

Development and Investigation of  $N\equiv W(OR)_3$ ,  $N\equiv M(OR)_3$ , and  $Mo_2(OR)_6$  Complexes  
for Triple-Bond Metathesis

by

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## List of Abbreviations

|                                 |  |
|---------------------------------|--|
| ACM                             | alkyne cross-metathesis                |
| AM                              | alkyne-metathesis                      |
| Anal                            | elemental analysis                     |
| Ar                              | aryl                                   |
| asym                            | asymmetrical                           |
| br                              | broad                                  |
| C <sub>6</sub> D <sub>6</sub>   | benzene- <i>d</i> <sub>6</sub>         |
| Calcd                           | calculated                             |
| CD <sub>2</sub> Cl <sub>2</sub> | dichloromethane- <i>d</i> <sub>2</sub> |
| CDCl <sub>3</sub>               | chloroform- <i>d</i>                   |
| cm <sup>-1</sup>                | wavenumber                             |
| d                               | day(s), doublet                        |
| DCE                             | 1,2-dichloroethane                     |
| DFT                             | density functional theory              |
| DME                             | 1,2-dimethoxyethane                    |
| EI MS                           | electron impact mass spectrometry      |
| Et                              | ethyl                                  |
| EtCN                            | propionitrile                          |

|                       |                                      |
|-----------------------|--------------------------------------|
| $\text{Et}_2\text{O}$ | diethyl ether                        |
| equiv                 | equivalent(s)                        |
| g                     | grams                                |
| GC/MS                 | gas chromatography mass spectrometry |
| GOF                   | goodness of fit                      |
| $^1\text{H}$          | proton                               |
| h                     | hour(s)                              |
| Hz                    | hertz                                |
| HOMO                  | highest occupied molecular orbital   |
| HOTF                  | trifluoromethanesulfonic acid        |
| kcal                  | kilocalorie(s)                       |
| $K_{\text{eq}}$       | equilibrium constant                 |
| LA                    | Lewis acid                           |
| LAs                   | Lewis acids                          |
| LUMO                  | lowest unoccupied molecular orbital  |
| m                     | multiplet                            |
| M                     | molar                                |
| Me                    | methyl                               |
| MeCN                  | acetonitrile                         |
| mg                    | milligrams                           |
| min                   | minute(s)                            |
| mL                    | milliliters                          |
| mol                   | moles                                |

|                       |   |
|-----------------------|---|
| <b>mM</b>             | millimolar                                |
| <b>mmol</b>           | millimoles                                |
| <b>NACM</b>           | nitrile-alkyne cross metathesis           |
| <b>NAX</b>            | nitrogen atom exchange                    |
| <b>NCAr</b>           | aryl nitrile                              |
| <b>NM</b>             | nitrile metathesis                        |
| <b>NMR</b>            | nuclear magnetic resonance                |
| <b>ORTEP</b>          | oak ridge thermal ellipsoid plot          |
| <b>q</b>              | quartet                                   |
| <b>Q</b>              | reaction endpoint                         |
| <b>Ph</b>             | phenyl                                    |
| <b>POSS</b>           | polyhedral oligomeric silsesquioxane      |
| <b>ppm</b>            | parts per million                         |
| <b>s</b>              | singlet                                   |
| <b>sym</b>            | symmetrical                               |
| <b>t</b>              | triplet                                   |
| <b><sup>t</sup>Bu</b> | CMe <sub>3</sub>                          |
| <b>THF</b>            | tetrahydrofuran                           |
| <b>TS</b>             | transition state                          |
| <b>XRD</b>            | x-ray diffraction                         |
| <b>°C</b>             | degrees Celsius                           |
| <b>Å</b>              | angstroms                                 |
| <b>δ</b>              | chemical shift in ppm downfield from zero |

$\mu\text{L}$  microliter(s)

$\equiv$  triple bond