-\text{CH}_3$ | - H |
| L2 | $-(\text{CH}_2)_7\text{CH}_3$ | - H |
| L3 | $-(\text{CH}_2)_{11}\text{CH}_3$ | - H |
| L4 | - H | $-(\text{CH}_2)_{14}\text{CH}_3$ |
| L5 | ^tBu | - H |
| L6 | $\text{C}(\text{CH}_3)_2\text{CH}_2^t\text{Bu}$ | - H |

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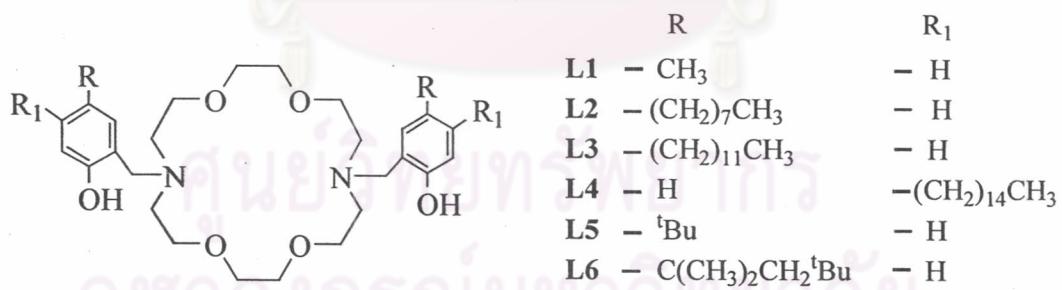
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Stability constants of complexes between phenolic diazacrown ether derivatives (**L1-L6**) and transition and heavy metal ions, i.e. Co^{2+} , Ni^{2+} , Cu^{2+} , Zn^{2+} , Cd^{2+} , Hg^{2+} and Pb^{2+} , were determined by UV spectrophotometric titration. Most of the systems studied exhibited ML complex formation, where a stability of complex increases, once substituents (R) extend, except for complexes of Ni^{2+} and Hg^{2+} . It has been additionally observed that Pb^{2+} forms very strong complexes with every ligand studied ($\log \beta_{\text{ML}} \geq 7$) and exact stability constants cannot be determined except that of **L6** ($\log \beta_{\text{ML}} = 6.17$). A stability order for first-row transition metal ions is: $\text{Zn}^{2+} < \text{Ni}^{2+} < \text{Co}^{2+}$ and that for IIB metal ions is: $\text{Zn}^{2+} < \text{Hg}^{2+} < \text{Cd}^{2+}$. The exception is found in the system of **L1** where Hg^{2+} forms the most stable complex. It has been compared and found that cation selectivity order shows the same trend as the stability orders.



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