

## Appendix I

# Supplemental Data for Table 13-3. 12-Vertex Transition Element Metallacarboranes

Compound <sup>a</sup>	Information <sup>b</sup>	References
Synthesis and Characterization		
<b>Scandium</b>		
<i>Closo-ScC<sub>2</sub>B<sub>9</sub> clusters</i>		
[3,1,2-Cp*(H)Sc(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ] <sup>2-</sup>	S, X, H, C	[541]
3,1,2-Cp*[(Me <sub>3</sub> Si) <sub>2</sub> CH]Sc(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> Li <sup>+</sup>	S, H, B, C	[541]
	X	[542]
{3,1,2-Cp*[(Me <sub>3</sub> Si) <sub>2</sub> CH]Sc(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> Li·Li(THF) <sub>3</sub> }	X	[541]
3,1,2-[H <sub>2</sub> C <sub>2</sub> (DIPP) <sub>2</sub> N <sub>2</sub> C=N](THF) <sub>2</sub> Sc(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) imidazolin-2- iminato complex	S, X, H, B, C	[1423]
3,1,2-[H <sub>2</sub> C <sub>2</sub> (DIPP) <sub>2</sub> N <sub>2</sub> C=N](THF)Sc[(NH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Sc-N imidazolin-2-iminato complex	S, X, H, B, C	[1423]
3,1,2-( <i>cyclo</i> -C <sub>3</sub> N <sub>2</sub> Ar <sub>2</sub> )=N-Sc(THF)[(H <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Sc-N Ar = 2,6-C <sub>6</sub> H <sub>3</sub> (CHMe <sub>2</sub> ) <sub>2</sub>	S, X, H, B, C	[1430]
<b>Yttrium</b>		
<i>Closo-YC<sub>2</sub>B<sub>9</sub> clusters</i>		
3,1,2-(THF) <sub>2</sub> Na(μ-Cl) <sub>2</sub> (THF) <sub>2</sub> Y(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR	[543]
3,1,2-(THF) <sub>2</sub> K(μ-Cl) <sub>2</sub> Y[(Me <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> )(MeOCH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] Y-N, Y-O	S, X, H, B, C, IR	[544]
3,1,2-[H <sub>2</sub> C <sub>2</sub> (DIPP) <sub>2</sub> N <sub>2</sub> C=N](THF) <sub>2</sub> Y(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) imidazolin-2- iminato complex	S, X, H, B, C	[1423]
3,1,2-( <i>cyclo</i> -C <sub>3</sub> N <sub>2</sub> Ar <sub>2</sub> )=N-Y(THF)[(H <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Y-N Ar = 2,6-C <sub>6</sub> H <sub>3</sub> (CHMe <sub>2</sub> ) <sub>2</sub>	S, X, H, B, C	[1430]
3,1,2-( <i>cyclo</i> -C <sub>3</sub> N <sub>2</sub> Ar <sub>2</sub> )=N-Y[(MeOCH <sub>2</sub> CH <sub>2</sub> )(H <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> )- C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] Y-O, Y-N Ar = 2,6-C <sub>6</sub> H <sub>3</sub> (CHMe <sub>2</sub> ) <sub>2</sub>	S, X, H, B, C	[1430]
3,1,2-( <i>cyclo</i> -C <sub>3</sub> N <sub>2</sub> Ar <sub>2</sub> )=N-Y[(H <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] 2 Y-N Ar = 2,6-C <sub>6</sub> H <sub>3</sub> (CHMe <sub>2</sub> ) <sub>2</sub>	S, X, H, B, C	[1430]
3,1,2-(thf) <sub>2</sub> ( <i>cyclo</i> -CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> NMe <sub>2</sub> )Y[( <i>cyclo</i> -CH <sub>2</sub> OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, C, IR	[1571]

Compound	Information	References
3,1,2-(thf) <sub>2</sub> [OCPh(C <sub>5</sub> H <sub>4</sub> N)(CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> NMe <sub>2</sub> )( <i>cyclo</i> -CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> NMe <sub>2</sub> )Y [( <i>cyclo</i> -CH <sub>2</sub> OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] N—Y	S, X, H, B, C, IR	[1571]
3,1,2-[(Me <sub>3</sub> C) <sub>3</sub> C <sub>6</sub> H <sub>2</sub> NC][CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> NMe <sub>2</sub> )Y( <i>cyclo</i> -CH <sub>2</sub> OCH <sub>2</sub> - C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) Y—N	S, H, B, C, IR	[1571]
3,1,2-(MeOCH <sub>2</sub> CH <sub>2</sub> OMe) <sub>2</sub> Y[(C <sub>9</sub> H <sub>6</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Y—C <sub>9</sub> H <sub>6</sub> constrained-geometry	S, X, H, B, C, IR, MS	[1588]
3,1,2-(thf) <sub>2</sub> Y{[Ph(NC <sub>5</sub> H <sub>4</sub> )C(O)C <sub>9</sub> H <sub>6</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Y—O, Y—N constrained-geometry	S, X, H, B, C, IR, MS	[1588]
3,1,2-(thf) <sub>3</sub> Y{[Ph <sub>2</sub> C(O)C <sub>9</sub> H <sub>6</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Y—O constrained-geometry	S, X, H, B, C, IR, MS	[1588]
3,1,2-(MeOCH <sub>2</sub> CH <sub>2</sub> OMe)Y{[Ph <sub>2</sub> C=C(O)C <sub>9</sub> H <sub>6</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Y—O constrained-geometry	S, H, B, C, IR, MS	[1588]
3,1,2-(thf) <sub>3</sub> Y{[(NC <sub>5</sub> H <sub>4</sub> ) <sub>2</sub> C(O)C <sub>9</sub> H <sub>6</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Y—O, Y—N constrained-geometry	S, H, B, C, IR, MS	[1588]
<b>Lanthanum</b>		
<i>Closo</i> -LaC <sub>2</sub> B <sub>9</sub> clusters		
3,1,2-(THF) <sub>2</sub> La(MeOCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sup>−</sup>	S, H, B, IR	[545]
<b>Lanthanide Elements</b>		
<i>LnC<sub>2</sub>B<sub>9</sub></i> clusters		
3,1,2-(THF) <sub>4</sub> Ln(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) Ln = Sm, Yb	S, X(Sm), H, B, IR, MAG	[547,548]
3,1,2-(HCONMe <sub>2</sub> ) <sub>4</sub> Yb(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, IR	[547]
	X	[547,548]
3,1,2-(THF) <sub>2</sub> (Me <sub>2</sub> HSi) <sub>2</sub> Ln[(Me <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Ln-N Ln = Sm, Er	S, X, H, B, C, IR	[544]
3,1,2-[H <sub>2</sub> C <sub>2</sub> (DIPP) <sub>2</sub> N <sub>2</sub> C=N] (THF) <sub>2</sub> Lu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) imidazolin-2- iminato complex	S, X, H, B, C	[1423]
3,1,2-[H <sub>2</sub> C <sub>2</sub> (DIPP) <sub>2</sub> N <sub>2</sub> C=N](THF)Lu[(NH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] M—N imidazolin-2-iminato complex	S, X, H, B, C	[1423]
3,1,2-( <i>cyclo</i> -C <sub>3</sub> N <sub>2</sub> Ar <sub>2</sub> )=N—Lu(THF) <sub>2</sub> (C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) Ar = 2,6- C <sub>6</sub> H <sub>3</sub> (CHMe <sub>2</sub> ) <sub>2</sub>	S, X, H, B, C	[1430]
3,1,2-( <i>cyclo</i> -C <sub>3</sub> N <sub>2</sub> Ar <sub>2</sub> )=N—Lu(THF)[(H <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] Lu—N Ar = 2,6-C <sub>6</sub> H <sub>3</sub> (CHMe <sub>2</sub> ) <sub>2</sub>	S, X, H, B, C	[1430]
{3,1,2-(thf)Cl <sub>2</sub> Gd[( <i>cyclo</i> -CH <sub>2</sub> OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]} <sub>2</sub> Na <sub>2</sub> (thf) <sub>4</sub>	S, X, H, B, C, IR	[1571]
<b>Uranium</b>		
<i>Closo</i> -UC <sub>2</sub> B <sub>9</sub> clusters		
3-U[1,2-(MeOCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub>	S, X, H, B	[545]
3,1,2-(THF) <sub>2</sub> I <sub>2</sub> U(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>−</sup>	S, X, H	[552]
3-Br <sub>2</sub> U(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub>	S, E	[553]
<b>Titanium</b>		
<i>Closo</i> -TiC <sub>2</sub> B <sub>9</sub> clusters		
3,1,2-Cp*MeTi(PMe <sub>3</sub> )(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, P	[554]
3,1,2-Cp*(N=CMe <sub>2</sub> )Ti(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, IR	[554]
3,1,2-Cp*RTi(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R = Et, CMe=CMe <sub>2</sub>	S, H, B, C	[554]
3,1,2-Cp*Ti(MeCN)(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-CHMeO-)	S, X, H(2d), B, C	[555]
2,1,7-(η <sup>8</sup> -C <sub>8</sub> H <sub>8</sub> )Ti(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>−</sup>	S, B, IR, E	[557]

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Compound	Information	References
3,1,2-(Me <sub>4</sub> C <sub>5</sub> CH <sub>2</sub> —)Ti(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR	[554]
	X	[556]
3,1,2-[ <i>cyclo</i> -(C <sub>13</sub> H <sub>9</sub> )P[N(CHMe <sub>2</sub> ) <sub>2</sub> —O—](NMe <sub>2</sub> ) <sub>2</sub> Ti(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) C <sub>13</sub> H <sub>9</sub> =fluorenyl Ti—O (FF)	S, X, H, B, C, P, IR	[558]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> (HNMe <sub>2</sub> )Ti(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, IR	[559]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> (HNMe <sub>2</sub> )Ti[(C <sub>9</sub> H <sub>7</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, H, B, C	[560]
3,1,2-Ln(Me <sub>2</sub> N)Ti[(C <sub>9</sub> H <sub>6</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] L=DME, THF, C <sub>5</sub> H <sub>5</sub> N	S, X(DME), H, B, C	[560]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> Ti[(Me <sub>2</sub> NCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, C	[561]
[(MeN)Ti[(Me <sub>2</sub> NCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> Ti(NMe <sub>2</sub> ) <sub>2</sub> (μ-O)	S, X, H, B	[561]
3,1,2-ClTi[(H <sub>2</sub> NCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, H, B, C, IR	[562]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> Ti[(R <sub>2</sub> C <sub>6</sub> H <sub>3</sub> -N=CH)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] R=CHMe <sub>2</sub> , Me Ti—N imido catalyst for polymerization of C <sub>2</sub> H <sub>4</sub> with MAO (R=CHMe <sub>2</sub> )	S, X, H, B, C, IR, MS	[563]
3,1,2-ClTi[Me(C <sub>6</sub> H <sub>10</sub> -O)C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] Ti—O C <sub>6</sub> H <sub>10</sub> =cyclohexyl	S, H, B, C, IR	[564]
3,1,2-Cl <sub>2</sub> Ti{[(PhCH <sub>2</sub> ) <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> ]RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> } Ti—N R=H, Me	S, H, B, C, IR, MS	[565]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> Ti{[(PhCH <sub>2</sub> ) <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> ]RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> } Ti—N R=H, Me	S, H, B, C, IR, MS	[565]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> (HMe <sub>2</sub> N)Ti{[(PhCH <sub>2</sub> ) <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> ]RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> } R=H, Me	S, X(H), H, B, C, IR, MS	[565]
3,1,2-L <sub>2</sub> Ti{[(NC <sub>5</sub> H <sub>4</sub> CH <sub>2</sub> )RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] Ti—N L=NMe <sub>2</sub> , Cl; R=H, Me picolyl	S, X(NMe <sub>2</sub> ), H, B, C, IR	[567]
3,1,2-(OCHMe <sub>2</sub> ) <sub>2</sub> Ti{[(NC <sub>5</sub> H <sub>4</sub> CH <sub>2</sub> )RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] Ti—N R=H, Me picolyl	S, X(Me), H, B, C, IR	[567]
3,1,2-(OPh) <sub>2</sub> Ti{[(NC <sub>5</sub> H <sub>4</sub> CH <sub>2</sub> )RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] (Ti—N) picolyl	S, X, H, B, C, IR	[567]
3-Ti[1,2-(Me <sub>2</sub> NCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> Ti—N	S, H	[566]
3,1,2-(NHR) <sub>2</sub> (NR <sub>2</sub> ) <sub>2</sub> Ti(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R=Me, Et	S, H, B, C	[568]
3,1,2-[ <i>cyclo</i> -CyNC(NMe <sub>2</sub> )NCy]Ti[(CH <sub>2</sub> O)(Me <sub>2</sub> NCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] Cy= <i>p</i> -MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> Ti—N, Ti—O reaction with Me <sub>2</sub> NH→ (Me <sub>2</sub> N)Ti[(Me <sub>2</sub> NCH <sub>2</sub> )CH <sub>2</sub> O]C <sub>2</sub> B <sub>9</sub> H <sub>9</sub>	S, X, H, B, C	[570]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> (Me <sub>2</sub> NH)Ti[(CH <sub>2</sub> OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B, C, IR	[571]
3,1,2-(R <sub>2</sub> N)Ti[(Me <sub>2</sub> NCH <sub>2</sub> )(OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] R=Me, Et	S, X, H, B, C, IR	[571]
3,1,2-(MeOC <sub>6</sub> H <sub>4</sub> NH)Ti[(Me <sub>2</sub> NCH <sub>2</sub> )(OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B, C, IR	[571]
3,1,2-(NC <sub>5</sub> H <sub>3</sub> Me—NH)(NC <sub>5</sub> H <sub>3</sub> Me—NH <sub>2</sub> )Ti[(Me <sub>2</sub> NCH <sub>2</sub> )(OCH <sub>2</sub> )- C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, C, IR	[571]
[3,1,2-(MeO)Ti[Et <sub>2</sub> NCH <sub>2</sub> )(OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> 2 Ti—O—Ti	S, X, H, B, C, IR	[571]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> Ti[(Me <sub>2</sub> NCH <sub>2</sub> )RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] R=H, Me	S, X, H, B, C, IR	[572]
3-Ti[1,2-(Me <sub>2</sub> NCH <sub>2</sub> )RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> R=H, Me	S, X(H), H, B, C, IR	[572]
3,1,2-(RO) <sub>2</sub> Ti[(Me <sub>2</sub> NCH <sub>2</sub> )RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] R=Ph, MeC <sub>6</sub> H <sub>4</sub> , OCHMe <sub>2</sub>	S, X, H, B, C, IR, C <sub>2</sub> H <sub>4</sub> polymerization	[572]
3,1,2-(Me <sub>2</sub> N)Ti [(Me <sub>2</sub> NCH <sub>2</sub> )(OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] N→Ti, O→Ti catalyst for T-amide catalyzed synthesis of cyclic guanidines from di- and triamines and carbodiimides		[1454]
<b>Zirconium</b>		
<i>Closo</i> -ZrC <sub>2</sub> B <sub>9</sub> clusters		
3,1,2-(Me <sub>2</sub> N)(HMe <sub>2</sub> N)Zr[(HMe <sub>4</sub> C <sub>5</sub> —CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, C	[1434]
3,1,2-(Me <sub>3</sub> SiCH <sub>2</sub> ) <sub>2</sub> Zr[(HMe <sub>4</sub> C <sub>5</sub> —CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>−</sup>	S, X, H, B, C	[1434]

Compound	Information	References
3,1,2-[( $\sigma, \sigma$ -CH <sub>2</sub> (NMe <sub>2</sub> )- <i>o</i> -C <sub>6</sub> H <sub>4</sub> ) Zr[(HMe <sub>4</sub> C <sub>5</sub> -CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]] <sup>-</sup>	S, X, H, B, C	[1434]
3,1,2-[MeN(CH <sub>2</sub> ) <sub>n</sub> NMeH]Zr[(CpCMe <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <i>n</i> =2,3 Zr—Cp	S, X, H, B	[1445]
{3,1,2-[MeN(CH <sub>2</sub> ) <sub>2</sub> NMe]Zr[(CpCMe <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]} <sub>2</sub> Li <sub>2</sub> Zr—Cp	S, X, H, B	[1445]
3,1,2-(THF)ClZr[(H <sub>2</sub> NCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, H, B, C, IR	[562]
3-(THF) <sub>2</sub> Zr[(1,2- <i>o</i> -C <sub>6</sub> H <sub>4</sub> CH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub>	S, H, B, C, IR	[574]
3,1,2-Cp*(MeC=CMe <sub>2</sub> )Zr(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X	[575]
3,1,2-Cl <sub>2</sub> (NEt <sub>2</sub> ) <sub>2</sub> Zr(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C	[568]
(THF) <sub>n</sub> Na <sub>3</sub> <sup>3+</sup> 3,1,2-Cl <sub>2</sub> Zr[(C <sub>5</sub> H <sub>4</sub> -CMe <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>3-</sup>	S, X, H, B, C, IR	[577]
3,1,2-CpClZr[(C <sub>5</sub> H <sub>4</sub> -CMe <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup> Na(DME) <sub>2</sub> <sup>+</sup>	S, X, H, B, C, IR	[577]
3,1,2-(CH <sub>2</sub> Ph)Zr[(C <sub>5</sub> H <sub>4</sub> -CMe <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup> Na(DME) <sub>3</sub> <sup>+</sup> ethylene polymerization catalyst	S, X, H, B, C, IR	[577]
3,1,2-[ <i>cyclo</i> -(C <sub>13</sub> H <sub>9</sub> )P{N(CHMe <sub>2</sub> ) <sub>2</sub> O-}](NMe <sub>2</sub> ) <sub>2</sub> Zr(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) C <sub>13</sub> H <sub>9</sub> =fluorenyl Zr—O (FF)	S, X, H, B, C, P, IR	[558]
3,1,2-[C <sub>6</sub> H <sub>3</sub> (CHMe <sub>2</sub> ) <sub>2</sub> -NH](THF)Zr[(C <sub>5</sub> H <sub>4</sub> -CMe <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, C, IR	[577]
3,1,2-(THF)(NEt <sub>2</sub> ) <sub>2</sub> Zr(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C	[568]
3,1,2-L <sub>n</sub> (Me <sub>2</sub> N)Zr[(C <sub>9</sub> H <sub>6</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] L = DME, THF, C <sub>5</sub> H <sub>5</sub> N	S, X(THF), H, B, C	[560]
3,1,2-Cl <sub>2</sub> Zr[(C <sub>9</sub> H <sub>6</sub> -Me <sub>2</sub> C)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	S, H, B, C, IR	[578]
3,1,2-LClZr[(C <sub>9</sub> H <sub>6</sub> -Me <sub>2</sub> C)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup> L = PhCH <sub>2</sub> , Cp	S, X, H, B, C, IR	[578]
3,1,2-(THF)LZr[(C <sub>9</sub> H <sub>6</sub> -Me <sub>2</sub> C)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] L = Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub> NH, <i>n</i> -C <sub>4</sub> H <sub>9</sub> O	S, X, H, B, C, IR	[578]
3,1,2-(THF)(Me <sub>3</sub> SiCH <sub>2</sub> )Zr[(CH <sub>2</sub> ) <sub>2</sub> NMe( $\mu$ -CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, C	[579]
{3,1,2-[MeO(CH <sub>2</sub> ) <sub>2</sub> O]Zr[(CH <sub>2</sub> ) <sub>2</sub> NMe( $\mu$ -CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]} <sub>2</sub>	S, X, H, B, C	[579]
3,1,2-(Me <sub>2</sub> NH)(Me <sub>2</sub> N) <sub>2</sub> Zr{[(CHMe <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>3</sub> -N=CH]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> } R = CHMe <sub>2</sub> , Me imido catalyst for polymerization of C <sub>2</sub> H <sub>4</sub> with MAO R = CHMe <sub>2</sub>	S, X, H, B, C, IR, MS	[563]
3,1,2-ClZr[Me(C <sub>6</sub> H <sub>10</sub> -O-)C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] Zr—O C <sub>6</sub> H <sub>10</sub> =cyclohexyl catalytic polymerization of C <sub>2</sub> H <sub>4</sub> with MAO	S, H, B, C, IR	[564]
(polystyryl){3,1,2-ClZr[Me(C <sub>6</sub> H <sub>10</sub> -O-)C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]} <sub>n</sub> M—O (C <sub>6</sub> H <sub>10</sub> =cyclohexyl) catalytic polymerization of C <sub>2</sub> H <sub>4</sub> with MAO	S, IR	[564]
3,1,2-Cl <sub>2</sub> Zr{[(PhCH <sub>2</sub> ) <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> ]RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> } Zr—N R = H, Me	S, H, B, C, IR, MS	[565]
3-Zr[1,2-(Me <sub>2</sub> NCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> Zr—N	S, X, H	[566]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> (Me <sub>2</sub> NH)Zr[(CH <sub>2</sub> OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B, C, IR	[571]
3,1,2-(Cl)Zr[(polystyryl)(C <sub>6</sub> H <sub>10</sub> -2'-NH)C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] Zr—N catalyst for polymerization. of C <sub>2</sub> H <sub>4</sub> and vinyl chloride with MAO	S, H, B, C, IR, TGA, far IR	[569]
3,1,2-[(Me <sub>3</sub> Si) <sub>2</sub> C <sub>5</sub> H <sub>3</sub> ]Zr{[MeN(CH <sub>2</sub> )CH <sub>2</sub> CH <sub>2</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> } alkyne insertion	S, X, H, B, C	[1595]
<b>Hafnium</b>		
<i>Closo-HfC<sub>2</sub>B<sub>9</sub> clusters</i>		
3-Cl(THF)Hf[1,2-(PhCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> ·Na(THF) <sub>3</sub>	S, H, B, C, IR	[574]
3,1,2-Cp*(MeC=CMe <sub>2</sub> )Hf(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S	[575]
[3-Cp*Hf(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> ( $\mu$ -CH <sub>2</sub> )	S	[575]
3-Hf[1,2-(Me <sub>2</sub> NCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> Ti—N	S	[566]
3,1,2-(Me <sub>2</sub> NH)(Me <sub>2</sub> N) <sub>2</sub> Hf{[(CHMe <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>3</sub> -N=CH]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> } R = CHMe <sub>2</sub> , Me imido catalyst for polymerization of C <sub>2</sub> H <sub>4</sub> with MAO R = CHMe <sub>2</sub>	S, X, H, B, C, IR, MS	[563]

Continued

Compound	Information	References
3,1,2-(THF)(Me <sub>3</sub> SiCH <sub>2</sub> )Hf[(CH <sub>2</sub> ) <sub>2</sub> NMe(μ-CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, H, B, C	[579]
3,1,2-Cl <sub>2</sub> Hf[(C <sub>9</sub> H <sub>6</sub> -Me <sub>2</sub> C)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	S, H, B, C, IR	[578]
3,1,2-Cp*Hf(7-CHMe)(3,5-Me <sub>2</sub> C <sub>5</sub> H <sub>3</sub> )(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -7-)	S, H, C	[582]
Li(THF) <sub>2</sub> <sup>+</sup> (μ-Cl) <sub>2</sub> Hf[(Me <sub>4</sub> C <sub>5</sub> -CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, C	[1434]
<b>Niobium</b>		
<i>Closo-NbC<sub>2</sub>B<sub>9</sub> clusters</i>		
2,1,7-(Me <sub>2</sub> N) <sub>3</sub> Nb(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C	[584]
3,1,2-(Me <sub>2</sub> N) <sub>3</sub> Nb(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C	[584]
<b>Tantalum</b>		
<i>Closo-TaC<sub>2</sub>B<sub>9</sub> clusters</i>		
3,1,2-(MeC <sub>5</sub> H <sub>4</sub> )Me <sub>2</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C	[585]
3,1,2-(MeC <sub>5</sub> H <sub>4</sub> )Cl <sub>2</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C	[585]
3-ClMeTa(μ-H)(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub>	S, X, H, B(2d), C, MS	[587]
3,1,2-(Me <sub>2</sub> N) <sub>3</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, IR, MS	[583]
2,1,7-(Me <sub>2</sub> N) <sub>3</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -3-Me)	S, H, B, C	[588]
3,1,2-(Me <sub>2</sub> N) <sub>3</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-Me)	S, H, B, C	[588]
[3,1,2-(Me <sub>2</sub> N) <sub>2</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> (μ-O)]	X	[589]
3,1,2-(Me <sub>2</sub> N) <sub>3</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C	[584]
2,1,12-Br(Me <sub>2</sub> N) <sub>2</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR	[589]
3,1,2-[p-FC <sub>6</sub> H <sub>4</sub> C(NMe <sub>2</sub> )=N] <sub>2</sub> ClTa(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, MS	[590]
3,1,2-(Me <sub>2</sub> N) <sub>2</sub> [(C <sub>6</sub> H <sub>13</sub> )N=C-NMe <sub>2</sub> ]Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, MS	[590]
3,1,2-[2,4-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub> -O] <sub>3</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, MS	[590]
3,1,2-(PhS) <sub>4</sub> Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, MS	[590]
3,1,2-[(C <sub>6</sub> H <sub>11</sub> ) <sub>7</sub> Si <sub>7</sub> O <sub>12</sub> ]Ta(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) sesquioxane	S, H, C, IR, MS	[591]
3,1,2-(C <sub>6</sub> H <sub>3</sub> Me <sub>2</sub> )-N=Ta[(CH <sub>2</sub> NHCH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -n-CHMe <sub>2</sub> ] N → Ta n = 3, 6, 8	S, X(n=8), H, C, B	[1499]
3,1,2-(C <sub>6</sub> H <sub>3</sub> Me <sub>2</sub> )-N=Ta-[N(=CH <sub>2</sub> )C <sub>6</sub> H <sub>3</sub> Me <sub>2</sub> ][Me <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> -C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -n-CHMe <sub>2</sub> ] N → Ta n = 3, 8	S, X(n=8), H, C, B	[1499]
3,1,2-(C <sub>5</sub> H <sub>5</sub> N)(Me <sub>2</sub> N)Ta[(Me <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ] N → Ta	S, H, B, C	[1540]
<b>Chromium</b>		
<i>Closo-CrC<sub>2</sub>B<sub>9</sub> clusters</i>		
3,1,2-CpCr(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R = H, Me, Ph	S, IR, UV	[594]
3,1,2-CpCr(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R = H, Me	S, IR, UV	[594]
3,1,2-Cr(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(1,1')-(CH <sub>2</sub> ) <sub>4</sub> <sup>-</sup>	S, H, B, C, IR	[598]
3,1,2-Cl(NC <sub>5</sub> H <sub>4</sub> -CH <sub>2</sub> -)Cr(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, IR, MS	[604]
[3,1,2-(μ-Cl)ClCr[R(Me <sub>2</sub> NCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> R = H, Me	S, X, IR, MS	[1590]
<b>Molybdenum</b>		
<i>Closo-MoCB<sub>10</sub> clusters</i>		
2,1-(CO) <sub>4</sub> Mo(OH)CB <sub>10</sub> H <sub>10</sub> <sup>-</sup>	S, B, IR	[606]
2,1-(CO) <sub>3</sub> (Ph <sub>3</sub> P)Mo(CB <sub>10</sub> H <sub>10</sub> -7-NMe=CHMe) ( <i>cis, trans</i> )	S, H, B, C, P, IR	[608]

Compound	Information	References
2,1-(CO) <sub>3</sub> LMo(CB <sub>10</sub> H <sub>10</sub> -7-L) L = [OCH <sub>2</sub> ] <sub>4</sub> , OEt <sub>2</sub> ; L' = PPh <sub>3</sub> , CNCMe <sub>3</sub>	S, X([OCH <sub>2</sub> ] <sub>4</sub> ), H, B, C	[609]
2,1-(RC≡CR')(CO) <sub>2</sub> Mo(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup> R, R' = H, Ph, CMe <sub>3</sub>	S, H, B, C	[609]
2,1-(Ph <sub>3</sub> P) <sub>2</sub> (CO) <sub>2</sub> Mo(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, C	[609]
2,1-(CO) <sub>3</sub> (Ph <sub>3</sub> P)Mo(CB <sub>10</sub> H <sub>8</sub> )-(μ-H) <sub>2</sub> -Ag(PPh <sub>3</sub> )	S, X, H, B, C, P	[610]
2,1-(CO) <sub>3</sub> (Ph <sub>3</sub> P)Mo(CB <sub>10</sub> H <sub>8</sub> )-(μ-H) <sub>3</sub> -Ru LL'L'' L, L', L'' = Cl, PPh <sub>3</sub>	S, X(Cl, 2PPh <sub>3</sub> ), H, B, C, P	[610]
2,1-(Me <sub>3</sub> CNC) <sub>4</sub> XMo(CB <sub>10</sub> H <sub>11</sub> ) X = I, Br	S, X(I), H, B, C	[611]
Related derivatives		[611]
2,1-(CO) <sub>3</sub> Mo[(Me <sub>3</sub> CHN)CB <sub>10</sub> H <sub>10</sub> ] <sup>-</sup> Cu(PPh <sub>3</sub> ) <sup>+</sup>	S, H, B, C, P	[592]
2,1-(CO) <sub>3</sub> (Ph <sub>3</sub> PAu)Mo[(Me <sub>3</sub> CHN)CB <sub>10</sub> H <sub>10</sub> ]	S, H, B, C, P	[592]
2,1-(CO) <sub>3</sub> Mo[(μ-NHCR=NH)CB <sub>10</sub> H <sub>10</sub> ] <sup>-</sup> R = Me, CMe=CH <sub>2</sub> , CH <sub>2</sub> Ph, Et	S, X(Me), H, B, C, IR	[612]
Related derivatives		[612]
2,1-(CO) <sub>2</sub> (HC≡CCMe <sub>3</sub> )Mo(CB <sub>10</sub> H <sub>10</sub> -3-NMe <sub>3</sub> )	S, X, H, B, C	[613]
2,1-(Me <sub>3</sub> CNC) <sub>4</sub> BrMo(CB <sub>10</sub> H <sub>10</sub> -3-X) X = Br, I	S, X(Br), H, B, C	[613]
Exo-[Cp* <sub>2</sub> Rh <sub>2</sub> (μ-CO)]-2,1-(CO) <sub>3</sub> Mo(CB <sub>10</sub> H <sub>10</sub> )	S, X, H, B, C	[613]
Related derivatives		[613]
2,1-(CO) <sub>3</sub> Mo[(μ-NHCMe=CO → Mo)CB <sub>10</sub> H <sub>10</sub> ] <sup>-</sup>	S, B, C, IR	[614]
Related derivatives		[614]
2,1-(CO) <sub>3</sub> Mo[R(CB <sub>10</sub> H <sub>9</sub> )]-μ(2,3)-N(CH <sub>2</sub> ) <sub>5</sub> <sup>-</sup> R = H, NH <sub>2</sub>	S, X(H), H, B, C, IR	[615]
Related derivatives		[615]
<b>Closo-MoC<sub>2</sub>B<sub>9</sub> clusters (no exo-polyhedral metals)</b>		
1,2,4-(η <sup>7</sup> -C <sub>7</sub> H <sub>7</sub> )Mo(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, IR	[616]
2,1,8-(η <sup>7</sup> -C <sub>7</sub> H <sub>7</sub> )Mo(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, IR	[616]
2,1,8-(CO) <sub>3</sub> IMo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -11-CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> Me) <sup>-</sup>	S, H, B, C, IR	[617]
2,1,8-(CO) <sub>3</sub> BrMo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, B, C, IR	[1320]
2,1,8-(CO) <sub>3</sub> IMo[Ph(OH)C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sup>-</sup>	S, X, H, B, C	[353]
2,1,8-(Me <sub>3</sub> CNC) <sub>2</sub> (O)Mo[Ph(NHCMe <sub>3</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, C	[353]
3,1,2-(CO) <sub>3</sub> Mo(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>2-</sup> R = H, Me	S, IR	[618]
3,1,2-(CO) <sub>3</sub> RMo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> R = H, Me	S, H, IR	[618]
3,1,2-(CO) <sub>3</sub> Mo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-1-CH <sub>2</sub> -[cyclo-N <sub>3</sub> P <sub>3</sub> ](C <sub>5</sub> H <sub>10</sub> N) <sub>4</sub> Me	S, P, IR	[619]
3,1,2-[(MeO) <sub>3</sub> P](CO)(MeC <sub>6</sub> H <sub>4</sub> C)Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, B, C, P	[620]
3,1,2-(CO) <sub>2</sub> (μ-MeC <sub>6</sub> H <sub>4</sub> C)Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, C, IR	[621]
3,1,2-μ-MeC <sub>6</sub> H <sub>4</sub> C=Mo(CO) <sub>2</sub> (PMe <sub>3</sub> )(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, C	[623]
3,1,2-(CO) <sub>2</sub> (C <sub>4</sub> H <sub>6</sub> )Mo(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, P, IR	[624]
Related derivatives		[624]
3,1,2-(R <sub>2</sub> C <sub>2</sub> )(Ph <sub>3</sub> P)(CO)Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[625]
3,1,2-(Ph <sub>3</sub> P)(CO) <sub>2</sub> Mo[(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -8-PhC <sub>2</sub> PhC(→ Mo))]	S, H, B, C, IR	[625]
Related derivatives		[625]
3,1,2-(C <sub>3</sub> H <sub>5</sub> )(CO) <sub>2</sub> Mo(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -10-L) <sup>-</sup> L = OEt <sub>2</sub> , O(CH <sub>2</sub> ) <sub>4</sub> F	S, H, B, C, IR	[626]
3,1,2-(CO) <sub>4</sub> Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -10-L) L = OEt <sub>2</sub> , O(CH <sub>2</sub> ) <sub>4</sub> F	S, H, B, C, IR	[626]

Continued

Compound	Information	References
3,1,2-(Me <sub>3</sub> CNC) <sub>3</sub> (CO)Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[627]
3,1,2-(Me <sub>3</sub> CNC) <sub>4</sub> Mo(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=Me, H	S, H, B, C, IR	[627]
3,1,2-(Me <sub>3</sub> CNC) <sub>2</sub> {=C[N(H)CMe <sub>3</sub> ]-C(Me)=C(Me)[C(O)N(H)-(CMe <sub>3</sub> )]}Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR	[627]
3,1,2-(Me <sub>3</sub> CNC) <sub>4</sub> MeMo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>+</sup>	S, H, B, C, IR	[627]
<i>Exo,endo</i> -3,1,2-(C <sub>3</sub> H <sub>5</sub> )(CO)Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR	[628]
3,1,2-(CH <sub>2</sub> =CHR)(CO)Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> R=Me, Et	S, H, B, C, IR	[628]
Related derivatives		[628]
3,1,2-(R <sub>2</sub> C <sub>2</sub> ) <sub>2</sub> [P(OMe) <sub>3</sub> ] <sub>2</sub> Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R=Me, Ph	S, H, B, C, P	[629]
3,1,2-(RR'C <sub>2</sub> ) <sub>2</sub> [P(OMe) <sub>3</sub> ]Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R=Me; R'=Me, Ph	S, H, B, C, P	[629]
3,1,2-(Ph <sub>2</sub> C <sub>2</sub> ) <sub>2</sub> [P(OMe) <sub>3</sub> ] <sub>2</sub> Mo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8- <i>cis</i> -Ph <sub>2</sub> C <sub>2</sub> H)	S, X, H, B, C, P	[629]
3,1,2-(Ph <sub>2</sub> S) <sub>2</sub> (CO) <sub>2</sub> Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>2-</sup>	S, X, H, B, IR	[630]
3,1,2-O[O <sub>2</sub> Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> <sup>2-</sup>	S, X, H, B, IR	[630]
5,1,2-(C <sub>3</sub> H <sub>5</sub> )(CO) <sub>2</sub> Mo(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> (non-icosahedral)	S, X, H, B, IR	[631]
[(PhS) <sub>2</sub> Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> <sup>2-</sup>	S, X, H(2d), B	[632]
5,1,2-(C <sub>3</sub> H <sub>5</sub> )(CO) <sub>2</sub> Mo(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -3-SMe <sub>2</sub> )	S, X, H, B, IR	[634]
2,1,8-(C <sub>3</sub> H <sub>5</sub> )(CO) <sub>2</sub> Mo(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> - <i>n</i> -SMe <sub>2</sub> ) <i>n</i> =6, 11	S, X, H, B, IR	[634]
5,1,2-(CO) <sub>2</sub> (C <sub>3</sub> H <sub>5</sub> )Mo(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-SMe <sub>2</sub> )	S, X, H, B(2d)	[635]
8,1,2-(CO) <sub>2</sub> (C <sub>3</sub> H <sub>5</sub> )Mo(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -5-SMe <sub>2</sub> )	S, X, H, B(2d)	[635]
[3,1,2-(CO) <sub>3</sub> Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> <sup>2-</sup> N(PPh <sub>3</sub> ) <sub>2</sub> <sup>2+</sup> (Mo-Mo)	S, X, H, B, IR	[637]
[3,1,2-L <sub>2</sub> (μ-S)Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> <sup>2-</sup> N(PPh <sub>3</sub> ) <sub>2</sub> <sup>2+</sup> L=O, S Mo-Mo	S, X, H, B, IR	[637]
3,1,2-(CO) <sub>2</sub> (R <sub>2</sub> C <sub>9</sub> H <sub>5</sub> )Mo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) indenyl R=H, Me	S, X(Me), H, IR	[1491]
<b><i>Closo-MoC<sub>2</sub>B<sub>9</sub> clusters with exo-polyhedral metals</i></b>		
3,1,2-(Ph <sub>3</sub> Sn)(CO) <sub>3</sub> Mo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, B, IR, Sn	[599]
{3,1,2-(CO)Mo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-1-CH <sub>2</sub> PMeN[Mo(CO) <sub>4</sub> ]-P(C <sub>5</sub> H <sub>10</sub> N) <sub>2</sub> N} <sub>n</sub> <sup>2-</sup>	S, P, IR	[619]
3,1,2-(C <sub>9</sub> H <sub>7</sub> )Mo[P(OMe) <sub>3</sub> ](μ-MeC <sub>6</sub> H <sub>4</sub> C)(CO) <sub>2</sub> Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, P, IR	[621]
3,1,2-(C <sub>7</sub> H <sub>7</sub> )Mo(μ-MeC <sub>6</sub> H <sub>4</sub> C)(CO)(L)Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L=CO, PMe <sub>3</sub>	S, H, B, C, P, IR	[621]
3,1,2-(CO) <sub>7</sub> Mo <sub>2</sub> (μ-MeC <sub>6</sub> H <sub>4</sub> C)(CO) <sub>2</sub> Mo(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> R=Me, H	S, H, B, C, P, IR	[639]
3,1,2-(CO) <sub>2</sub> [(C <sub>6</sub> H <sub>4</sub> MeC≡W(CO)Cp]Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, P, IR	[624]
3,1,2-(C <sub>4</sub> Me <sub>4</sub> )(CO)Co(μ-MeC <sub>6</sub> H <sub>4</sub> C)(CO) <sub>2</sub> Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[641]
3,1,2-(μ-CO)[(Ph <sub>3</sub> ) <sub>2</sub> Rh][P(OMe) <sub>3</sub> ](MeC <sub>6</sub> H <sub>4</sub> C)Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, P	[620]
Cp* <sub>2</sub> Rh <sub>2</sub> (μ-H)(μ-CO)-3,1,2-(CO) <sub>2</sub> Mo(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=H, Me	S, H, B, C, IR	[642]
3,1,2-(Et <sub>3</sub> P) <sub>2</sub> Pt(CO) <sub>2</sub> Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -CH <sub>2</sub> -C <sub>6</sub> H <sub>4</sub> Me)	S, IR	[644]
3,1,2-(PMe <sub>2</sub> Ph)LPt(μ-CC <sub>6</sub> H <sub>4</sub> Me)(CO) <sub>2</sub> Mo(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) L=CO, PMe <sub>2</sub> Ph	S, H, B, C, P, IR	[644]
Related derivatives		[644]

Compound	Information	References
<b>Tungsten</b>		
<i>Closo-WCB<sub>10</sub> clusters</i>		
2,1-(CO) <sub>3</sub> (Ph <sub>3</sub> P)W(CB <sub>10</sub> H <sub>10</sub> -7-L) L = OEt <sub>2</sub> , O(CH <sub>2</sub> ) <sub>4</sub> , SMe <sub>2</sub> , S(CH <sub>2</sub> ) <sub>4</sub> , cyclic ethers and thioethers, NCCMe <sub>3</sub> , CNCMe <sub>3</sub> , CNC <sub>6</sub> H <sub>3</sub> Me <sub>2</sub>	S, H, B, C, P	[608]
2,1-(Me <sub>3</sub> CNC) <sub>3</sub> IW[(Me <sub>3</sub> CHN)CB <sub>10</sub> H <sub>10</sub> ]	S, H, B, C, P	[592]
2,1-(Me <sub>3</sub> CNC) <sub>2</sub> (CO)LW(CB <sub>10</sub> H <sub>11</sub> ) L = Me <sub>3</sub> C, CO	S, H, B, C, P	[611]
2,1-(CO) <sub>2</sub> (HC≡CCMe <sub>3</sub> )W(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup>	S, X, H, B, C, P, IR	[608]
2,1-(CO) <sub>3</sub> (Ph <sub>3</sub> P)W(CB <sub>10</sub> H <sub>10</sub> -7-OMeR) R = Et, <i>n</i> -C <sub>4</sub> H <sub>9</sub>	S, X( <i>n</i> -C <sub>4</sub> H <sub>9</sub> ), H, B, C, P, IR	[608]
Related WCB <sub>10</sub> complexes		[608,645]
<i>Closo-WC<sub>2</sub>B<sub>9</sub> clusters (no exo-polyhedral metals)</i>		
3,1,2-(CO) <sub>3</sub> MeW(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> R = H, Me	S, H, IR	[618]
3,1,2-(CO) <sub>3</sub> W(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-1-CH <sub>2</sub> -[cyclo-N <sub>3</sub> P <sub>3</sub> ](C <sub>5</sub> H <sub>10</sub> N) <sub>4</sub> Me	S, P, IR	[619]
3,1,2-μ-MeC <sub>6</sub> H <sub>4</sub> C=WC(O) <sub>2</sub> (PMe <sub>3</sub> )(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, C	[623]
3,1,2-(CO)L(RC≡CR)W(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L = Me <sub>2</sub> C <sub>2</sub> , Ph <sub>2</sub> C <sub>2</sub> , Ph <sub>3</sub> P, PMe <sub>2</sub> Ph; R = Me, Ph	S, H, B, C, P, IR	[624]
Related derivatives		[624]
2,1,8-(CO) <sub>3</sub> IW(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, B, C, IR	[617]
Related derivatives		[617]
2,1,8-(CO) <sub>2</sub> (PhC=CR)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-11-CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> Me R = Me, Ph	S, X(Me), H, B, C, IR	[646]
2,1,8-(CO) <sub>4</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-11-CH <sub>2</sub> -R R = <i>p</i> -C <sub>6</sub> H <sub>4</sub> Me, Me	S, H, B, C, IR	[647]
2,1,8-(CO) <sub>2</sub> (RC=CPh)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-11-CH <sub>2</sub> -C <sub>6</sub> H <sub>4</sub> Me R = Ph, Me	S, X(Me), H, B, C, IR	[647]
2,1,8-(CO) <sub>3</sub> XW(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> X = Cl, I	S, X(I), H, B, C, IR	[647]
3,1,2-(C <sub>3</sub> H <sub>5</sub> )(CO) <sub>2</sub> W(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -10-L) R = H, Me; L = OEt <sub>2</sub> , OC <sub>4</sub> H <sub>6</sub> , SMe <sub>2</sub> , Ph <sub>3</sub> P, NC <sub>5</sub> H <sub>5</sub> , NC <sub>5</sub> H <sub>4</sub> C <sub>5</sub> H <sub>4</sub> N	S, H, B, C, IR	[626]
3,1,2-(C <sub>3</sub> H <sub>5</sub> )(CO) <sub>2</sub> W(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -10-L) <sup>-</sup> L = OEt <sub>2</sub> , O(CH <sub>2</sub> ) <sub>4</sub> F	S, H, B, C, IR	[626]
3,1,2-(CO) <sub>4</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -10-L) L = OEt <sub>2</sub> , O(CH <sub>2</sub> ) <sub>4</sub> F	S, H, B, C, IR	[626]
3,1,2-(Me <sub>3</sub> CNC) <sub>4</sub> W(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R = Me, H	S, H, B, C, IR	[627]
3,1,2-(MeC≡CMe)(CO)W(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR	[627]
N(PPh <sub>3</sub> ) <sub>2</sub> <sup>+</sup> 3,1,2-(CO) <sub>2</sub> (C <sub>3</sub> H <sub>5</sub> )W(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, C, IR	[627]
3,1,2-(μ-CC <sub>6</sub> H <sub>4</sub> Me)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, C, IR	[648,649]
3,1,2-(μ-CC <sub>6</sub> H <sub>4</sub> Me)(CO) <sub>2</sub> W(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, C, IR	[649]
3,1,2-R(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> R = <i>o/p</i> -MeC <sub>6</sub> H <sub>4</sub> C, Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub> C	S, H, C, P, IR	[652]
3,1,2-(CO) <sub>2</sub> (μ-L)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> L = <i>n</i> -C <sub>4</sub> H <sub>9</sub> -C <sub>2</sub> C, MeC <sub>6</sub> H <sub>4</sub> C	S, H, B, C, IR	[621]
3,1,2-(PhC=CPh) <sub>2</sub> (CO)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-CH <sub>2</sub> R) R = C <sub>6</sub> H <sub>4</sub> Me, Me	S, H, B, C, P, IR	[653]
Related derivatives		[653]
3,1,2-(CO) <sub>2</sub> (RC)W(R'MeC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> R = C <sub>6</sub> H <sub>4</sub> Me, Me; R' = H, Me	S, H, C, IR	[654]
3,1,2-(CO) <sub>2</sub> (R <sub>2</sub> PCH <sub>2</sub> CH <sub>2</sub> PR <sub>2</sub> )W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-Et) R = Ph, Me	S, H, B, C, P, IR	[654]
3,1,2-(CO) <sub>2</sub> L <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -CH <sub>2</sub> R) L = CO, Ph <sub>3</sub> P, PHPh <sub>2</sub> , CN- <i>n</i> -C <sub>4</sub> H <sub>9</sub> , PhC <sub>2</sub> Ph; R = C <sub>6</sub> H <sub>4</sub> Me, C <sub>6</sub> H <sub>4</sub> OMe	S, H, B, C, P, IR	[655]

Continued



Compound	Information	References
3,1,2-(CO)(PhC=CPh) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -CH <sub>2</sub> R) R=C <sub>6</sub> H <sub>4</sub> Me, C <sub>6</sub> H <sub>4</sub> OMe	S, H, B, C, P, IR	[655]
Related derivatives		[655]
3,1,2-(CO) <sub>2</sub> [CpM(CO)](μ-CR)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -CH <sub>2</sub> R') M=W, Mo; R=Me, C <sub>6</sub> H <sub>4</sub> Me; R'=Me, C <sub>6</sub> H <sub>4</sub> Me	S, H, B, C, IR	[656]
3,1,2-(CO) <sub>2</sub> [CpW(CO)PMe <sub>3</sub> ](μ-CMe)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> Et)	S, H, B, C, P, IR	[656]
Related derivatives		[656]
3,1,2-(CO) <sub>2</sub> (RC)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> <sup>-</sup> R=Me, MeC <sub>6</sub> H <sub>4</sub> , MeOC <sub>6</sub> H <sub>4</sub> , MeOCH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )	S, H, B, C, P	[657]
3,1,2-(CO) <sub>2</sub> (Ph <sub>2</sub> PCH <sub>2</sub> PPh <sub>2</sub> )W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -μ-(CH <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )	S, H, B, C, P, IR	[657]
3,1,2-(CO)(PhC=CPh) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -μ-(CH <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )	S, H, B, C, P, IR	[657]
2,1,8-(CO) <sub>3</sub> (I)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -μ-(CH <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>4</sub> <sup>-</sup> )	S, X, H, B, C, P	[657]
2,1,8-(CO) <sub>2</sub> (Me <sub>3</sub> CC=CH)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -μ-(CH <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>4</sub> <sup>-</sup> )	S, H, B, C, P	[657]
2,1,8-(CO)(Me <sub>3</sub> CC=CH) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -μ-(CH <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>4</sub> <sup>-</sup> )	S, H, B, C, P	[657]
Related derivatives		[657]
3,1,2-μ-MeC <sub>6</sub> H <sub>4</sub> C=W(CO) <sub>2</sub> (PMe <sub>3</sub> )(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, C	[623]
3-W(CO) <sub>2</sub> (μ-H)[μ-C <sub>2</sub> (C <sub>6</sub> H <sub>4</sub> Me) <sub>2</sub> ](CO) <sub>2</sub> W(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup>	S, H, C, IR	[658]
3,1,2-(CO) <sub>2</sub> (μ-H)[μ-C(C <sub>6</sub> H <sub>4</sub> Me)]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, C, IR	[658]
3,1,2-(PMe <sub>3</sub> )(CO) <sub>2</sub> [μ-C(C <sub>6</sub> H <sub>4</sub> Me)]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, C, P, IR	[658]
3,1,2-(PMe <sub>3</sub> )(CO)[ROC=C(C <sub>6</sub> H <sub>4</sub> Me)]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=H, Me	S, H, C, P, IR	[658]
3,1,2-(CO) <sub>2</sub> (Me <sub>3</sub> C-C≡C-C)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, C, IR	[659]
2,1,8-(CO) <sub>2</sub> (Me <sub>3</sub> C-C≡C-C)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, C, IR	[659]
3,1,2-(NCMe <sub>3</sub> )(NHCMe <sub>3</sub> )(2,6-Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub> -O)W(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, MS	[660]
[3,1,2-(NCMe <sub>3</sub> )(NHCMe <sub>3</sub> ) <sub>2</sub> W(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> O]	S, H, B, C, MS	[660]
Related derivatives		[660]
<i>Closo</i> -WC <sub>2</sub> B <sub>9</sub> clusters with <i>exo</i> -polyhedral metal atoms		
{3,1,2-(CO)W(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-1-CH <sub>2</sub> -PMeN[Mo(CO) <sub>4</sub> ]P(C <sub>5</sub> H <sub>10</sub> N) <sub>2</sub> N <sub>2</sub> } <sub>n</sub> <sup>2-</sup>	S, P, IR	[619]
<i>Nido</i> [slipped]-(CO)(Et <sub>2</sub> C <sub>2</sub> )[(CO) <sub>2</sub> (indenyl)(μ-CHC <sub>6</sub> H <sub>4</sub> Me)Mo]-W-(η <sup>3</sup> -Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )	S, X, H, B, C, IR	[645]
Related derivatives		[645]
3,1,2-(C <sub>7</sub> H <sub>7</sub> )Mo(μ-MeC <sub>6</sub> H <sub>4</sub> C)(CO)(L)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L=CO, PMe <sub>3</sub>	S, H, B, C, P, IR	[621]
3,1,2-μ-C <sub>6</sub> H <sub>4</sub> Me-(CO)CpW-(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[640]
2,1,7-μ-C <sub>6</sub> H <sub>4</sub> Me-(CO)CpM-(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) M=Mo, W	S, X(W), H, B, C, IR	[640]
3,1,2-(CO) <sub>2</sub> [(C <sub>9</sub> H <sub>7</sub> )(CO)(μ-CC <sub>6</sub> H <sub>4</sub> Me)M]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) M=Mo, W; C <sub>9</sub> H <sub>7</sub> =indenyl	S, X, H, B, C, IR	[661]
3,1,2-(CO) <sub>2</sub> [(C <sub>9</sub> H <sub>7</sub> )(CO)(μ-CC <sub>6</sub> H <sub>4</sub> Me)Mo]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> C <sub>9</sub> H <sub>7</sub> =indenyl	S, H, B, C, IR	[661]
3,1,2-(CO) <sub>2</sub> [(C <sub>9</sub> H <sub>7</sub> )(CO)(μ-CC <sub>6</sub> H <sub>4</sub> Me)Mo]W(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) C <sub>9</sub> H <sub>7</sub> =indenyl	S, H, B, C, IR	[661]

Compound	Information	References
3,1,2-(CO) <sub>2</sub> [(C <sub>9</sub> H <sub>7</sub> )(PMe <sub>3</sub> )(CO)(μ-CC <sub>6</sub> H <sub>4</sub> Me)M]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) M=Mo, W; C <sub>9</sub> H <sub>7</sub> =indenyl	S, H, B, C, P, IR	[661]
3,1,2-(CO)(PMe <sub>3</sub> )[(C <sub>9</sub> H <sub>7</sub> )(CO)(μ-CC <sub>6</sub> H <sub>4</sub> Me)Mo]W(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) C <sub>9</sub> H <sub>7</sub> =indenyl	S, H, B, C, P, IR	[661]
3,1,2-(CO) <sub>2</sub> [(C <sub>9</sub> H <sub>7</sub> )L(μ-CC <sub>6</sub> H <sub>4</sub> Me)W]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) C <sub>9</sub> H <sub>7</sub> =indenyl L=PMe <sub>3</sub> , N <sub>2</sub> (C <sub>6</sub> H <sub>4</sub> Me) <sub>2</sub>	S, H, B, C, P, IR	[661,664]
Related derivatives		[661]
3,1,2-(C <sub>9</sub> H <sub>7</sub> )(PHPh <sub>2</sub> )Mo(μ-MeCC <sub>6</sub> H <sub>4</sub> C)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[662]
3,1,2-(C <sub>7</sub> H <sub>7</sub> )Mo(μ-HO-C≡C-C <sub>6</sub> H <sub>4</sub> Me)(μ-Ph <sub>2</sub> P)(CO)W(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=H, Me	S, X(R=H), H, B, C, P, IR	[662]
3,1,2-(C <sub>7</sub> H <sub>7</sub> )Mo(μ-Me-C≡C-C <sub>6</sub> H <sub>4</sub> Me)(μ-CH <sub>2</sub> )(μ-O)(OEt)W- (Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, H, B, C	[663]
3,1,2-(C <sub>9</sub> H <sub>7</sub> )Mo(CO)(μ- <i>n</i> -C <sub>4</sub> H <sub>9</sub> C <sub>2</sub> C)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[621]
3,1,2-(C <sub>7</sub> H <sub>7</sub> )Mo(μ-RC)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=MeC <sub>6</sub> H <sub>4</sub> , <i>n</i> -C <sub>4</sub> H <sub>9</sub> C <sub>2</sub> C	S, H, B, C, IR	[621]
Related derivatives		[621]
3,1,2-(CO) <sub>7</sub> Mo <sub>2</sub> (μ-MeC <sub>6</sub> H <sub>4</sub> C)(CO) <sub>2</sub> W(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> R=Me, H	S, H, B, C, P, IR	[639]
3,1,2-(CO) <sub>6</sub> (Me <sub>3</sub> P)Mo <sub>2</sub> (μ-MeC <sub>6</sub> H <sub>4</sub> C)(CO) <sub>2</sub> W(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, X, H, B, C, IR	[639]
Related derivatives		[639]
3,1,2-Cp(CO)W(MeC <sub>6</sub> H <sub>4</sub> )C(CO) <sub>2</sub> W(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R=H, Me	S, H, B, C, IR	[665]
3,1,2-μ-Cp(CO)W=CR-(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -CH <sub>2</sub> -R') R=C≡CCMe <sub>3</sub> , Me, C≡CC <sub>6</sub> H <sub>4</sub> Me; R'=C <sub>6</sub> H <sub>4</sub> Me, Me	S, H, B, C	[666]
3,1,2-[Cp(CO) <sub>2</sub> M=CR']Au[RC=](CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) M=Mo, W; R,R'=Me, C <sub>6</sub> H <sub>4</sub> Me	S, H, C, P, IR	[667]
3,1,2-[Cp(CO) <sub>2</sub> M=CR']Au[L <sub>2</sub> Pt(RC)(CO) <sub>2</sub> ]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) M=Mo, W; R, R'=Me, C <sub>6</sub> H <sub>4</sub> Me; L <sub>2</sub> =η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> , 2PMe <sub>2</sub> Ph	S, H, C, P, IR	[667]
3,1,2-[Cp*Co(CO) <sub>2</sub> CpW(RC)]Au[MeC=](CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=C <sub>6</sub> H <sub>4</sub> Me	S, H, C, IR	[668]
Related derivatives		[664]
3,1,2-μ-RC-(C <sub>7</sub> H <sub>7</sub> )Mo(CO)(RC≡CR')W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=C <sub>6</sub> H <sub>4</sub> Me, Me <sub>3</sub> CC≡C; R'=Me, Et	S, H, C, IR	[669]
Related derivatives		[669]
3,1,2-[MeC <sub>6</sub> H <sub>4</sub> CH-C <sub>3</sub> H(Me <sub>3</sub> C)]CpM(CO) <sub>2</sub> (CO) <sub>2</sub> W(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) M=Mo, W	S, X(W), H, C, IR	[659]
3,1,2-(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )(μ-RC)(CO)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -CH <sub>2</sub> -R) <sup>-</sup> R=C <sub>6</sub> H <sub>4</sub> Me, Me W-W	S, H, B, C, IR	[670]
Related derivatives		[670]
[3,1,2-(μ-SPh <sub>2</sub> ) <sub>2</sub> W(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ] <sup>n-</sup> W-W n=2,1,0	S, X, H, B, UV, IR, E, XPS	[636]
3,1,2-(CO) <sub>6</sub> Fe <sub>2</sub> (CR)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> ) <sup>-</sup> R=Me, Ph, <i>p</i> -MeC <sub>6</sub> H <sub>4</sub>	S, X(Ph), H, B, C, IR	[671]
3,1,2-(CO) <sub>6</sub> Fe <sub>2</sub> (μ-AuPPh <sub>3</sub> )(CMe)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> )	S, H, B, C, IR	[671]
3,1,2-[μ-(CO) <sub>6</sub> Fe <sub>2</sub> ](CO) <sub>2</sub> (μ-MeC <sub>6</sub> H <sub>4</sub> C)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> ) <sup>-</sup>	S, H, B, C, P	[620]
3,1,2-[μ-(CO) <sub>3</sub> Fe](μ-RCH)(μ-CO)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) <sup>-</sup> R=C <sub>6</sub> H <sub>3</sub> Me <sub>2</sub> , C <sub>6</sub> H <sub>4</sub> Me	S, X, H, IR	[672]
3,1,2-(μ-CO)(μ-R)(CO) <sub>2</sub> [μ-(CO) <sub>3</sub> Fe]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> R= <i>o/p</i> -MeC <sub>6</sub> H <sub>4</sub> C	S, H, B, C, IR	[652]

Continued

Compound	Information	References
Related derivatives		[652]
3,1,2-(CO) <sub>6</sub> Fe <sub>2</sub> (μ-CH=C=Cn-C <sub>4</sub> H <sub>9</sub> )(CO) <sub>2</sub> (μ-CO)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) <sup>-</sup>	S, X, H, B, C, IR	[673]
3,1,2-(CO) <sub>6</sub> Fe <sub>2</sub> (μ-CH=C=Cn-C <sub>4</sub> H <sub>9</sub> )(CO) <sub>2</sub> (μ-COH)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )	S, X, H, B, C, IR	[673]
3,1,2-(μ-R)(CO) <sub>2</sub> [Cp(L)Ru]W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=C <sub>6</sub> H <sub>3</sub> Me <sub>2</sub> ; L=CO, PMe <sub>2</sub> Ph	S, H, B(2d), C, P, IR	[652]
3,1,2-CpRu(μ-H)(CO)(μ-CMe)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )	S, H, B, C, IR	[674]
3,1,2-CpRu(CO)(μ-CC <sub>6</sub> H <sub>4</sub> Me)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )	S, X, H, B, C	[674]
Related derivatives		[674]
3,1,2-(C <sub>4</sub> Me <sub>4</sub> )(CO)Co(μ-RC)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=MeC <sub>6</sub> H <sub>4</sub> , Me	S, H, B, C, IR	[641]
3,1,2-[μ-(CO) <sub>4</sub> Co <sub>2</sub> ](μ-PhC)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, B, C, IR	[672]
3,1,2-[μ-(CO) <sub>6</sub> Co <sub>2</sub> ](μ-PhC)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, X, H, IR	[672]
(μ-CR)(CO) <sub>6</sub> Co <sub>2</sub> -3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> 2 B—H—Co R=Me, Ph, <i>m/p</i> -MeC <sub>6</sub> H <sub>4</sub>	S, H, B, C, IR	[675]
(Ph <sub>3</sub> P) <sub>2</sub> Rh(CO) <sub>2</sub> -3,1,2-(μ-CC <sub>6</sub> H <sub>4</sub> Me)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-R) R=H, C <sub>7</sub> H <sub>9</sub>	S, X, H, C, P, IR	[648]
(μ-CO)[(Ph <sub>3</sub> ) <sub>2</sub> Rh](CO)(MeC <sub>6</sub> H <sub>4</sub> C)-3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, P	[620]
(Ph <sub>2</sub> PCH <sub>2</sub> ) <sub>2</sub> Rh(CO) <sub>2</sub> (μ-CC <sub>6</sub> H <sub>4</sub> Me)-3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, C, P, IR	[649]
Cp* <sub>2</sub> Rh <sub>2</sub> (μ-H)(μ-CO)-(CO) <sub>2</sub> -3,1,2-W(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=H, Me	S, X(Me), H, B, C, IR	[642]
L <sub>3</sub> (H)Ir(μ-MeC <sub>6</sub> H <sub>4</sub> C)(CO) <sub>2</sub> -3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) L=Me <sub>3</sub> P, (MeO) <sub>3</sub> P	S, X, H, B, C, P, IR	[676]
Related derivatives		[676]
3,1,2-[μ-(PEt <sub>3</sub> ) <sub>2</sub> Ir](μ-RC)(μ-CO)(CO)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R= <i>p</i> -C <sub>6</sub> H <sub>4</sub> Me	S, H, B, P	[672]
3,1,2-[μ-(CO)(H)Ir](Ph <sub>2</sub> P <sub>2</sub> C <sub>2</sub> H <sub>4</sub> )(μ-RC)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) R= <i>p</i> -C <sub>6</sub> H <sub>4</sub> Me	S, H, B, C, P	[672]
3,1,2-[μ-(CO)(H)Ir](N <sub>2</sub> [C <sub>5</sub> H <sub>4</sub> ] <sub>2</sub> )(μ-RC)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) R= <i>p</i> -C <sub>6</sub> H <sub>4</sub> Me	S, H, B, C, P	[672]
Related derivatives		[672]
3,1,2-L <sub>2</sub> Pt=W(CO) <sub>3</sub> (Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L=Ph <sub>3</sub> P, Et <sub>3</sub> P	S, H, B, C, P, IR	[600]
"Hypercloso"-3,1,2-(Et <sub>3</sub> P) <sub>2</sub> (μ-CO) <sub>2</sub> Pt=W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -R) R=H, Et	S, H, B, C, P, IR	[677]
<i>Exo</i> -(Et <sub>3</sub> P) <sub>2</sub> Pt(B)(μ-H)(CO) <sub>2</sub> (PMe <sub>3</sub> )-3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> - <i>p</i> -CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> Me)	S, X, H, B, C, P, IR	[677]
<i>Exo</i> -(Et <sub>3</sub> P) <sub>2</sub> Pt(B)(μ-H)(CO) <sub>3</sub> -3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -R) R=H, Et	S, X, H, B, C, P, IR	[677]
Related derivatives		[677]
<i>Exo</i> -(μ-CR)(CO) <sub>2</sub> (η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Pt-3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> R=Ph, <i>p</i> -MeC <sub>6</sub> H <sub>4</sub>	S, H, C, P, Pt, IR	[650]
<i>Exo</i> -[μ-(C-C <sub>6</sub> H <sub>3</sub> Me <sub>2</sub> )(Et <sub>3</sub> P)(H)Pt](CO) <sub>2</sub> -3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) Pt—H—B	S, H, B, C, P, IR	[678]
<i>Exo</i> -[μ-(C-C <sub>6</sub> H <sub>3</sub> Me <sub>2</sub> )(Et <sub>3</sub> P)Pt](CO) <sub>2</sub> -3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) Pt—B two isomers	S, X(1 isomer), H, B, C, P, IR	[678]
<i>Exo</i> -(μ-CO)(μ-R)(CO)Pt-3,1,2-LW(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> R= <i>o/p</i> -MeC <sub>6</sub> H <sub>4</sub> C, Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub> C; L=η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> , (PMePh <sub>2</sub> ) <sub>2</sub>	S, H, C, P, Pt, IR	[652]
<i>Exo</i> -μ-(PhMe <sub>2</sub> P) <sub>2</sub> Pt-μ-MeC <sub>6</sub> H <sub>4</sub> C-3,1,2-(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) two isomers	S, X, H, B, C, P, IR	[679]

Compound	Information	References
<i>Exo</i> - $\mu$ -(PhMe <sub>2</sub> P) <sub>2</sub> Pt- $\mu$ -MeC <sub>6</sub> H <sub>4</sub> C-2,1,7-(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) two isomers	S, X, H, B, C, P, IR	[679]
$\mu$ -MCl- $\mu$ -RC-3,1,2-(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> M=Cu, Au; R=C <sub>6</sub> H <sub>4</sub> Me, Me <sub>3</sub> CC $\equiv$ C	S, H, C, IR	[680]
$\mu$ -(CuCl) <sub>2</sub> - $\mu$ -RC-3,1,2-(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> R=C <sub>6</sub> H <sub>4</sub> Me, Me <sub>3</sub> CC $\equiv$ C	S, H, B, C, IR	[680]
Related derivatives		[680]
3,1,2-( $\mu$ -Ph <sub>3</sub> PAu)(CO) <sub>2</sub> (MeC <sub>6</sub> H <sub>4</sub> C)W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, P	[620]
Related derivatives		[620]
3,1,2-( $\mu$ -Ph <sub>3</sub> PAu-MeC)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, H, C, P	[649]
3,1,2- $\mu$ , $\mu'$ -(CC <sub>6</sub> H <sub>4</sub> Me) <sub>2</sub> Au-[(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )] <sub>2</sub> <sup>-</sup>	S, H, C, IR	[649]
Related derivatives		[649]
3,1,2-(Ph <sub>3</sub> P)Au( $\mu$ -CC <sub>6</sub> H <sub>4</sub> Me)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, H, C, P	[648]
3,1,2-( $\mu$ -Ph <sub>3</sub> PAu)R(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R = <i>o/p</i> -MeC <sub>6</sub> H <sub>4</sub> C, Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub> C	S, H, C, IR	[652]
3,1,2-( $\mu$ -CO)( $\mu$ -R)(CO)( $\mu$ -Ph <sub>3</sub> PAu)PtLW(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R = Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub> C; L = $\eta^4$ -C <sub>8</sub> H <sub>12</sub>	S, H, C, P, Pt, IR	[652]
3,1,2-(Ph <sub>3</sub> P)Au( $\mu$ -RC)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R = <i>n</i> -C <sub>4</sub> H <sub>9</sub> C <sub>2</sub> , MeC <sub>6</sub> H <sub>4</sub>	S, H, B, C, IR	[639]
3,1,2-L <sub>2</sub> Pt(L')Au( $\mu$ -RC)(CO) <sub>2</sub> W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R = Me <sub>3</sub> CC $\equiv$ C, MeC <sub>6</sub> H <sub>4</sub> ; L <sub>2</sub> = (PhMe <sub>2</sub> P) <sub>2</sub> , $\eta^4$ -C <sub>8</sub> H <sub>12</sub> ; L' = Ph <sub>3</sub> P, Me <sub>2</sub> PhP	S, H, B, C, P, IR	[639]
[Ph <sub>2</sub> P-Au( $\mu$ -RC)(CO) <sub>2</sub> -3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )] <sub>2</sub> (CH=CH) two isomers	S, H, B, C, P, IR	[681]
[Ph <sub>2</sub> P-Au( $\mu$ -RC)( $\eta^4$ -C <sub>8</sub> H <sub>12</sub> )Pt(CO) <sub>2</sub> -3,1,2-W(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )] <sub>2</sub> CH <sub>2</sub> ) <sub>n</sub> n = 2-6; R = <i>p</i> -C <sub>6</sub> H <sub>4</sub> Me	S, H, B, C, P, IR	[681]
<b>Manganese</b>		
<i>Closo</i> -MnCB <sub>10</sub> clusters		
2,1-(H <sub>3</sub> NCB <sub>10</sub> H <sub>10</sub> ) <sub>2</sub> Mn <sup>2-</sup>	S, UV	[593]
2,1-(CO) <sub>3</sub> Mn(CB <sub>10</sub> H <sub>11</sub> ) <sup>2-</sup>	S	[682]
[2,1-(CO) <sub>3</sub> Mn](Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>10</sub> ) <sup>2-</sup>	S, H, B, C, IR	[683]
[2,1-(CO) <sub>3</sub> Mn](Me <sub>3</sub> CNH <sub>2</sub> )CB <sub>10</sub> H <sub>10</sub> ) <sup>-</sup>	S, H, B, C, IR	[683]
[2,1-(CO) <sub>3</sub> Mn](Me <sub>3</sub> CNH <sub>2</sub> )CB <sub>10</sub> H <sub>9</sub> ]- <i>exo</i> -Pt(Ph <sub>2</sub> PCH <sub>2</sub> CH <sub>2</sub> PPh <sub>2</sub> )	S, X, H, B, C, IR	[683]
<i>Closo</i> -MnMCB <sub>9</sub> clusters		
8,2,1-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>2</sub> PPh <sub>2</sub> ]Pt(CO) <sub>3</sub> Mn(PhCB <sub>9</sub> H <sub>9</sub> )	S, H, B, C, P	[359]
<i>Closo</i> -MnPCB <sub>9</sub> clusters		
2,1,7-(CO) <sub>3</sub> Mn(MePCB <sub>9</sub> H <sub>10</sub> )	S, H, IR	[684]
2,1,7-(Et <sub>3</sub> P)(CO) <sub>2</sub> Mn(MePCB <sub>9</sub> H <sub>10</sub> )	S, IR	[685]
<i>Closo</i> -MnC <sub>2</sub> B <sub>9</sub> clusters		
3,1,2-(CO) <sub>3</sub> Mn[(CH=CH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sup>-</sup>	S, H, IR	[686]
3,1,2-(CO) <sub>3</sub> Mn[(1,2-C <sub>4</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sup>-</sup> benzodicarbollide complex	S, H, UV	[687]
3,1,2-(CO) <sub>3</sub> Mn[(1,2-C <sub>4</sub> H <sub>6</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sup>-</sup> dihydrobenzodicarbollide complex	S, H, UV	[687]
{3,1,2-(CO) <sub>2</sub> BrMn[(1,2-C <sub>4</sub> H <sub>6</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]} <sub>2</sub> <sup>2-</sup> dihydrobenzodicarbollide complex	S, H, UV	[687]

Continued

Compound	Information	References
3,1,2-(CO) <sub>3</sub> Mn(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -1-C <sub>6</sub> H <sub>4</sub> R) <sup>-</sup> R=H, <i>m/p</i> -F	S, F, IR	[688]
3,1,2-(CO) <sub>3</sub> Mn(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-SMe <sub>2</sub> )	S	[707]
3,1,2-(CO) <sub>3</sub> Mn(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -OC <sub>4</sub> H <sub>8</sub> L) L=PPh <sub>3</sub> , NEt <sub>3</sub> , NC <sub>5</sub> H <sub>4</sub> Me, I <sup>-</sup>	S, H, B, C	[691]
3,1,2-(CO) <sub>3</sub> Mn(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-8-L R=H, L=O(CH <sub>2</sub> ) <sub>4</sub> , SMe <sub>2</sub> , NMe=CHMe, NMe=CHEt, NHMeEt, NH <sub>2</sub> Me, NMe=CHMe, NMe=CHEt; R=Me, L=O(CH <sub>2</sub> ) <sub>4</sub>	S, X(H, NMe=CHMe), B, C	[692]
3,1,2-(CO) <sub>3</sub> Mn(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-8-X <sup>-</sup> R=H, X=Cl, Br, I; R=Me, X=I	S, H, B, C	[692]
3,1,2-(NO)(CO) <sub>2</sub> Mn(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C	[692]
Related derivatives		[692]
<b>Technetium</b>		
<i>Closo-TcC<sub>2</sub>B<sub>9</sub> clusters</i>		
2,1,8-(CO) <sub>3</sub> <sup>99m</sup> Tc(8-RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=H, C <sub>5</sub> H <sub>4</sub> N, CH <sub>2</sub> -C <sub>5</sub> H <sub>4</sub> N	S (facile isomerization of 3,1,2 isomer)	[1418]
3,1,2-(CO) <sub>3</sub> <sup>99m</sup> Tc{[(CH <sub>2</sub> ) <sub>2</sub> C(O)OH](C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> radiopharmaceutical applications	S, H, B, C, IR, MS	[695]
	S(aqueous F <sup>-</sup> ), H, B, C, MS, radiochromatograms	[696]
3,1,2-(CO) <sub>3</sub> <sup>99m</sup> Tc{[(HO(O)C(CH <sub>2</sub> ) <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	S(microwave), H, B, C	[697]
2,1,7-(CO) <sub>3</sub> M{[HO(C(O))(CH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup> M= <sup>99</sup> Tc, <sup>99m</sup> Tc	S(aqueous F <sup>-</sup> ), H, B, C, IR, MS	[698]
3,1,2-(CO) <sub>3</sub> <sup>99m</sup> Tc{[Me <sub>2</sub> HN(CH <sub>2</sub> ) <sub>3</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	S(aqueous F <sup>-</sup> ), H, B, C, IR, MS	[698]
2,1,8-(CO) <sub>2</sub> (NO) <sup>99m</sup> Tc(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR, MS, UV	[699]
2,1,8-(CO) <sub>2</sub> (NO) <sup>99m</sup> Tc(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) in vivo/in vitro imaging	S, H, B, C, IR, MS, UV	[699]
2,1,8-(CO) <sub>2</sub> (NO) <sup>99m</sup> Tc{(PhCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> } R=H, CH <sub>2</sub> Ph in vivo/in vitro imaging	S, H, B, C, IR, MS, UV	[699]
2,1,7-(CO) <sub>3</sub> <sup>99m</sup> Tc (7-RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=(CH <sub>2</sub> ) <sub>n</sub> N(C <sub>2</sub> H <sub>4</sub> ) <sub>2</sub> NC <sub>6</sub> H <sub>4</sub> OMe <i>n</i> =1, 3 probe for α-adrenergic receptors	S, H, B, C, IR, MS	[1438]
2,1,8-(CO) <sub>3</sub> <sup>99m</sup> Tc(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> M=Re, <sup>99m</sup> Tc R=CH <sub>2</sub> - <i>cyclo</i> -1',4'-NC <sub>4</sub> H <sub>8</sub> N-C <sub>6</sub> H <sub>4</sub> - <i>o</i> -OMe, C(O)ONH(CH <sub>2</sub> ) <sub>4</sub> -1',4'-NC <sub>4</sub> H <sub>8</sub> N-C <sub>6</sub> H <sub>4</sub> - <i>o</i> -OMe molecular imaging probes	S	[1465]
<b>Rhenium</b>		
<i>Closo-ReCB<sub>10</sub> clusters</i>		
2,1-(CO) <sub>3</sub> Re[(Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>10</sub> (N-Re)	S, H, B, C, P, IR	[683]
[2,1-(CO) <sub>3</sub> Re[(Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>10</sub> ] <sup>2-</sup>	S, H, B, C, IR	[683]
2,1-(CO) <sub>2</sub> (L)[(Ph <sub>2</sub> PCH <sub>2</sub> CH <sub>2</sub> PPh <sub>2</sub> )M]Re(CB <sub>10</sub> H <sub>11</sub> ) M=Pt, L=CO; M=Pd, L=CO, CCM <sub>3</sub> , CNC <sub>6</sub> H <sub>3</sub> Me <sub>2</sub>	S, X(Pd, CO), H, B, C, P	[700]
2,1-Cp*M(CO) <sub>3</sub> Re(CB <sub>10</sub> H <sub>11</sub> ) M=Rh, Ir	S, X(Rh), H, B, C, P, IR	[701]
2,1-(CO) <sub>3</sub> Re[(Me <sub>3</sub> CHN)CB <sub>10</sub> H <sub>10</sub> ]	S, H, B, C, IR	[702]
2,1-(CO) <sub>2</sub> LRe[(Me <sub>3</sub> CHN)CB <sub>10</sub> H <sub>10</sub> ] L=Me <sub>3</sub> CC≡C, NMe <sub>3</sub>	S, H, B, C, IR	[702]
2,1-(CO) <sub>2</sub> LRe[(Me <sub>3</sub> CHN)CB <sub>10</sub> H <sub>9</sub> -3-C(=CHCMe <sub>3</sub> )/CH=CHCMe <sub>3</sub> ]	S, X, H, B, C, IR	[702]
2,1-(CO) <sub>3</sub> Re[(Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>10</sub> ] <sup>2-</sup>	S, H, B, C, IR	[703]
2,1-(CO) <sub>2</sub> Re[(Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>9</sub> ]-3-C=CH(CMe <sub>3</sub> )-CH=CHCMe <sub>3</sub> , Ph	S, H, B, C, IR	[703]

Compound	Information	References
<i>Closo-ReC<sub>2</sub>B<sub>9</sub> clusters (no exo-polyhedral metals)</i>		
3,1,2-(CO) <sub>3</sub> Re(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> R=CH <sub>2</sub> C <sub>5</sub> H <sub>4</sub> N, CH <sub>2</sub> Ph	S, H, B, C, IR, MS	[705]
2,1,8-(CO) <sub>3</sub> Re(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> R=H, CH <sub>2</sub> C <sub>5</sub> H <sub>4</sub> N, Ph	S, H, B, C, IR, MS	[705]
2,1,8-(CO) <sub>3</sub> Re(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=CH <sub>2</sub> C <sub>5</sub> H <sub>4</sub> NMe, CH <sub>2</sub> C <sub>5</sub> H <sub>4</sub> NH	S, X, H, B, C, IR, MS	[705]
2,1,8-(CO) <sub>3</sub> Re(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, H, B, C, binding affinity for estrogen receptor	[694]
2,1,8-(NO)(CO) <sub>2</sub> Re[(HOC <sub>6</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, C, binding affinity for estrogen receptor	[694]
3,1,2-(CO) <sub>3</sub> Re[( <i>cyclo</i> -C <sub>6</sub> H <sub>11</sub> )CH <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup>	S, H, B, C, IR, MS	[705]
3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-SMe <sub>2</sub> )	S	[707]
3,1,2-(CO) <sub>3</sub> Re[(glucosyl-CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup>	S, H, B, C, IR, MS	[708]
3,1,2-(CO) <sub>3</sub> Re[(HNC <sub>5</sub> H <sub>4</sub> -CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, H, B, C, IR, MS	[708]
3,1,2-(CO) <sub>3</sub> Re[[Me <sub>2</sub> HN(CH <sub>2</sub> ) <sub>3</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, C, IR, MS	[708]
3,1,2-(CO) <sub>3</sub> Re[(C <sub>6</sub> H <sub>5</sub> )( <i>p</i> -HOC <sub>6</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, B, C, IR, MS	[708]
2,1,7-(CO) <sub>3</sub> Re{[HO(CO)] <sub>2</sub> (C <sub>6</sub> H <sub>5</sub> CH <sub>2</sub> ) <sub>2</sub> H <sub>2</sub> N <sub>2</sub> }C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup>	S, H, B, C, IR, MS	[708]
3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-R R=OC <sub>4</sub> H <sub>8</sub> , OC <sub>4</sub> H <sub>8</sub> O	S, H, B(2d), C, IR	[1594]
3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-[O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> X <sup>-</sup> X=Cl, Br, I, OCH <sub>2</sub> Ph, N <sub>3</sub> , OH, NH <sub>3</sub> <sup>+</sup>	S, H, B(2d), C, IR	[1594]
3,1,2-(NO)(CO) <sub>2</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-[O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> X X=Cl, Br, I, OCH <sub>2</sub> Ph, N <sub>3</sub> , OH, NH <sub>3</sub> <sup>+</sup> peptide bioconjugates for drug delivery	S, X(X=I), H, B(2d), C, IR	[1594]
2,1,8-(CO) <sub>3</sub> Re(RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> R=H, <i>p</i> -C <sub>6</sub> H <sub>4</sub> OH; R'= <i>p</i> -C <sub>6</sub> H <sub>4</sub> OH, Ph	S(microwave), H, B, C	[697]
3,1,2-(CO) <sub>3</sub> <sup>186,188</sup> Re{[(CH <sub>2</sub> ) <sub>2</sub> C(O)OH](C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> radiopharmaceutical applications	S, H, B, C, IR, MS	[695]
3,1,2-(CO) <sub>3</sub> Re(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> R=H, (CH <sub>2</sub> ) <sub>2</sub> C(O)OH, (CH <sub>2</sub> ) <sub>3</sub> NHMe	S(aqueous F <sup>-</sup> ), H, B, C, MS, radiochromatograms	[696]
2,1,8-(CO) <sub>2</sub> (NO)Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR, MS, UV, fluorescence	[699]
2,1,8-(CO) <sub>2</sub> (NO)Re(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) in vivo/in vitro imaging	S, H, B, C, IR, MS, UV, fluorescence	[699]
2,1,8-(CO) <sub>2</sub> (NO)Re[(PhCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] R=H, CH <sub>2</sub> Ph in vivo/in vitro imaging	S, H, B, C, IR, MS, UV, fluorescence	[699]
3,1,2-(NO)(CO) <sub>2</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR	[709]
3,1,2-(NO)LL'Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L, L'=CO, PPh <sub>3</sub> , PMe <sub>3</sub> , CNCMe <sub>3</sub> , CNxyl	S, H, B, C, IR	[709]
Other related derivatives		[709]
3,1,2-(ON)(OC)BrRe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, C, IR, UV	[710]
3,1,2-(CO) <sub>3</sub> Re[(CHR)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] R=H, Me	S	[711]
3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -1-C <sub>6</sub> H <sub>4</sub> R) <sup>-</sup> R=H, <i>m</i> , <i>p</i> -F	S, F, IR	[688]
3,1,2-(NO)(CO) <sub>2</sub> Re(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=H, Me	S, IR, UV	[712]
3,1,2-(NO)LL'Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=CO; L'=NCC <sub>6</sub> H <sub>3</sub> Me <sub>2</sub> , PMe <sub>3</sub> ; L, L'=NCC <sub>6</sub> H <sub>3</sub> Me <sub>2</sub>	S, IR, UV	[712]
3,1,2-(CO) <sub>2</sub> (NO)Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, UV, photoluminescence	[365,713]
2,1,7-(CO) <sub>3</sub> Re(7-RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=(CH <sub>2</sub> ) <sub>n</sub> N(C <sub>2</sub> H <sub>4</sub> ) <sub>2</sub> NC <sub>6</sub> H <sub>4</sub> OMe <i>n</i> =1, 3 probe for α-adrenergic receptors	S, H, B, C, IR, MS	[1438]
2,1,8-(CO) <sub>3</sub> Re(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> R=CH <sub>2</sub> - <i>cyclo</i> -1',4'-NC <sub>4</sub> H <sub>8</sub> N-C <sub>6</sub> H <sub>4</sub> - <i>o</i> -OMe, C(O)ONH(CH <sub>2</sub> ) <sub>4</sub> -1',4'-NC <sub>4</sub> H <sub>8</sub> N-C <sub>6</sub> H <sub>4</sub> - <i>o</i> -OMe molecular imaging probes	S, X	[1465]

Continued

Compound	Information	References
3,1,2-(CO) <sub>3</sub> Re{[RNH—C(=NH <sub>2</sub> )NHCH <sub>2</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> } <sup>-</sup> R=H, Et guanidiny derivatives	S, H, B, C, IR, MS	[1470]
<b>Closo-ReC<sub>2</sub>B<sub>9</sub> clusters with exo-polyhedral metal atoms</b>		
3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-exo-(μ-H) <sub>3</sub> -RuCl(PPh <sub>3</sub> ) <sub>2</sub>	S, X, H, B, C, P, IR	[714]
3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-exo-(μ-H) <sub>2</sub> -RhL <sub>2</sub> L <sub>2</sub> =(PPh <sub>3</sub> ) <sub>2</sub> , Fe(C <sub>5</sub> H <sub>4</sub> PPh <sub>2</sub> ) <sub>2</sub>	S, H, B, C, P, IR	[714]
3,1,2-(CO) <sub>3</sub> (Ph <sub>3</sub> PM)Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) M=Cu, Ag	S, H, B, C, P, IR	[714]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>2</sub> PPh <sub>2</sub> ]Pt—(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>+</sup> Pt—H—B 2 isomers	S, X, H, B, C, P, IR	[715]
Related derivatives		[715]
3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )—Ag(C <sub>3</sub> H <sub>3</sub> N <sub>2</sub> ) <sub>2</sub> CH <sub>2</sub> (pyrazolyl)	S, X, H, B, C	[716]
[3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )—Ag] <sub>2</sub> (C <sub>3</sub> H <sub>3</sub> N <sub>2</sub> ) <sub>3</sub> CH (pyrazolyl)	S, X, H, B, C	[716]
[3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )—Ag] <sub>4</sub>	S, X, H, B, C	[716]
Related derivatives		[716]
<b>Iron</b>		
<b>Closo-FeCB<sub>10</sub> clusters</b>		
2,1-(H <sub>3</sub> NCB <sub>10</sub> H <sub>10</sub> ) <sub>2</sub> Fe <sup>-</sup>	S, UV	[593]
2,1-(CO) <sub>3</sub> Fe(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, C	[609]
2,1-(CO) <sub>2</sub> LFe(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup> L=Ph <sub>3</sub> P, CNCMe <sub>3</sub>	S, H, B, C, P	[609]
2,1-CpFe[(THF)CB <sub>10</sub> H <sub>10</sub> ]	S, B, IR, MS	[318]
2,1-CpFe[(Et <sub>2</sub> O)CB <sub>10</sub> H <sub>10</sub> ]	S, X, B, IR, MS	[318]
2,1-(CO) <sub>2</sub> Fe(CB <sub>10</sub> H <sub>10</sub> )-3-(CH=CHCMe <sub>3</sub> ) <sup>-</sup>	S, H, B, C, P, IR	[717]
2,1-(CO) <sub>2</sub> Fe(CB <sub>10</sub> H <sub>9</sub> )-3-(CH=CHCMe <sub>3</sub> )-5-(NMe=CHMe)	S, X, H, B, C, P, IR	[717]
Related derivatives		[717]
2,1-(CO) <sub>3</sub> Fe(CB <sub>10</sub> H <sub>8</sub> )-(μ-H) <sub>3</sub> -Cu(PPh <sub>3</sub> )	S, H, B, C, P	[610]
2,1-(CO) <sub>3</sub> Fe(CB <sub>10</sub> H <sub>8</sub> )-(μ-H) <sub>2</sub> -Ag(PPh <sub>3</sub> )	S, H, B, C, P	[610]
2,1-(CO) <sub>2</sub> LFe(CB <sub>10</sub> H <sub>10</sub> )-(μ-H)-Fe(CO) <sub>2</sub> Cp* L=CO, PPh <sub>3</sub>	S, H, B, C, P	[610]
<b>Closo-FeC<sub>2</sub>B<sub>9</sub> clusters</b>		
<b>Mono(dicarbollyl) Fe complexes</b>		
3,1,2-CpFe(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-4-SMe <sub>2</sub>	S, H, B, E, Mössbauer	[718]
3,1,2-CpFe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) R=OH, OCOCF <sub>3</sub>	S, X(OCOCF <sub>3</sub> ), IR	[724]
3,1,2-CpFe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-Br)	S	[725]
3,1,2-CpFe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-HgCl)	S, B, IR	[725]
3,1,2-CpFe(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -8-R-9-R'-12-R'') R, R', R''=H, H, H; OH, H, H; Cl, H, H; Br, H, H; H, Br, H; Br, Br, H; H, Br, Br	MS(detailed)	[726]
3,1,2-CpFe[(C <sub>3</sub> H <sub>3</sub> R)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] R=H, Ph	S, H, C, IR	[727]
3,1,2-CpFe[(CH=CHCH <sub>2</sub> Ph)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, IR, MS	[727]
3,1,2-CpFe(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> R=CH=CHMe, CH <sub>2</sub> CH=CH <sub>2</sub>	S, H, IR	[727]
3,1,2-CpFe[C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> (SMe <sub>2</sub> )Br <sub>2</sub> ]	S, X	[729]
3,1,2-CpFe[(LCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] L=Ph <sub>3</sub> P, Me <sub>2</sub> S, NC <sub>5</sub> H <sub>5</sub>	S	[1368]
3,1,2-CpFe(MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, IR, MS	[732]

Compound	Information	References
3,1,2-CpFe(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=CH <sub>2</sub> OH, COOH, CH <sub>2</sub> Cl, CHO, CH <sub>2</sub> COOH, CH <sub>2</sub> OAc, CH <sub>2</sub> OEt, CH <sub>2</sub> OCH <sub>2</sub> CH <sub>2</sub> OMe	S, IR	[732]
3,1,2-CpFe[CH(OH)R]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> R=Me, Ph	S, IR	[732]
3,1,2-CpFe[[C(O)Me]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, IR, MS	[732]
3,1,2-CpFe[[C(O)CH <sub>2</sub> D]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, MS	[732]
3,1,2-CpFe[(HC≡C)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, IR	[734]
3,1,2-CpFe[(HC≡C)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup>	S, IR	[734]
3,1,2-CpFe[(MeCO)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, IR	[734]
3,1,2-CpFe[[CH <sub>2</sub> =C(OMe)]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, IR	[734]
3,1,2-CpFe(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=H <sub>2</sub> C=C, CMeCl, CMeBr	S, H, IR	[734]
3,1,2-CpFe[( <i>m/p</i> -FC <sub>6</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S	[735]
3,1,2-CpFe(C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-4-SMe <sub>2</sub> -8-OC(O)CF <sub>3</sub> , <sub>1,2</sub> -12-HgCl	S, H, B(2d), F	[736]
3,1,2-CpFe(C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-4-SMe <sub>2</sub> -8,9-[OC(O)CF <sub>3</sub> ] <sub>2</sub>	S, H, B(2d), F	[736]
3,1,2-CpFe(C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> )-4-SMe <sub>2</sub> -7,8-[OC(O)CF <sub>3</sub> ] <sub>2</sub> -12-HgCl	S, X	[736]
3,1,2-Cp*Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+L</sup> L = 2,3-dichloro-5,6-dicyano- <i>p</i> -benzoquinone and 7,7',8,8'-tetracyano-2,3,5,6-tetrafluoroquinodimethane charge-transfer salts	S, X, IR, E, MAG	[738]
3,1,2-(η <sup>6</sup> -C <sub>5</sub> R <sub>5</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) R=H, Me	S, H, B	[739]
3,1,2-[η <sup>6</sup> -CH <sub>2</sub> =C(Me)C <sub>5</sub> H <sub>4</sub> ]Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) carbonium ion-carborane complex	S, IR	[740]
3,1,2-[η <sup>6</sup> -CH <sub>2</sub> =C(Me)C <sub>5</sub> H <sub>4</sub> ]Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> carbonium ion-carborane complex	S, H, IR	[740]
3,1,2-[η <sup>6</sup> -Me <sub>2</sub> (O)Me <sub>2</sub> CC <sub>5</sub> H <sub>4</sub> ]Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, IR	[740]
3,1,2-[η <sup>6</sup> -Me <sub>2</sub> (S)Me <sub>2</sub> CC <sub>5</sub> H <sub>4</sub> ]Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, IR	[740]
1,3-[3,1,2-(η <sup>6</sup> -C <sub>5</sub> H <sub>4</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> -R R=C(Me) <sub>2</sub> CH=CMe, C(Me) <sub>2</sub> CH <sub>2</sub> C(Me)SMe <sub>2</sub>	S, IR	[740]
3,1,2-(η <sup>5</sup> -C <sub>6</sub> H <sub>7</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )	S, H, B	[742]
3,1,2-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+</sup>	S, H, B	[742]
3,1,2-(η <sup>5</sup> -C <sub>6</sub> H <sub>7</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-L R=H, Me; L=SMe <sub>2</sub> , NMe <sub>3</sub> )	S, X(H,SMe <sub>2</sub> ), H, B, E, Mössbauer (H,SMe <sub>2</sub> )	[743]
3,1,2-L <sub>3</sub> Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+</sup> L=Me <sub>3</sub> CNC, P(OMe) <sub>3</sub>	S, X(H, SMe <sub>2</sub> ; H, NMe <sub>3</sub> ), H, B, P, E, Mössbauer (H, SMe <sub>2</sub> )	[743]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sub>2</sub>	S, X, H, NMe <sub>3</sub> , H, B, P, E, Mössbauer	[743]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-NMe <sub>2</sub> )	S, X, H, NMe <sub>3</sub> , H, B, P, E, Mössbauer	[743]
3,1,2-(η <sup>6</sup> -MeRC <sub>6</sub> H <sub>4</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R=H, Me	S, X, H, B	[744]
3,1,2-(η <sup>6</sup> -1,3,5-C <sub>6</sub> Me <sub>6</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, IR, MS	[745]
2,1,7-(η <sup>6</sup> -L)Fe(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L=C <sub>6</sub> H <sub>6</sub> , 1,3-C <sub>6</sub> H <sub>4</sub> Me <sub>2</sub> , C <sub>10</sub> H <sub>8</sub>	S, H, B, IR	[746]
CpFe( <i>cyclo</i> -P <sub>5</sub> )Fe[2,4,12-(N-C <sub>4</sub> H <sub>9</sub> NH)C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ] <sup>+</sup> 3,1,2-Co(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> triple-decker P <sub>5</sub> complex	S, X, H, B	[1422]
3,1,2-(η <sup>5</sup> -C <sub>9</sub> H <sub>7</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) C <sub>9</sub> H <sub>7</sub> =indenyl	S, X	[748]
3,1,2-(CO)LL'Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L,L'=CO, PPh <sub>3</sub> ; PPh <sub>3</sub> , MeCN; CO, MeCN; CO, P(OMe) <sub>3</sub> ; P(OMe) <sub>3</sub> ; CO	S, X[PPh <sub>3</sub> , PPh <sub>3</sub> /MeCN, P(OMe) <sub>3</sub> , CO], H, B, IR	[750]
3,1,2-(Me <sub>3</sub> CNC) <sub>3</sub> Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+</sup>	S, X, H, B	[742]

Continued



Compound	Information	References
3,1,2-(CO) <sub>2</sub> Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>2-</sup>	S, B	[751]
3,1,2-(CO) <sub>2</sub> Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> L = Me, PhCH <sub>2</sub> , Ph <sub>3</sub> Sn, COMe	S, X(Ph <sub>3</sub> Sn), H, B, IR	[751]
3,1,2-(CO)(η <sup>3</sup> -CH <sub>2</sub> CRCH <sub>2</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> R = H, Me	S, X(H), H, B, IR	[751]
3,1,2-(CO)(COMe)(PMe <sub>3</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, X, H, B, IR	[751]
3,1,2-L <sub>2</sub> Fe[(Me <sub>2</sub> NCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] (N → Fe) L = CO, CNCMe <sub>3</sub> , PMe <sub>3</sub> ; L <sub>2</sub> = COD	S, X(CO), H, B, C, P(PMe <sub>3</sub> ), IR	[752]
2,1,7-(CO) <sub>3</sub> Fe(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, IR	[746,747]
2,1,7-[(MeO) <sub>3</sub> P] <sub>3</sub> Fe(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, P	[746,747]
3,1,2-LFe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L = 1',3',5'-C <sub>6</sub> H <sub>3</sub> Me <sub>3</sub> , C <sub>5</sub> Me <sub>5</sub> H	thermal rearrangement	[1449]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>n</sub> PPh <sub>2</sub> ]ClFe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) n = 2,3 catalysis of polymerization of methyl methacrylate (MMA) and styrene (n = 3)	S, X, ESR, IR	[1505]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>n</sub> PPh <sub>2</sub> ]ClFe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) n = 2,3 paramagnetic	E, MS(MALDO-TOF)	[1558]
<b>Bis(dicarbollyl) Fe complexes</b>		
3-Fe(1,2 - C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> Na <sup>+</sup> ·2L L = bipyridyl, 1,10-phenanthroline	S, IR	[757]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-C <sub>4</sub> H <sub>8</sub> O <sub>2</sub> ) dioxane	S, X, H, B, C	[758]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OCH <sub>2</sub> OCH <sub>2</sub> -L) L = NC <sub>5</sub> H <sub>5</sub> , PPh <sub>3</sub> , OH, 2-O(OC <sub>6</sub> H <sub>4</sub> Me)	S, X[NC <sub>5</sub> H <sub>5</sub> , PPh <sub>3</sub> , 2-O(OC <sub>6</sub> H <sub>4</sub> Me)], H, B, C	[758]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SEt <sub>2</sub> ) <sub>2</sub>	S, H, B, IR, UV, E	[759]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SEt <sub>2</sub> ) <sub>2</sub>	S, H, B, IR, UV	[759]
3-Fe[1,2 - (C <sub>4</sub> H <sub>3</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> <sup>-</sup>	S, H, B, IR, UV, MS, E, COND, MAG	[760]
3-Fe[1,2 - (m/p - FC <sub>6</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> <sup>-</sup>	F	[761]
3-Fe(1,2 - Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> <sup>-</sup>	S, IR, UV, MAG, E	[618]
	ESR	[723]
[3-(μ - CO) <sub>2</sub> Fe(1,2 - C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> <sup>2-</sup>	S, B, IR	[597]
	X	[765]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )-3-Fe(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> N <sub>3</sub> <sup>-</sup> nucleoside conjugate	S(dipolar addition [chemical ligation]), H, B, IR, MS, UV	[766]
3-Fe(1,2 - R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> <sup>2-</sup> R = Me, Ph	H, UV	[618]
3-CpFe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )-3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, B, H, IR, UV, E	[767]
[Cp* <sub>2</sub> ThR] <sub>2</sub> <sup>2+</sup> Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>2-</sup> R = H, Me, SiMe <sub>3</sub>	S, X, H, C	[768]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-C <sub>5</sub> H <sub>4</sub> NCOMe) <sub>2</sub>	S, H, B, MS	[762]
2-Fe(1,7-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -11-C <sub>5</sub> H <sub>5</sub> N) <sub>2</sub>	S, H, B, IR, MS	[762]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-NEt <sub>3</sub> ) <sub>2</sub>	S, X, H, B, IR, MS	[762]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-Cl) <sub>2</sub> <sup>-</sup> extraction agent for Cs <sup>+</sup> and Sr <sup>2+</sup>	S	[769]
	X	[770]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sub>2</sub> (meso, dd, ll)	S, X, H, B, IR, UV, E	[773]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sub>2</sub> <sup>+</sup> DDQ <sup>-</sup> DDQ = 2,3-Cl <sub>2</sub> -5,6-(CN) <sub>2</sub> -p-benzoquinone	S, X, IR, COND	[773]
Meso-commo-3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sub>2</sub>	S, X, IR	[774]

Compound	Information	References
<i>dd/ll</i> -3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sub>2</sub> <sup>+</sup>	S	[774]
(TTF) <sup>+</sup> [3-Fe[1,2-(C <sub>4</sub> H <sub>3</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> ] <sub>2</sub> <sup>-</sup> TTF=tetrathiafulvalenium	S, H, B, IR, UV, MAG, COND, MS	[760]
(ET) <sub>2</sub> <sup>+</sup> [3-Fe[1,2-(C <sub>4</sub> H <sub>3</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> ] <sub>2</sub> <sup>-</sup> ET=bis(ethylenedithio)tetrathiafulvalenium	S, X, IR, UV, MAG, COND	[603]
[3-Fe <sup>III</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ] <sub>2</sub> Fe <sup>II</sup> -4L L=bipyridyl, 1,10-phenanthroline	S, IR	[757]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )-3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SEt <sub>2</sub> )	S, B, ESR	[377]
(C <sub>5</sub> H <sub>4</sub> NMe) <sub>2</sub> <sup>2+</sup> 2[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Fe(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-X)] <sup>-</sup> X=H, I probe for sequential voltage tuning	S, E	[1580]
(C <sub>5</sub> H <sub>4</sub> NMe) <sub>2</sub> <sup>2+</sup> 2[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11-n</sub> ) <sub>2</sub> Fe <sup>-</sup> n=0,1 probe for sequential voltage tuning	S, E	[1580]
2-Fe(1,7-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(8,8')-N <sub>2</sub> C <sub>3</sub> H <sub>3</sub>	S, X, H, B, MS, E	[775]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ-CHS <sub>2</sub>	S, IR	[776]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(8,8')-OMe	X	[777]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -4-SMe <sub>2</sub> -12-HgCl) <sub>2</sub>	S, X, H, B	[778]
3-Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -8,8'-Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>2</sub> PPh <sub>2</sub>	S, X, H, B, P, E	[1505]
3-Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -8,8'-Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>2</sub> PPh <sub>2</sub> paramagnetic	E, MS(MALDO-TOF)	[1558]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-Br) <sub>2</sub> <sup>-</sup> diffuse redox mediators for glucose oxidase bioanodes	S, H, B, MS	[1519]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-Br) <sub>2</sub> <sup>-</sup> diffuse redox mediators for glucose oxidase bioanodes	S, H, B, MS	[1519]
Na <sup>+</sup> [3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-[O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> [O-(CH <sub>2</sub> ) <sub>4</sub> ] <sub>n</sub> OH)] <sup>-</sup> n=1-22 redox-active polymers for surface modification	S, H, B, C, IR, MS, E	[1551]
Na <sup>+</sup> [3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(8,8')-(CH <sub>2</sub> ) <sub>4</sub> [O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> O] <sup>-</sup>	S, X, H, B, C, IR, MS, E	[1551]
3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-R) <sub>2</sub> <sup>-</sup> R=OC <sub>4</sub> H <sub>4</sub> <sup>+</sup> , O-(CH <sub>2</sub> ) <sub>4</sub> NHC <sub>4</sub> H <sub>8</sub> O <sup>+</sup> , [O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> NC <sub>4</sub> H <sub>8</sub> O, O-(CH <sub>2</sub> ) <sub>4</sub> PPh <sub>2</sub> (CH <sub>2</sub> ) <sub>2</sub> PPh <sub>2</sub> <sup>+</sup>	S, X, H, B, C, IR, MS	[1554]
MeC <sub>6</sub> H <sub>4</sub> N-C <sub>5</sub> H <sub>11</sub> <sup>+</sup> 3-Fe(1,2-Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> <sup>-</sup> cocatalyst with Pd(0) nanoparticles for oxidation of benzyl alcohol and lignin to form aromatic aldehydes	S, H, B, C, IR, MS	[1593]
<b>Closo-FePCB<sub>9</sub> clusters</b>		
3-Fe(1,2-MePCB <sub>9</sub> H <sub>10</sub> ) <sub>2</sub>	S, H, IR, UV, E	[684]
2-Fe(1,7-MePCB <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> two isomers	S, H, B, IR, UV, E	[684]
	X	[779]
2-Fe(1,7-PCB <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> <sup>2-</sup>	S, H, IR, UV, E	[684]
(1,7-MePCB <sub>9</sub> H <sub>10</sub> )-2-Fe(1,7-PCB <sub>9</sub> H <sub>10</sub> ) <sup>-</sup>	S, H, IR, E	[684]
2,1,7-CpFe[(Me)PCB <sub>9</sub> H <sub>10</sub> ]	S, H, UV, MS	[684]
2-Fe[1,7-[(CO) <sub>5</sub> M]PCB <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> <sup>2-</sup> M=Cr, Mo	S, H, B, IR	[780]
<b>Closo-FeSCB<sub>9</sub> clusters</b>		
1,2,4-(C <sub>4</sub> Me <sub>4</sub> )Fe(SCB <sub>9</sub> H <sub>9</sub> )-4-NHCMe <sub>3</sub>	S, X, H, B	[392]
<b>Closo-FePC<sub>2</sub>B<sub>8</sub> clusters</b>		
1,2,3,4-Cp*Fe(PC <sub>2</sub> B <sub>8</sub> H <sub>10</sub> )	S, H, B, P, E, Mössbauer	[782]
1,2,3,4-Cp*Fe(PC <sub>2</sub> B <sub>8</sub> H <sub>10</sub> ) <sup>+</sup>	ESR	[782]

Continued

Compound	Information	References
1,2,4,5-Cp*Fe(PC <sub>2</sub> B <sub>8</sub> H <sub>10</sub> )	S, H, B, P, E, Mössbauer	[782]
1,2,4,5-Cp*Fe(PC <sub>2</sub> B <sub>8</sub> H <sub>10</sub> ) <sup>+</sup>	ESR	[782]
1,2,4,5-CpFe(PC <sub>2</sub> B <sub>8</sub> H <sub>10</sub> )	S, X, H	[783]
1,2,4,8-CpFe(PC <sub>2</sub> B <sub>8</sub> H <sub>10</sub> )	S, H, B(2d), C, IR	[783]
CpFe(PC <sub>2</sub> B <sub>8</sub> H <sub>10</sub> ) 3 isomers	S, H, B(2d), C, P, IR, MS	[784]
<b>Closo-FeC<sub>3</sub>B<sub>8</sub> clusters</b>		
<i>Mono(tricarbollyl) Fe complexes</i>		
2,1,8,10-CpFe(C <sub>3</sub> B <sub>8</sub> H <sub>11</sub> )	S, H, B, MS	[404]
1,2,3,9-CpFe[(2,3-cyclo-CH <sub>2</sub> OCH <sub>2</sub> )(9-HNCMe <sub>3</sub> )C <sub>3</sub> B <sub>8</sub> H <sub>8</sub> ] room temperature polyhedral rearrangement	S, X, H, B	[786]
1,2,4,10-CpFe( <i>n</i> -RC <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ) <i>n</i> = 2, 10 R = Ph, 1-C <sub>10</sub> H <sub>7</sub> , 2-C <sub>10</sub> H <sub>7</sub>	S, H, B, C	[1569]
1,2,4,10-CpFe[(2-CpFeC <sub>5</sub> H <sub>4</sub> )(10-Me <sub>3</sub> CNH)C <sub>3</sub> B <sub>8</sub> H <sub>9</sub> ]	S, H, B, E	[787]
1,2,4,12-CpFe(12-RC <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ) R = Ph, 1-C <sub>10</sub> H <sub>7</sub> , 2-C <sub>10</sub> H <sub>7</sub>	S, H, B, C	[1569]
1,2,4,12-CpFe[(2-CpFeC <sub>5</sub> H <sub>4</sub> )(12-Me <sub>3</sub> CNH)C <sub>3</sub> B <sub>8</sub> H <sub>9</sub> ]	S, H, B, E	[787]
1,2,4,10-CpFe(RR'-10-R''C <sub>3</sub> B <sub>8</sub> H <sub>8</sub> ) R,R' = H, Me; R'' = NHCMe <sub>3</sub>	S, H, B(2d), E	[1421]
2,1,7,10-CpFe(C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> )- <i>n</i> -NHCMe <sub>3</sub> <i>n</i> = 1, 4, 8	S, X, H, B, MS	[404]
2,1,7,9-CpFe(C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> )-9-NH-CH <sub>2</sub> -C <sub>6</sub> H <sub>4</sub> Me	S, H, B, C, MS	[788]
<i>Nido</i> -CpFe(C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> )-NHCMe <sub>3</sub> <sup>2-</sup>	S, B(2d)	[789]
2,1,7,9-CpFe(RC <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ) R = H <sub>2</sub> N, Me <sub>2</sub> N, Me <sub>3</sub> CHN, Me <sub>3</sub> CMeN, MeHN	S, X(Me <sub>3</sub> CHN), H, B(2d)	[793]
[1',7',9'-(HNC <sub>4</sub> H <sub>9</sub> )C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ](3-Fe)[1,2-LC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] L = SMe <sub>2</sub> , NMe <sub>3</sub>	S, X, H, B	[1442]
1,2,4,12-CpFe[12-(2'-C <sub>10</sub> H <sub>7</sub> )C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ]	S, X, H, B, C	[1481]
1,2,4,12-CpFe(2-PhC <sub>3</sub> B <sub>8</sub> H <sub>10</sub> )	S, X, H, B, C	[1481]
1,2,4,12-(C <sub>6</sub> H <sub>6</sub> )Fe{[12-(Me <sub>3</sub> C)HN]C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> }	X	[1530]
1,2,4,12-CpFe[(12-Me <sub>3</sub> CNH)C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ]	S, H, B, C, MS	[423]
	E, Mössbauer	[718]
<i>Bis(tricarbollyl) Fe complexes</i>		
[9,1,7-(NHCMe <sub>3</sub> )C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ] <sub>2</sub> Fe	S, X, H, B(2d), MS	[792]
(1,7,9-RC <sub>3</sub> B <sub>8</sub> H <sub>10</sub> )Fe(1,7,10-RC <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ) R = NHCMe <sub>3</sub> , NH <sub>2</sub>	S, X(NHCMe <sub>3</sub> ), H, B(2d)	[795]
Related derivatives		[795]
2,1,7,9-CpFe(C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> )-9-N(CH <sub>2</sub> -C <sub>6</sub> H <sub>4</sub> -CH <sub>2</sub> ) <sub>2</sub> N-9-(C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> )FeCp	S, X, H, B, C, MS	[788]
2,1,7,9-CpFe(C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> )-9-NH-CH <sub>2</sub> -C <sub>6</sub> H <sub>4</sub> -CH <sub>2</sub> -NH-9-(C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> )-FeCp	S, X, H, B, C, MS	[788]
<i>Nido-FeC<sub>4</sub>B<sub>7</sub> clusters</i>		
(MeC <sub>6</sub> H <sub>5</sub> )Fe(Et <sub>4</sub> C <sub>4</sub> B <sub>7</sub> H <sub>7</sub> )	S, MS	[59]
CpFe[[8-MeOC(O)]C <sub>4</sub> B <sub>7</sub> H <sub>11</sub> ]	S, X, H, B, C, IR, MS	[797]
<b>Ruthenium</b>		
<i>Closo-RuCB<sub>10</sub> clusters</i>		
2,1-(CO) <sub>3</sub> Ru(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup>	S, B	[799]
2,1-[(CO) <sub>6</sub> Ru <sub>2</sub> (μ-H)](CO) <sub>2</sub> Ru(CB <sub>10</sub> H <sub>11</sub> )	S, H, B, C, IR	[801]

Compound	Information	References
2,1-[(CO) <sub>6</sub> Ru <sub>2</sub> ](CO) <sub>2</sub> Ru(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, C, IR	[801]
2,1-[(CO) <sub>6</sub> Ru <sub>2</sub> (μ-H)(Ph <sub>3</sub> P)Au](CO) <sub>2</sub> Ru(CB <sub>10</sub> H <sub>10</sub> )	S, X, H, B, C, P, IR	[801]
Related derivatives		[802]
2,1-[η <sup>6</sup> -C <sub>6</sub> H <sub>5</sub> PPh <sub>2</sub> -RuCl(PPh <sub>3</sub> )](μ-H) <sub>2</sub> Ru(CB <sub>10</sub> H <sub>8</sub> R) R=H, OMe	S, X(OMe), H, B, P, IR	[803]
Exo-Ru <sub>2</sub> (CO) <sub>4</sub> (Ph <sub>2</sub> PCH <sub>2</sub> PPh <sub>2</sub> )-2,1-(CO) <sub>2</sub> Ru[(Me <sub>2</sub> S)CB <sub>10</sub> H <sub>10</sub> ]	S, H, B, C, P, IR	[804]
Exo-Ru <sub>2</sub> (CO) <sub>4</sub> (Ph <sub>2</sub> PCH <sub>2</sub> PPh <sub>2</sub> )-2,1-(CO) <sub>2</sub> Ru[(MeS)CB <sub>10</sub> H <sub>10</sub> ] <sup>-</sup>	S, X, H, B, C, P, IR	[804]
Exo-Ru <sub>2</sub> (CO) <sub>4</sub> (Ph <sub>2</sub> PCH <sub>2</sub> PPh <sub>2</sub> )-2,1-(CO) <sub>2</sub> Ru[(Ph <sub>3</sub> PAuMeS)CB <sub>10</sub> H <sub>10</sub> ]	S, X, H, B, C, P, IR	[804]
Exo-Ru <sub>2</sub> (CO) <sub>4</sub> (μ-H){(PPh <sub>2</sub> ) <sub>2</sub> [μ-Fe(C <sub>5</sub> H <sub>4</sub> ) <sub>2</sub> ]}-2,1-(CO) <sub>2</sub> Ru[(Me <sub>2</sub> S)-CB <sub>10</sub> H <sub>9</sub> ]	S, H, B, C, P, IR	[800]
Exo-Ru <sub>2</sub> (CO) <sub>4</sub> (μ-H){(PPh <sub>2</sub> ) <sub>2</sub> (μ-1',2'-C <sub>2</sub> B <sub>10</sub> H <sub>10</sub> )}-2,1-(CO) <sub>2</sub> Ru-[(Me <sub>2</sub> S)CB <sub>10</sub> H <sub>9</sub> ]	S, H, B, C, P, IR	[800]
Exo-Ru <sub>2</sub> (CO) <sub>4</sub> (μ-SCMe <sub>3</sub> )-(μ-SCMe <sub>3</sub> )-2,1-(CO) <sub>2</sub> Ru[(Me <sub>2</sub> S)CB <sub>10</sub> H <sub>9</sub> ]	S, X, H, B, C, P, IR	[800]
Related derivatives		[800]
<b>Closo-Ru<sub>3</sub>CB<sub>8</sub> clusters</b>		
Exo-Ru <sub>3</sub> (CO) <sub>9</sub> -2,7,11,1-(CO) <sub>6</sub> Ru <sub>3</sub> (CB <sub>8</sub> H <sub>9</sub> ) <sup>-</sup> planar Ru <sub>6</sub>	S, X, H, B, C, P, IR	[805]
Exo-Ru <sub>2</sub> (CO) <sub>4</sub> L <sub>2</sub> -2,3,4,1-(CO) <sub>6</sub> Ru <sub>3</sub> (PhCB <sub>8</sub> H <sub>8</sub> ) <sup>-</sup> L=CO, PPh <sub>3</sub> planar Ru <sub>5</sub>	S, X(CO), H, B, C, P, IR	[805]
<b>Closo-Ru<sub>2</sub>B<sub>9</sub> clusters</b>		
<i>Mono(dicarbollyl) Ru complexes</i>		
3,1,2-LRu[(Me <sub>2</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] L=Cp, Cp*	S, H, B	[806]
Related derivatives		[807]
3,1,2-Cp* <b>Ru</b> [(ClAuPPh <sub>2</sub> )PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B, P	[808]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh <sub>2</sub> ]HCl <b>Ru</b> (Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	Controlled synthesis of poly(methyl methacrylate) with amines	[809,810]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh <sub>2</sub> ]Cl <b>Ru</b> (Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	Controlled synthesis of poly(methyl methacrylate) with amines	[809,810]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh-μ-C <sub>6</sub> H <sub>4</sub> -]Cl <b>Ru</b> (Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -10-)	Controlled synthesis of poly(methyl methacrylate) with amines	[809,810]
2,1,7-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> ) <b>Ru</b> (C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S	[812]
3,1,2-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> ) <b>Ru</b> (C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -10-R) <sup>-</sup> R=SMe <sub>2</sub> , SH	S, X(SH), H, B, C	[813]
3,1,2-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> ) <b>Ru</b> (C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -10-SMe)	S, X, H, B, C	[813]
3,1,2-(η <sup>6</sup> -L) <b>Ru</b> (C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6-SMe) <sup>+</sup> L=C <sub>6</sub> H <sub>6</sub> , 1,3,5-C <sub>6</sub> Me <sub>3</sub> H <sub>3</sub>	S, X(C <sub>6</sub> H <sub>6</sub> ), H, B	[818]
3,1,2-(η <sup>6</sup> -arene) <b>Ru</b> (RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -8-L) <sup>+</sup> R=H, Me; arene=C <sub>6</sub> H <sub>6</sub> , MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> ; L=SMe <sub>2</sub> , SC <sub>4</sub> Me <sub>4</sub> ,SEtPh	S, X(C <sub>6</sub> H <sub>6</sub> , H, SMe <sub>2</sub> ), H, B, C	[819]
3,1,2-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> ) <b>Ru</b> (RC <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -8-SMe <sub>2</sub> ) <sup>+</sup>	S, H, B, C	[819]
3,1,2-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> ) <b>Ru</b> (RC <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -8-SMe)	S, H, B, C	[819]
3,1,2-(η <sup>6</sup> -MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> ) <b>Ru</b> [(ClAuPPh <sub>2</sub> )PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, P	[808]
3,1,2-(η <sup>6</sup> -MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> ) <b>Ru</b> [(PhS) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, IR	[820]
3,1,2-(η <sup>6</sup> -MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> ) <b>Ru</b> [(PhC≡C)PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, IR	[821]
3,1,2-(η <sup>6</sup> -MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> ) <b>Ru</b> [CpFe(C <sub>5</sub> H <sub>4</sub> )-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B	[822]
3,1,2-(η <sup>6</sup> -MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> ) <b>Ru</b> [(MeOCH <sub>2</sub> )RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] R=(CH <sub>2</sub> OMe, H	S, X(CH <sub>2</sub> OMe), H, B, IR	[545]
3,1,2-(η <sup>6</sup> -MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> ) <b>Ru</b> (RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R, R'=H, H; H, Ph; Me, Ph	S, X(Me, Ph), H, B, IR	[823]

Continued

Compound	Information	References
2,1,8-(MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> )Ru[8-(1'-1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, MS	[1570]
2,1,12-(η <sup>6</sup> -MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> )Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B	[1419]
3,1,2-(η <sup>6</sup> -1',3',5'-C <sub>6</sub> H <sub>3</sub> Me <sub>3</sub> )Ru(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, X, H, B, IR	[823]
2,1,8-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Ru[(8-SMe <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -11-SMe <sub>2</sub> ] <sup>+</sup> BF <sub>4</sub> <sup>-</sup>	S, X, B	[824]
3,1,2-(CO) <sub>3</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )-0.5C <sub>6</sub> H <sub>6</sub>	S, H, B, IR, MS, UV	[827]
3,1,2-(CO) <sub>3</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, P, IR	[828]
3,1,2-(CO) <sub>2</sub> (CO <sub>2</sub> Me)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, IR	[826]
Related derivatives		[826]
3,1,2-(CO) <sub>2</sub> Ru[(MeOCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	S, H, B, IR	[545]
3,1,2-(CO) <sub>2</sub> (THF)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR	[825]
3,1,2-(CO) <sub>2</sub> (MeC=CHMe)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, C, IR	[829]
3,1,2-(CO) <sub>2</sub> (MeC=CMePPh <sub>3</sub> )Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, IR	[829]
3,1,2-(CO) <sub>2</sub> Ru[C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -7-CH(PMe <sub>2</sub> Ph)-CH <sub>2</sub> -] Ru-CH <sub>2</sub>	S, X, H, B, C, IR	[829]
Related derivatives		[829]
μ-Ru <sub>2</sub> (CO) <sub>6</sub> -3,1,2-(CO) <sub>2</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -μ-H <sub>2</sub> )	S, H, B, C, P, IR	[828]
μ-Ru <sub>2</sub> (CO) <sub>5</sub> (PR <sub>3</sub> )H-3,1,2-(CO) <sub>2</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -μ-H) R=PPh <sub>3</sub> , P(MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> ) <sub>3</sub>	S, H, B, C, P, IR	[828]
μ-Ru <sub>2</sub> (CO) <sub>4</sub> (PR <sub>3</sub> ) <sub>2</sub> H]-3,1,2-(CO) <sub>2</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -μ-H) R=PPh <sub>3</sub> , PMe <sub>2</sub> Ph	S, X(PMe <sub>2</sub> Ph), H, B, C, P, IR	[828]
Related derivatives		[828]
3,1,2-(CO) <sub>2</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[830]
3,1,2-(CO) <sub>2</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -8-CH=CHSiMe <sub>3</sub> )	S, H, B, C, IR	[830]
3,1,2-(MeC≡CPh)(CO) <sub>2</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, P, IR	[831]
3,1,2-L(CO) <sub>2</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=AsPh <sub>3</sub> , SbPh <sub>3</sub> , S=P(Ph <sub>2</sub> )CH <sub>2</sub> P-(Ph <sub>2</sub> )=S	S, H, B, C, P, IR	[831]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> H <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, IR	[832,833]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HXRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) X=Cl, CO	S, IR	[832]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRu(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-7-NC <sub>5</sub> H <sub>5</sub>	S, H, IR	[832,835]
	B	[832]
	P	[835]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRu(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-PPh <sub>2</sub> R R=Ph, Me	S, X(Ph), H, B, C, P, IR	[836]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRu(RR'C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -8-SR''R''') R, R'=H, Me; R'', R'''=Me, Et, Ph, (CH <sub>2</sub> ) <sub>4</sub> catalysts for radical polymerization of styrene and <i>n</i> -butyl acrylate	S	[1366]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> (CO)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, IR	[827]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, X, H, B, P, IR	[838]
3,1,2-(CO)L(Ph <sub>3</sub> P)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=CO, PPh <sub>3</sub>	S, H, P, IR	[838]
(Ph <sub>3</sub> P)(CO)Rh(μ-H)-3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, P, IR, MS	[838]
(η <sup>4</sup> -L)Rh(μ-H)-3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=C <sub>8</sub> H <sub>12</sub> , C <sub>7</sub> H <sub>8</sub>	S, X(C <sub>8</sub> H <sub>12</sub> ), H, B, C, P, IR, MS	[838]
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	H, B	[832,833]
	IR	[832]

Compound	Information	References
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> H <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	B, IR	[832,833]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> (NO <sub>3</sub> )Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )·2.2CH <sub>2</sub> Cl <sub>2</sub>	S, X, H, B, P	[1581]
3,1,2-ClH(cyclo-Ph <sub>2</sub> PCHMeCH <sub>2</sub> CHMePPh <sub>2</sub> )Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P, IR	[839]
3,1,2-Cl(cyclo-Ph <sub>2</sub> PCHMeCH <sub>2</sub> CHMePPh <sub>2</sub> )Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, IR	[839]
3,1,2-[cyclo-Ph(C <sub>6</sub> H <sub>4</sub> )(PCHMeCH <sub>2</sub> CHMePPh <sub>2</sub> )-Cl]Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) B-C <sub>6</sub> H <sub>4</sub>	S, X, IR	[839]
3,1,2-[Ph(C <sub>6</sub> H <sub>4</sub> )P(CH <sub>2</sub> ) <sub>3</sub> PPh <sub>2</sub> ]ClRu(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -8-) B-C <sub>6</sub> H <sub>4</sub> -P catalyst for methyl methacrylate polymerization	S, X, ESR	[840]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> H(Cl)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) catalyst for methyl methacrylate polymerization	S	[840]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh <sub>2</sub> ]H(Cl)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) catalyst for methyl methacrylate polymerization	S	[840]
3,1,2-Cl(H)(Ph <sub>2</sub> PCHMeCH <sub>2</sub> CHMe)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P, IR	[841]
3,1,2-(Ph <sub>3</sub> P)(C <sub>5</sub> H <sub>3</sub> N) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S (phosphine displacement)	[1411]
3,1,2-(Ph <sub>3</sub> P)(2-phen)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S (phosphine displacement)	[1411]
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> (CO)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, P	[835]
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> HRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R=Cl, CO	S, IR	[832]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>n</sub> PPh <sub>2</sub> ](H)ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) n=3, 4	S, H, B, C, P, IR	[842]
3,1,2-(CO)(Ph <sub>3</sub> P)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-CH=CHR) R=Ph, H	S, X(Ph), H, B, C, P, IR	[843]
3,1,2-L(CO)(Ph <sub>3</sub> P)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=CNCMe <sub>3</sub> , THF	S, H, B, C, P, IR	[843]
3,1,2-(CO)(Ph <sub>3</sub> P)(CNCMe <sub>3</sub> )Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-CH=CH <sub>2</sub> )	S, X, H, B, C, P, IR	[843]
Related derivatives		[843]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>n</sub> PPh <sub>2</sub> ]ClHRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) n=3-5	S, H, B, P	[1579]
3,1,2-[PhP(CH <sub>2</sub> ) <sub>n</sub> PPh <sub>2</sub> ]BrRu(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -10-C <sub>6</sub> H <sub>4</sub> ) P-C <sub>6</sub> H <sub>4</sub> n=3-5	S, X (n=4), H, B, P	[1579]
3,1,2-[P(CH <sub>2</sub> ) <sub>n</sub> PPh <sub>2</sub> ]BrRu[C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -9,10-(C <sub>6</sub> H <sub>4</sub> ) <sub>2</sub> 2 P-C <sub>6</sub> H <sub>4</sub> n=3,4	S, X(n=3), H, B, P	[1579]
3,1,2-LH(Ph <sub>3</sub> P)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> L=Ph <sub>3</sub> P, CO	S, H, B, C, P	[844]
3,1,2-[S <sub>3</sub> (C <sub>2</sub> H <sub>4</sub> ) <sub>3</sub> ]Ru(PhRC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=H, Ph	S, X, H, B, C	[845]
Exo-[(CO)(Ph <sub>3</sub> P)Rh(μ-H)]-3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, P	[844]
Exo-[(Ph <sub>3</sub> P) <sub>2</sub> RuH(μ-H)]-3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, P	[844]
Exo-[(CO) <sub>3</sub> (Ph <sub>3</sub> P)Ru(μ-H)]-3,1,2-(Ph <sub>3</sub> P)(CO)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, X, H, B, C, P	[844]
Exo-[(Ph <sub>3</sub> P)Cu(μ-H)]-3,1,2-(Ph <sub>3</sub> P)LRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=Ph <sub>3</sub> P, CO	S, X, H, B, C, P	[844]
Exo-[(Ph <sub>3</sub> P)Au(μ-H)]-3,1,2-(Ph <sub>3</sub> P)LRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=Ph <sub>3</sub> P, CO	S, X, H, B, C, P	[844]
Related derivatives		[844]
Exo-(Ph <sub>3</sub> P)Cu(μ-H)-3,1,2-[(Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh <sub>2</sub> )Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B	[1429]
Exo-Cp(CO) <sub>2</sub> W=C(C <sub>6</sub> H <sub>4</sub> Me)-3,1,2-(CO) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	X	[846]
3,1,2-L(N <sub>2</sub> C <sub>10</sub> H <sub>6</sub> R <sub>2</sub> )Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=CO, NCMe; R=H, (CH <sub>2</sub> ) <sub>8</sub> Me, CMe <sub>3</sub> N <sub>2</sub> C <sub>10</sub> H <sub>6</sub> =bipyridine	S, X(CO, H), H, B, C, IR, UV, ESR(CO, Me), E	[847]
Exo-HB(pz) <sub>3</sub> (CO) <sub>2</sub> M=C(C <sub>6</sub> H <sub>4</sub> Me)-3,1,2-(CO) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) M=Mo, W; pz=pyrazolyl	S, X(W), H, B, C, IR	[846]
3,1,2-Cl(Ph <sub>3</sub> P) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-10-(9-Hg-1,2-C <sub>2</sub> B <sub>10</sub> H <sub>11</sub> )	S, X, H, B, P	[848]
3,1,2-(CO)LRu[(Me <sub>2</sub> NCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] N→Ru L=CO, MeCN	S, X, H, B, C, IR	[752]

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Compound	Information	References
3,1,2-[Cp(CO) <sub>2</sub> M=CC <sub>6</sub> H <sub>4</sub> Me](CO) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) M=W, Mo	S, H, B, C, IR	[849]
3,1,2-[Cp(O)W-CHC <sub>6</sub> H <sub>4</sub> Me](CO) <sub>2</sub> Ru(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) R=H, Me; two isomers each	S, X(H), H, B, C, IR	[849]
Related derivatives		[849]
<i>Exo</i> -(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Ir(μ-H)-3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, P, IR	[850]
3,1,2-(η <sup>6</sup> -C <sub>5</sub> H <sub>5</sub> BMe)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) boratabenzene complex	S, X	[1425]
3,1,2-(C <sub>6</sub> H <sub>6</sub> )Ru(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>+</sup> R=H, Me C <sub>6</sub> H <sub>6</sub> exchange with arenes		[1437]
3,1,2-H[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh <sub>2</sub> ]Cl(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	Thermal rearrangement	[1451]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh <sub>2</sub> ]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) catalyst for radical polymerization of methyl methacrylate	S, X, H, ESR	[1451]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh(C <sub>6</sub> H <sub>4</sub> )]Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) C <sub>6</sub> H <sub>4</sub> -B(8) catalyst for radical polymerization of methyl methacrylate	S, X, H, ESR	[1451]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> P(C <sub>6</sub> H <sub>4</sub> ) <sub>2</sub> ]Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) 2 C <sub>6</sub> H <sub>4</sub> -B catalyst for radical polymerization of methyl methacrylate	S, X, H, IR, ESR	[1451]
3,1,2-Cl[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> P(C <sub>6</sub> H <sub>4</sub> ) <sub>2</sub> ]Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) C <sub>6</sub> H <sub>4</sub> -B catalyst for controlled radical polymerization of Me methacrylate to poly(methyl methacrylate) (PMMA)	S(as in ref. [1477])	[1523]
3,1,2-(H)[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>5</sub> PPh <sub>2</sub> ]ClRu(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P	[1477]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>5</sub> PPh <sub>2</sub> ]ClRu(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P, ESR	[1477]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>5</sub> PPh(C <sub>6</sub> H <sub>4</sub> )]ClRu(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -R) R=H, Cl	S, H, B, P, ESR	[1477]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>5</sub> P(C <sub>6</sub> H <sub>4</sub> ) <sub>2</sub> ]ClRu(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) 2 Ph-B	S, H, B, P, ESR	[1477]
3,1 2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>n</sub> PPh <sub>2</sub> ]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) n=2-5	E(catalysis in methylmethacrylate polymerization)	[1515]
3,1 2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>n</sub> PPh(8-C <sub>6</sub> H <sub>4</sub> )]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) n=3-5	E(catalysis in methylmethacrylate polymerization)	[1515]
3,1 2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>n</sub> P(7-C <sub>6</sub> H <sub>4</sub> )(8-C <sub>6</sub> H <sub>4</sub> )]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) n=4,5	E(catalysis in methylmethacrylate polymerization)	[1515]
3,1 2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>n</sub> PPh(8-C <sub>6</sub> H <sub>4</sub> )]ClRu(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) n=3,4	E(catalysis in methylmethacrylate polymerization)	[1515]
3,1 2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh <sub>2</sub> ]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	Matrix-assisted laser desorption/ionization TOF MS	[1516]
3,1 2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh(8-C <sub>6</sub> H <sub>4</sub> )]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	Matrix-assisted laser desorption/ionization TOF MS	[1516]
3,1 2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> P(7-C <sub>6</sub> H <sub>4</sub> )(8-C <sub>6</sub> H <sub>4</sub> )]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	Matrix-assisted laser desorption/ionization TOF MS	[1516]
3,1 2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh(8-C <sub>6</sub> H <sub>4</sub> )]ClRu(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )	Matrix-assisted laser desorption/ionization TOF MS	[1516]
3,1 2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>4</sub> PPh(8-C <sub>6</sub> H <sub>4</sub> )]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -7-Cl)	Matrix-assisted laser desorption/ionization TOF MS	[1516]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>5</sub> PPh <sub>2</sub> ]ClRu(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S(as in N74), catalyst for controlled radical polymerization of Me methacrylate to poly(methyl methacrylate) (PMMA)	[1523]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>5</sub> PPh(C <sub>6</sub> H <sub>4</sub> )]ClRu(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -R) R=H, Cl	S(as in N74), catalyst for controlled radical polymerization of Me methacrylate to poly(methyl methacrylate) (PMMA)	[1523]

Compound	Information	References
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>5</sub> P(C <sub>6</sub> H <sub>4</sub> ) <sub>2</sub> ]ClRu(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) 2 Ph—B	S(as in ref. 1477), catalyst for controlled radical polymerization of Me methacrylate to poly(methyl methacrylate) (PMMA)	[1523]
3,1,2-(O <sub>2</sub> )[Ph <sub>2</sub> Ph(CH <sub>2</sub> ) <sub>n</sub> PPh-C <sub>6</sub> H <sub>4</sub> ]ClRu(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) C <sub>6</sub> H <sub>4</sub> —B(8) <i>n</i> =3, 4	S, X, H(2d), P, C, IR	[1539]
3,1,2-(C <sub>6</sub> H <sub>6</sub> )Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )- <i>n</i> -SMe <sub>2</sub> <sup>+</sup> <i>n</i> =4, 8	S, H, B	[1592]
3,1,2-(C <sub>6</sub> H <sub>6</sub> )Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )- <i>n</i> -SMe <i>n</i> =4, 8	S, H, B	[1592]
<b>Bis(dicarbollyl) Ru complexes</b>		
Tl <sup>+</sup> [3,1,2-(CO) <sub>2</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )] <sub>2</sub> <sup>-</sup>	S, X, H, B, C, IR	[830]
[3,1,2-(CO) <sub>2</sub> Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> (Ph <sub>2</sub> P) <sub>2</sub> CH <sub>2</sub>	S, X, H, B, C, P, IR	[831]
<i>trans</i> -[(1,2-MeC <sub>6</sub> H <sub>4</sub> CHMe <sub>2</sub> ) <sub>3</sub> -Ru(1,2-PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )] <sub>2</sub> CH=CH	S, X, H, B	[851]
[3,1,2-(CO)(μ-CO)Ru(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> <sup>2-</sup>	S, H, B, IR	[826]
<b>Closo-RuC<sub>3</sub>B<sub>8</sub> clusters</b>		
[9,1,7-(NHCMe <sub>3</sub> )C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ] <sub>2</sub> Ru	S, X, H, B(2d), MS	[792]
1,2,4,12-Cp* <sup>+</sup> Ru[(12-Me <sub>3</sub> CNH)C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ]	S, X, H, B, C, MS	[423]
1,2,4,10-Cp* <sup>+</sup> Ru[(12-Me <sub>3</sub> CNH)C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ]	S	[423]
1,2,4,12-(C <sub>6</sub> H <sub>3</sub> Me <sub>3</sub> )Ru[(12-Me <sub>3</sub> CNH)C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ]	S, H, B	[693]
<b>Osmium</b>		
<b>Closo-OsCB<sub>10</sub> clusters</b>		
2,1-(CO) <sub>3</sub> Os(CB <sub>10</sub> H <sub>10</sub> )-5-NMe <sub>3</sub>	S, X, H, B, C, UV	[799]
<i>Exo</i> -Os <sub>2</sub> (CO) <sub>6</sub> -2,1-(CO) <sub>2</sub> Os(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, C, UV	[799]
<i>Exo</i> -Os <sub>2</sub> (CO) <sub>6</sub> (μ-H)-2,1-(CO) <sub>2</sub> Os(CB <sub>10</sub> H <sub>11</sub> )	S, H, B, C, UV	[799]
μ-[Os(CO) <sub>3</sub> ] <sub>2</sub> -2,1-(CO) <sub>2</sub> Os(CB <sub>10</sub> H <sub>10</sub> )-7-NMe <sub>3</sub>	S, X, H, B, IR	[852]
<b>Closo-OsC<sub>2</sub>B<sub>9</sub> clusters</b>		
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> (H) <sub>2</sub> Os(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X(Cl), H, B, P, IR	[853]
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> (H) <sub>2</sub> Os(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, P, IR	[854]
<b>Cobalt</b>		
<b>Closo-CoCB<sub>10</sub> clusters</b>		
2,1-Co(CB <sub>10</sub> H <sub>11</sub> ) <sub>2</sub> <sup>2-</sup>	S, H, IR, UV	[593]
2,1-Co[(H <sub>3</sub> N)CB <sub>10</sub> H <sub>10</sub> ] <sub>2</sub> <sup>-</sup>	S, UV	[593]
2,1-Co(CB <sub>10</sub> H <sub>10</sub> —NH <sub>2</sub> Et) <sub>2</sub> <sup>-</sup>	S	[682]
2,1-CpCo{[(Me <sub>3</sub> Si) <sub>2</sub> CH]CB <sub>10</sub> H <sub>9</sub> }- <i>n</i> -SMe <sub>2</sub> <i>n</i> =7,12	S, H, B, IR, MS	[860]
2,1-CpCo{[(Me <sub>3</sub> Si) <sub>2</sub> CH]CB <sub>10</sub> H <sub>9</sub> }-C <sub>5</sub> H <sub>4</sub> CoCp	S, H, B, MS	[860]
2,1-(MeC <sub>6</sub> H <sub>5</sub> )Co{[(Me <sub>3</sub> Si) <sub>2</sub> CH]CB <sub>10</sub> H <sub>10</sub> }	S, X, H, B, IR, MS	[860]
2,1-(MeC <sub>6</sub> H <sub>5</sub> )Co[(Me <sub>3</sub> CNH <sub>2</sub> )CB <sub>10</sub> H <sub>10</sub> ]	S, X, H, B, C, P	[861]
2,1-L(ON)Co[(Me <sub>3</sub> N)CB <sub>10</sub> H <sub>10</sub> ] L=CO, PPh <sub>3</sub> , PEt <sub>3</sub> , CNCMe <sub>3</sub>	S, X(CNCMe <sub>3</sub> ), H, B, C, IR	[862]
2,1-(OC)(ON)Co[(Me <sub>3</sub> N)CB <sub>10</sub> H <sub>10</sub> ]-2,7-Co(CO(NO))	S, X, H, B, C, IR	[862]
2,1-(Ph <sub>3</sub> PCu)(ON)LCo(HCB <sub>10</sub> H <sub>10</sub> ) (Co-Cu) L=CO, PPh <sub>3</sub>	S, X(PPh <sub>3</sub> ), H, B, C, IR	[863]
2,1-(Ph <sub>3</sub> PAG)(Et <sub>3</sub> P)(ON)Co(HCB <sub>10</sub> H <sub>9</sub> ) (Ag—H—B)	S, H, B, C, IR	[863]

Continued



Compound	Information	References
Related derivatives		[863]
<i>Closo-Co<sub>2</sub>CB<sub>9</sub> clusters</i>		
2,3,1-(CO) <sub>5</sub> Co <sub>2</sub> (CB <sub>9</sub> H <sub>10</sub> ) <sup>-</sup>	S, X, H, B, C, P, IR	[437]
2,3,1-(CO) <sub>5</sub> Co <sub>2</sub> (CB <sub>9</sub> H <sub>9</sub> -12-L) L = O(CH <sub>2</sub> ) <sub>4</sub> , O(CH <sub>2</sub> ) <sub>4</sub> PPh <sub>3</sub>	S, X[O(CH <sub>2</sub> ) <sub>4</sub> ], H, B, C, P, IR	[437]
<i>Closo-CoPCB<sub>9</sub> clusters</i>		
3,1,2-/2,1,7-Co(PCB <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> <sup>-</sup>	S, H, IR, UV	[684]
2,1,7-(MePCB <sub>9</sub> H <sub>10</sub> )Co(PCB <sub>9</sub> H <sub>10</sub> )	S, H, IR, MS	[684]
3,1,2-(η-C <sub>9</sub> H <sub>13</sub> )Co(PCB <sub>9</sub> H <sub>10</sub> )	S, X, H, B, P, IR, MS	[1493]
<i>Closo-CoC<sub>2</sub>B<sub>9</sub> clusters</i>		
<i>Mono(dicarbollyl) Co complexes</i>		
<i>CpCo(C<sub>2</sub>B<sub>9</sub>) and Cp*Co(C<sub>2</sub>B<sub>9</sub>) complexes</i>		
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>11-n</sub> D <sub>n</sub> )	S(D exchange), H, IR	[874]
2,1,12-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, B, IR	[875]
2,1,7 (4,1,7)-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S	[812]
3,1,2-CpCo(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R = H, Me, Ph, HOCH <sub>2</sub> , ClCH <sub>2</sub> , HO(O)C	MS (detailed)	[878]
8,9'-[3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )] <sub>2</sub>	S, X, H, B	[879]
3,1,2-CpCo(MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	H	[880]
3,1,2-CpCo[(CH <sub>2</sub> ) <sub>3</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B, IR, UV, E	[869]
3,1,2-CpCo(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	C	[870]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-Ph)	S	[881]
3,1,2-CpCo{[C(O)OH] <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> }	S, H, IR	[882]
3,1,2-CpCo(RMeC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R = C(O)OMe, CF=CFCF <sub>3</sub>	S, H, IR	[316]
3,1,2-CpCo{[CH(OH)R]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> } R = Me, Et, CMe <sub>3</sub>	S	[883]
3,1,2-CpCo{[(CH <sub>2</sub> R)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] R = Me, Et, CMe <sub>3</sub> }	S	[883]
3,1,2/2,1,7-CpCo{[(HOCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S	[884]
3,1,2-CpCo[( <i>m/p</i> -FC <sub>6</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	F	[761]
3,1,2-CpCo{[C(O)NMe <sub>2</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> }	S, H(rotational barrier)	[885]
3,1,2-CpCo{[(CH <sub>2</sub> NMe <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, H	[885]
3,1,2-CpCo{[ <i>p</i> -NO <sub>2</sub> C <sub>6</sub> H <sub>4</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]-R R = H, ONO <sub>2</sub> }	S	[886]
3,1,2-CpCo{[ <i>p</i> -NH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]-R R = H, NH <sub>2</sub> }	S	[886]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-6-CH=CH <sub>2</sub>	S	[887]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -R) R = OH, ONO <sub>2</sub>	S	[886]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) R = OH, OC(O)CF <sub>3</sub>	S	[724]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6-NH <sub>2</sub> )	S	[887]
(+/-)3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe)	S, H, B, MS, CD	[810]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -8,9-Br <sub>2</sub> ) supramolecular C—H...X—B bonds	S, X, H, B, C	[891]
3,1,2-CpCo(MeC <sub>2</sub> B <sub>9</sub> H <sub>9-n</sub> Br <sub>n</sub> ) n = 1,2	S, H	[880]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-8,9,12-Br <sub>3</sub>	E	

Compound	Information	References
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-I)	S	[881]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-9-HgCl	S, H, B	[892]
3,1,2-CpCo(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> )-9-Y-12-HgX R=H, CH <sub>2</sub> OH; X=Cl, Br, OC(O)CF <sub>3</sub> ; Y=H, HgCl, HgBr	E	[873]
3,1,2-CpCo(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> )-9,12-X <sub>2</sub> X=HgCl, HgOAc; R=H, CH <sub>2</sub> OH	S, H, B	[892]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-n-HgOC(O)Me n=8,9	S, H, B	[892]
	S, X	[893]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-9,12-[HgOC(O)Me] <sub>2</sub> n=8,9	S	[893]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-n-HgOC(O)CF <sub>3</sub> n=8,9	S, H, B	[892]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-9-Hg(10- <i>nido</i> -7,8-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B	[894]
3,1,2-(PhC <sub>5</sub> H <sub>4</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, IR	[896]
3,1,2-(RC <sub>5</sub> H <sub>4</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R= <i>n</i> -C <sub>4</sub> H <sub>9</sub> , C≡CPh	S, H, IR	[882]
3,1,2-[ <i>cyclo</i> -1,2- <i>o</i> -NC <sub>5</sub> H <sub>4</sub> -C(OH)] [ <i>o</i> -C <sub>5</sub> H <sub>4</sub> N-C(O)O]Co(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) 2Co-N, Co-O	S, X	[1472]
3,1,2-CpCo[1-(1'-1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, MS	[1570]
2,1,8-CpCo[8-(1'-1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, MS	[1570]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>n</sub> PPh <sub>2</sub> ]ClCo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) n=2,3	S, X(n=2), H, B, P	[1534]
8,1,2-CpCo[(F <sub>3</sub> C-C <sub>6</sub> F <sub>4</sub> )PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, MS	[1549]
8,1,2-CpCo[(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, MS	[1549]
8,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, MS	[1549]
<b>(<math>\eta^6</math>-benzene)Co(C<sub>2</sub>B<sub>9</sub>) complexes</b>		
3,1,2-( <i>p</i> -NO <sub>2</sub> C <sub>6</sub> H <sub>4</sub> -C <sub>5</sub> H <sub>4</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> R) R=H, ONO <sub>2</sub>	S	[886]
3,1,2-( <i>m/p</i> -FC <sub>6</sub> H <sub>4</sub> -C <sub>5</sub> H <sub>4</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, F	[896]
<b>Other (hydrocarbon)Co(C<sub>2</sub>B<sub>9</sub>) complexes</b>		
( $\eta^4$ -C <sub>4</sub> Me <sub>4</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )	S, H, B	[739]
3,1,2-( $\eta^1, \eta^3$ -C <sub>8</sub> H <sub>12</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -5-SMe <sub>2</sub> )	S, X, H, B	[1426]
3,1,2-( $\eta^5$ -C <sub>9</sub> H <sub>7</sub> )Co(RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) RR'=Ph, H; CH <sub>2</sub> OMe, H; CH <sub>2</sub> OMe, CH <sub>2</sub> OMe C <sub>9</sub> H <sub>7</sub> =indenyl	S, X, H, B, IR	[899]
<b>(<math>\eta^6</math>-pyrrolyl)Co(C<sub>2</sub>B<sub>9</sub>) complexes</b>		
3,1,11-(NC <sub>4</sub> H <sub>4</sub> )Co(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, H, B	[902]
3,1,2-(NC <sub>4</sub> H <sub>4</sub> )Co(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, X, H, B, C, IR	[903]
3,1,2-(NC <sub>4</sub> H <sub>4</sub> )Co[(CH <sub>2</sub> ) <sub>3</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H(variable temp), B, C	[904]
3,1,2-(NC <sub>4</sub> H <sub>4</sub> )Co[Me(MeS)C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, C, IR, E	[901]
3,1,2-(NC <sub>4</sub> H <sub>4</sub> )Co[(MeS) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, X, H, B, C, IR	[903]
3,1,2-(NC <sub>4</sub> Me <sub>2</sub> H <sub>2</sub> )Co(RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R,R'=Ph; R=Me, R'=SMe	S, H, B, C, IR, E(influence of substituents)	[901]
3,1,2-(NC <sub>4</sub> Me <sub>2</sub> H <sub>2</sub> )Co(MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, X, H, B, C	[906]
3,1,2-(NC <sub>4</sub> H <sub>4</sub> )Co[R[(CH <sub>2</sub> ) <sub>3</sub> NC <sub>4</sub> H <sub>4</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] R=Me, Ph	S, X(Me), H, B, C, IR	[909]
3,1,2-(NC <sub>4</sub> H <sub>4</sub> )Co[(C <sub>3</sub> H <sub>5</sub> )PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B, C, IR	[905]
Ag[3,1,2-(NC <sub>4</sub> H <sub>4</sub> )Co(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )] <sub>2</sub> <sup>+</sup> R=H, Ph	S, H(actual), B(actual), C, IR, MS, E	[910]

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Compound	Information	References
3,1,2-[(Me-aquocobaloxime)NC <sub>4</sub> H <sub>4</sub> )Co(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=H, Ph	S, H(actual), B(actual), C, IR, MS, E	[910]
<i>Other (ligand)Co(C<sub>2</sub>B<sub>9</sub>) complexes</i>		
3,1,2-(CO) <sub>2</sub> Co(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -10-CO)	S, H, B, C, IR	[911]
3,1,2-(PMe <sub>2</sub> Ph) <sub>2</sub> ClCo(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, IR	[911]
3,1,2-(PhCN <sub>4</sub> SCoCpS <sub>2</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-3,4-(PhCN <sub>4</sub> S)	S, X, H, B, C, IR	[913]
<i>Bis(dicarbollyl) Co complexes</i>		
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup>	COND, aggregation in H <sub>2</sub> O; dynamic and static light-scattering; atomic force microscopy; scanning electron microscopy	[932]
	Compartmentalized hybrid star-shaped metallacarborane-polyethylene oxide—poly (2-alkyl oxazoline nanoparticles)	[1547]
	Surface activity and molecular reorganization at water surface via B-H...H-C bonds NLO	[1559]
	Preferential chlorination rates, MS	[1504]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> ·HIV-1 protease complex HIV-1 protease inhibitor	X	[933]
(TMTSF) <sub>2</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> TMTSF=tetramethyltetraselenafulvalenium	S, X, COND	[935]
	Raman	[1417]
[3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> ] <sub>8</sub> Zn <sup>2+</sup> (phthalocynaine) intercellular accumulation in GL6 glioblastoma cells		[1553]
(ETTF) <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> ETTF=bis (ethylenedithiotetrathiafulvalenium)	S, X, COND	[935]
(BEDT-TTF) <sub>2</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> R=H, Br BEDT-TTF=bis (ethylenedilithio)tetrathiofulvalene	S, X, COND	[936]
(BEDT-TTF) <sub>2</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I) <sup>-</sup> BEDT-TTF=bis (ethylenedilithio)tetrathiofulvalene	S, X, COND	[936]
(BEDT-TTF) <sub>2</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-Cl) <sub>2</sub> <sup>-</sup> BEDT-TTF=bis (ethylenedilithio)tetrathiofulvalene	S, X, COND	[1568]
(BEDT-TTF) <sub>n</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-Br) <sub>2</sub> <sup>-</sup> n=1,2 BEDT-TTF=bis (ethylenedilithio)tetrathiofulvalene) molecular semiconductors	S, X, COND	[937]
(BMDT-TTF) <sub>4</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-Br) <sub>2</sub> <sup>-</sup> BMDT-TTF=bis (methylenedilithio)tetrathiofulvalene molecular semiconductors	S, X, COND	[937]
(BMDT-TTF) <sub>2</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-Cl) <sub>2</sub> <sup>-</sup> BMDT-TTF=bis (methylenedilithio)tetrathiofulvalene	S, X, COND	[1568]
(BMDT-TTF) <sub>4</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I) <sub>2</sub> <sup>-</sup> BMDT-TTF=bis (methylenedilithio)tetrathiofulvalene molecular semiconductor	S, X, COND	[938]
(BEDT-TTF) <sub>n</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I) <sub>2</sub> <sup>-</sup> n=1,2 BEDT-TTF=bis (ethylenedilithio)tetrathiofulvalene) molecular semiconductor	S, X, COND	[938]
BEDT-TTF <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-I) <sub>2</sub> <sup>-</sup>	S, X, COND	[1428]
TTF <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -9,12-I) <sub>2</sub> <sup>-</sup> radical cation salt	S, X, COND	[1428]
TMTTF <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OH) <sup>-</sup>	S, X, COND	[1427]
BEDT-TTF <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OH) <sup>-</sup>	S, X, COND	[1427]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>6</sub> D <sub>5</sub> ) <sub>2</sub> <sup>-</sup>	S, B	[923]
	S, IR	[939]

Compound	Information	References
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-[(O-CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> <sup>-</sup> NH <sub>2</sub> <sup>+</sup> (CH <sub>2</sub> ) <sub>3</sub> SiO <sub>3</sub> <sup>-</sup> silica zwitterionic HPLC stationary phase	S	[1566]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-[O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> -O-(phthalocyanine-OR)Zn R=Me, [(CH <sub>2</sub> ) <sub>2</sub> O] <sub>2</sub> (C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, MS, UV	[940]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> -O-NC <sub>5</sub> H <sub>4</sub> - <i>m</i> -OR <sup>-</sup> K <sup>+</sup> R=Zn phthalocyanine	S, H, C, MS, UV, fluorescence	[941]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> -1-R <sup>-</sup> R=1,2-C <sub>2</sub> B <sub>10</sub> H <sub>11</sub> , 1,7-C <sub>2</sub> B <sub>10</sub> H <sub>11</sub> , 1,12-C <sub>2</sub> B <sub>10</sub> H <sub>11</sub>	S, H, B, C, MS	[942,943]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> - <i>nido</i> -7,8-RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> <sup>2-</sup> R=H, Me, Ph	S, H, B, C, MS	[942]
<i>nido</i> -7,8-[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> <sup>3-</sup>	S, H, B, C, MS	[942]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> - <i>nido</i> -7,9-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> <sup>2-</sup>	S, H, B, C, MS	[942]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-l)Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-(OCH <sub>2</sub> CH <sub>2</sub> ) <sub>2</sub> X <sup>-</sup> X=Br, I	S, H, B, IR, MS	[944]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OCH <sub>2</sub> CH <sub>2</sub> OH	S, H, B, IR, MS	[944]
3K <sup>+</sup> 2,4,6-N <sub>3</sub> C <sub>3</sub> [ <i>p</i> -C <sub>6</sub> H <sub>4</sub> -O(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> -O-(8-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co-(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>3</sub> ] <sup>3-</sup> metallodendrimer	S, H, B, C, IR, MS, TGA	[1410]
6K <sup>+</sup> 1,3,5-C <sub>6</sub> H <sub>3</sub> [3,5-C <sub>6</sub> H <sub>3</sub> [O(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> -O-(8-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co-(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ] <sub>3</sub> ] <sup>6-</sup> metallodendrimer	S, H, B, C, IR, MS, TGA	[1410]
8K <sup>+</sup> C <sub>6</sub> H <sub>4</sub> [ <i>p</i> -CH <sub>2</sub> OC <sub>6</sub> H <sub>3</sub> -2',4'-[CH <sub>2</sub> O[C <sub>6</sub> H <sub>3</sub> -2',4'-[CH <sub>2</sub> OCH <sub>2</sub> CH <sub>2</sub> OCH <sub>2</sub> CH <sub>2</sub> O-8'-(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ] <sub>2</sub> ] <sub>2</sub> ] <sup>8-</sup> metallodendrimer	S, H, B, C, IR, UV	[1420]
Si[(CH <sub>2</sub> ) <sub>2</sub> SiMe <sub>2</sub> (CH <sub>2</sub> ) <sub>2</sub> -μ(SiMe)-1,1'-[3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ] <sub>4</sub> ] <sup>8-</sup> metallodendrimers	S, H, B, C, Si, IR, MS	[945]
Si[(CH <sub>2</sub> ) <sub>2</sub> SiMe[(CH <sub>2</sub> ) <sub>2</sub> SiMe <sub>2</sub> (CH <sub>2</sub> ) <sub>2</sub> -μ(SiMe)-1,1'-[3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ] <sub>4</sub> ] <sup>8-</sup> metallodendrimers	S, H, B, C, Si, IR, MS	[945]
<i>Cyclo</i> -O <sub>4</sub> Si[(CH <sub>2</sub> ) <sub>2</sub> SiMe[(CH <sub>2</sub> ) <sub>2</sub> -μ(SiMe)-1,1'-[3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ] <sub>4</sub> ] <sup>8-</sup> metallodendrimers	S, H, B, C, Si, IR, MS	[945]
Polyanionic aryl ether photoluminescent dendrimers with μ(8,8')-SiMeH-3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> units	S, H, UV, MS, fluorescence	[946]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) <sup>-</sup> R=NH=C(O)Me, NH=C(O)Ph	S, H, B, C, MS	[948]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-N=CINH(CH <sub>2</sub> R')(CH <sub>2</sub> R'')] R=Me, Ph; R'=H, Et; R''=Et, <i>n</i> -C <sub>4</sub> H <sub>9</sub>	S, X(Me, H, <i>n</i> -C <sub>4</sub> H <sub>9</sub> ), H, B, C, MS	[948]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) R=NH <sub>3</sub> , NH <sub>2</sub> Et, NH <sub>2</sub> CH <sub>2</sub> Ph, NH(CH <sub>2</sub> Ph) <sub>2</sub>	S, X[NH(CH <sub>2</sub> Ph) <sub>2</sub> ], H, B, C, MS	[948]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-NH <sub>2</sub> (CH <sub>2</sub> ) <sub>3</sub> SO <sub>3</sub> ] <sup>-</sup>	S, H, B, C, MS	[948]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> Cl <sub>3</sub> ) <sub>2</sub> <sup>-</sup>	Gamma irradiation effects on Cs <sup>+</sup> and Sr <sup>2+</sup> extraction in polyethylene glycol	[949]
M <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> M <sup>+</sup> =Ag <sup>+</sup> , Ph <sub>3</sub> C <sup>+</sup>	S, H, B, IR	[952]
Ag(MeCN) <sub>2</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> B—H...Ag	X	[953]
Ag[NC(CH <sub>2</sub> ) <sub>n</sub> CN] <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> <i>n</i> =3,4	S, X, IR	[954]
Ag(η <sup>2</sup> - <i>p</i> -MeC <sub>6</sub> H <sub>4</sub> Me) <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup>	X	[955]
Ag(η <sup>2</sup> - <i>p</i> -MeC <sub>6</sub> H <sub>4</sub> Me) <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> -8,8'-μ-C <sub>6</sub> H <sub>4</sub>	X	[955]

Continued

Compound	Information	References
Ni(TMTAA) <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> grafted crown ether N <sub>4</sub> -macrocyclic polymer	S, X, H, MS	[957]
M[2.2.2]cryptate <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> supramolecular assembly	S, X, H	[958]
Sr(H <sub>2</sub> O) <sub>8</sub> (MeCN)(CTV) <sub>4</sub> (H <sub>2</sub> O) <sub>4</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> (CTV=cyclotrimeratrylene) host-guest H-bonded adamantoid network	S, X	[961]
3- <sup>60</sup> Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> ionic associates of incapacitating agents; solvent extraction	S	[963]
[ZnCl(Hpz-5-CMe <sub>3</sub> ) <sub>3</sub> ] <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> Hpz=pyrazole cobaltacarborane in bowl-shaped host cavity	S, X, H, MS	[964]
[2,6L <sub>2</sub> C <sub>6</sub> H <sub>3</sub> ]SnPh <sub>2</sub> <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> L=MeO, Me <sub>3</sub> CO stabilization by pincer-type triaryltin ligand	S, X, H, B, Sn, IR, MS	[965]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-LiC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup>	S, B	[914]
3-Co(1,2-LiC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> <sup>-</sup>	S, B	[914]
3-Co(1,2-MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> <sup>-</sup>	S, H, B, C	[914]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H	[914]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1,2-( <i>n</i> -C <sub>6</sub> H <sub>13</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	S, H, B, C	[914]
3-Co[1,2-( <i>n</i> -C <sub>6</sub> H <sub>13</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> <sup>-</sup>	S, H	[914]
3-Co[1,2-(MeOCH <sub>2</sub> CH <sub>2</sub> OCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> <sup>-</sup>	S, X, H, B, C	[914]
3-Co(1,2-PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> <sup>-</sup> extraction agent for <sup>137</sup> Cs <sup>+</sup> , <sup>90</sup> Sr <sup>2+</sup> , <sup>152</sup> Eu <sup>3+</sup>	X	[966]
3-Co[1,2-(CH=CH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> <sup>-</sup>	S, H, IR	[686]
3-Co[μ-1,2-(CH=CH-CH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> <sup>-</sup>	S, H, IR	[686]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1,2-(C <sub>4</sub> H <sub>6</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sup>-</sup>	S, IR, E	[687]
3-Co[1,2-(C <sub>4</sub> H <sub>6</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> <sup>-</sup>	S, H, IR	[687]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) <sup>-</sup> R=Ph, dioxane	S, X, H, B	[969]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6-Ph) <sub>2</sub> <sup>-</sup>	S, H, B, IR	[923,970]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-R) <sub>2</sub> <sup>-</sup> R=Me, Et, <i>n</i> -C <sub>6</sub> H <sub>13</sub> , Ph	S, B	[971]
3-Co(1,2-RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> <sup>-</sup> R=Ph, Me, EtS	S, H, B, C, IR, E	[926]
3-Co(1,2-MeRC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> <sup>-</sup> R=(CH <sub>2</sub> ) <sub>3</sub> O(CH <sub>2</sub> ) <sub>2</sub> OMe, (CH <sub>2</sub> ) <sub>3</sub> OCH <sub>2</sub> CHMe <sub>2</sub> , (CH <sub>2</sub> ) <sub>6</sub> O(CH <sub>2</sub> ) <sub>3</sub> Me, (CH <sub>2</sub> ) <sub>3</sub> Me	S, B[(CH <sub>2</sub> ) <sub>3</sub> O(CH <sub>2</sub> ) <sub>2</sub> OMe]	[972]
3-Co[1,2-(EtS)PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> <sup>-</sup>	S, H, B, C, IR, E	[926]
3-Co[1,2-[(CH <sub>2</sub> ) <sub>2</sub> OMe]MeC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> <sup>-</sup>	S, H, B, C, IR, E	[926]
3-Co[1,2-(EtS)MeC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> <sup>-</sup>	S, H, B, C, IR, UV, E	[926]
3-Co[1,2-(PhC <sub>6</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>10</sub> H <sub>11</sub> ] <sub>2</sub> <sup>-</sup>	S, X, H, B, C, IR, S	[1586]
3-Co[1,2-(SC <sub>4</sub> H <sub>3</sub> ) <sub>n</sub> C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ] <sub>2</sub> <sup>-</sup> n=1-3	S, H, B, C, MS, UV, E	[974]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1,2-(HOOC)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	S, IR	[975]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1,2-(ClOC)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	S, IR	[975]
3-Co[1,2-(ClOC)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> <sup>-</sup>	S, IR	[975]
3-Co[C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -C <sub>5</sub> H <sub>5</sub> NC(O)OMe] <sub>2</sub> Cl	S, B, H, IR, MS	[762]

Compound	Information	References
M <sup>+</sup> 3-Co[1,2-(Ph <sub>2</sub> P)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> <sup>-</sup> analogues of BINAP M=Ph <sub>3</sub> PAg, Ph <sub>3</sub> PAu, Me <sub>2</sub> C(O)Ag, Rh	S, X[Ph <sub>3</sub> PAg, Ph <sub>3</sub> PAu, Me <sub>2</sub> C(O)Ag], H, B, C, P, IR, MS	[976]
(Ph <sub>3</sub> P)ClPd{3-Co[1,2-(Ph <sub>2</sub> P)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> } analogue of BINAP	S, H, B, C, P, IR, MS	[976]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(1,1')-SiMe <sub>2</sub> <sup>-</sup>	S, X, H, B(2d), C, Si, IR, UV, MS	[978]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1,2-(Me <sub>3</sub> Si)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	S, X, H, B(2d), C, Si, IR, UV, MS	[978]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(1,1')-SiMeH <sup>-</sup>	S, H, B([2]d), C, Si, IR, UV, MS	[978]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> -μ(1,1')-SiMeR-μ(8,8')-C <sub>6</sub> H <sub>4</sub> <sup>-</sup>	S, H, B(2d), C, Si, IR, UV, MS	[978]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )[1,2-(Me <sub>3</sub> Si)C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]-μ(8,8')-C <sub>6</sub> H <sub>4</sub> <sup>-</sup>	S, H, B(2d), C, Si, IR, UV, MS	[978]
3-Co[1,2-(Me <sub>3</sub> Si)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> <sup>-</sup>	S, X, H, B(2d), C, Si, IR, UV, MS	[978]
3-Co[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -5-O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> ) <sub>2</sub> NHR <sup>-</sup> R=H, n-C <sub>4</sub> H <sub>9</sub> HIV protease inhibitor	S, X	[979]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OH)(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) <sup>-</sup> R=Me, Et	S, H, B, C, MS	[980]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-O(CH <sub>2</sub> ) <sub>2</sub> NH[C(O)OEt] <sub>2</sub> -C(O)Me <sup>-</sup>	S, X, H, B, C, IR	[982]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-L)(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-L) L=pyr, morpholine, PPh <sub>3</sub> , OPPh <sub>3</sub>	S, H, B, C, P, IR	[1409]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-L)(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-L) <sup>-</sup> L=Ph, C <sub>6</sub> H <sub>2</sub> Me <sub>3</sub>	S, H, B, S	[1409]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -8,8'-I <sup>-</sup> reactions with hindered Lewis bases → C-H activation in arenes		[1409]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -ONOPh) <sup>-</sup>	S, H, B, C	[966]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-X) <sup>-</sup> X=OP(O)(OH) <sub>2</sub> , OP(O)(OH)Ph	S, H, B(2d), P	[984]
3-Co[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-CH <sub>2</sub> Ph)(CH <sub>2</sub> CH <sub>2</sub> -O) <sub>2</sub> ] <sub>3</sub> Ln <sup>3+</sup>	X	[985]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R R=O(CH <sub>2</sub> ) <sub>4</sub> O, OC <sub>5</sub> H <sub>10</sub> , O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> X X=NC <sub>4</sub> H <sub>4</sub> , NC <sub>8</sub> H <sub>6</sub> , NC <sub>12</sub> H <sub>8</sub>	S, H, B, C, IR, MS	[986]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-L)(CH <sub>2</sub> CH <sub>2</sub> O) <sub>2</sub> <sup>-</sup> L=OPPH <sub>2</sub> , OCCH <sub>2</sub> , NR, C <sub>4</sub> H <sub>9</sub> , C <sub>12</sub> H <sub>25</sub> , CH <sub>2</sub> Ph selective extraction agents for lanthanide and actinide cations from highly acidic nuclear waste	S, H, B	[985]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O <sub>2</sub> (C <sub>2</sub> H <sub>4</sub> ) <sub>2</sub>	S, H, B, C, MS	[987]
[O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>5</sub> -O-8,8'-[(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ] <sup>-</sup>	ED <sub>50</sub> effects on living cells	[1555]
3-Co(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I) <sub>2</sub> <sup>-</sup>	ED <sub>50</sub> effects on living cells	[1555]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-Me) <sup>-</sup>	ED <sub>50</sub> effects on living cells	[1555]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> -1-(1,2-C <sub>2</sub> B <sub>10</sub> H <sub>11</sub> ) <sup>-</sup> (FF)	S, H, B, MS	[943]
1,2-[(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> -O-8-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>10</sub> H <sub>10</sub> ] <sup>2-</sup> (FF)	S, H, B, C, MS	[942]
1',n-[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O <sub>3</sub> (C <sub>2</sub> H <sub>4</sub> ) <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>4</sub> ] <sup>2-</sup> n=2',3',4'	S, H, B, C, MS	[987]
1',3',5'-[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O <sub>3</sub> (C <sub>2</sub> H <sub>4</sub> ) <sub>2</sub> -[C(O)] <sub>n</sub> ) <sub>3</sub> C <sub>6</sub> H <sub>4</sub> ] <sup>3-</sup> n=0,1	S, H, B, C, MS	[987]
1',2'-[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O <sub>2</sub> (C <sub>2</sub> H <sub>4</sub> ) <sub>2</sub> S] <sub>2</sub> C <sub>2</sub> B <sub>10</sub> H <sub>10</sub> ] <sup>2-</sup>	S, H, B, C, MS	[987]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(8,8')-NC(O)(CH <sub>2</sub> ) <sub>n</sub> P(O)PhR n=1, 2; R=Ph, n-C <sub>8</sub> H <sub>17</sub> extraction agents for lanthanides and actinides from nuclear waste	S, X(Ph, n=1), H, B, C, P, MS	[988]

Continued

Compound	Information	References
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OMe)(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) R = NR'C(O)CH <sub>2</sub> P(O)PhR'' R' = CH <sub>2</sub> Ph, R'' = Ph, <i>n</i> -C <sub>8</sub> H <sub>17</sub> extraction agents for lanthanides and actinides from nuclear waste	S, X(Ph, CH <sub>2</sub> Ph), H, B, C, P, MS	[988]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-NR(O)CH <sub>2</sub> P(O)Ph <sub>2</sub> ] R = H, Et, CH <sub>2</sub> Ph extraction agents for lanthanides and actinides from nuclear waste	S, H, B, C, P, MS	[988]
(Porphyrin)R <sub>4</sub> X <sub>8</sub> -N,N'-(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> -O-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> R = H, Ph; X = H, Et	S, H, MS, UV, E	[989]
(Porphyrin)Ph <sub>4</sub> -C <sub>6</sub> H <sub>4</sub> -NH-(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> -O-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, C, MS, UV, E	[989]
(Porphyrin)Ph <sub>4</sub> -C <sub>6</sub> H <sub>4</sub> -N[(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> -O-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> <sup>2-</sup>	S, H, C, MS, UV, E	[989]
(Porphyrin)[C <sub>5</sub> H <sub>3</sub> N(CH <sub>2</sub> -O) <sub>2</sub> -1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]Co[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ] <sub>4</sub> <sup>4-</sup>	S, H, C, UV, fluorescence spectroscopy	[990]
(Porphyrin)[(C <sub>6</sub> H <sub>4</sub> )-O-(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> -O-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>4</sub> <sup>4-</sup>	S, H, C, UV, fluorescence spectroscopy, light scattering, atomic force microscopy, inhibition of HIV-1 protease	[991]
(Porphyrin)[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-O(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> ] <sub>n</sub> <i>n</i> = 1–4	S, X( <i>n</i> = 1), H, C, UV, MS	[992]
(Porphyrin)[(C <sub>6</sub> H <sub>3</sub> )-3',5'-[O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> O(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>n</sub> (C <sub>6</sub> H <sub>5</sub> ) <sub>4-n</sub>	S, X( <i>n</i> = 2), H, C, U, fluorescence	[993]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>3</sub> -Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-(chlorine) <sub>6</sub>	S, H, B, UV, MS	[994]
Calix[4]arene[OCH <sub>2</sub> OCH <sub>2</sub> O(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>4</sub> <sup>4-</sup>	S, H, B, UV, calorimetry	[996]
Other calix[4]arene (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) derivatives	S	[995]
Resorc[4]arene (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) derivatives	S	[995]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> -NC <sub>4</sub> H <sub>4</sub> <sup>-</sup>	X	[986]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-O(CH <sub>2</sub> ) <sub>2</sub> OL L = OMe, OEt, O(CH <sub>2</sub> ) <sub>2</sub> OMe, O(CH <sub>2</sub> ) <sub>2</sub> OEt	S, X(OEt), H, B, C, IR	[997]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) R = <i>p</i> -C <sub>6</sub> H <sub>4</sub> OMe, Et, 2'-C <sub>4</sub> H <sub>3</sub> S	S, H, B	[998]
3-Co(1,2-RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sub>2</sub> <sup>+</sup> DD/LL, <i>meso</i>	X, E, UV (electrochemical)	[897]
3-Co(1,2-RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sub>2</sub>	UV (electrochemical), ESR (electrochemical)	[897]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SMe <sub>2</sub> )	XPS	[928]
	S, H, IR	[759]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SH) <sub>2</sub> <sup>-</sup>	S, B	[923]
2-Co(1,7-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(6,6')-S <sup>-</sup>	UV (detailed; photometric detection)	[1000]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(8,8')-S <sub>2</sub> <sup>-</sup>	UV (detailed; photometric detection)	[1000]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(4,8';8,4')-R <sub>2</sub> <sup>-</sup> R = C <sub>6</sub> H <sub>4</sub> , C <sub>6</sub> H <sub>3</sub> Et, C <sub>6</sub> H <sub>3</sub> Me	UV (detailed; photometric detection)	[1000]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(8,8')-O <sub>2</sub> P(O)R <sup>-</sup> R = Cl, NEt <sub>2</sub> , Ph	UV (detailed; photometric detection)	[1000]
3-Co[1,2-Ph(RS) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> <sup>-</sup> R = Me, Et, <i>n</i> -C <sub>4</sub> H <sub>9</sub>	S, H, B, C, IR	[1001]
[1,2-Ph(RS) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ]Co[1',9'-PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]-μ(1',8)-RS R = Me, Et, <i>n</i> -C <sub>4</sub> H <sub>9</sub>	S, X(Et), H, B, C, IR	[1001]
{(C <sub>5</sub> H <sub>4</sub> )Co(μ-S) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> }[CH=CHC(O)OMe](Me <sub>2</sub> C=CS <sub>2</sub> H) <sub>2</sub>	S, X, H, B, IR, MS	[1582]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -8,9-Cl <sub>2</sub> ) <sub>2</sub> <sup>-</sup>	S, H, B, UV, IR	[916]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -9,12-Cl <sub>2</sub> ) <sub>2</sub> <sup>-</sup>	S, H, B, UV, IR	[916]

Compound	Information	References
$(1,2\text{-C}_2\text{B}_9\text{H}_{10}\text{-8-Cl})\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_9\text{-8,9-Cl}_2)^-$	S, H, B, UV, IR	[916]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_8\text{-8,9,12-Cl}_3)_2^-$	S, H, B, UV, IR	[916]
	Molecular dynamics of liquid-liquid extraction in octanol-H <sub>2</sub> O system	[1006]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_8\text{-8,9,12-Cl}_3)_2^- (n\text{-C}_8\text{H}_{17})_3\text{NH}^+$	IR( $\nu_{\text{NH}}$ as a measure of acid strength)	[1007]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2^- \text{-}\mu(8,8')\text{-Br}$	S(electrochemical)	[1004]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10}\text{-8-Br})_2^-$	S	[1005,1008]
	H, B	[1005]
$(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10}\text{-8-I})_2^-$	X	[1014]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10}\text{-9-I})_2^-$	S	[971]
$(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10}\text{-8-I})^-$	S, H, B, UV, IR	[916]
$(1,2\text{-C}_2\text{B}_9\text{H}_{10}\text{-8-L})\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10}\text{-8-L})$ L = Me <sub>2</sub> S, C <sub>5</sub> H <sub>5</sub> N, Me <sub>3</sub> N, H <sub>3</sub> N	S, H, B, MS	[1016]
$(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{-}3\text{-Co}(1',2'\text{-C}_2\text{B}_9\text{H}_{10})\text{-}8\text{-O}(\text{CH}_2)_2\text{-O}(\text{CH}_2)_2\text{N}_3^-$ nucleoside conjugate	S(dipolar addition [chemical ligation]), H, B, IR, MS, UV	[766]
$(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{-}3\text{-Co}(1',2'\text{-C}_2\text{B}_9\text{H}_{10})\text{-}8\text{-O}(\text{CH}_2)_2\text{-O}(\text{CH}_2)_2\text{-O}(\text{CH}_2)_n\text{C}\equiv\text{CH}$	S(dipolar addition [chemical ligation]), H, B, IR, MS, UV	[766]
$(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{-}3\text{-Co}(1',2'\text{-C}_2\text{B}_9\text{H}_{10})\text{-}8\text{-O}(\text{CH}_2)_2\text{-O}(\text{CH}_2)_2\text{S}(\text{CH}_2)_3\text{SH}$	S(dipolar addition [chemical ligation]), H, B, IR, MS, UV	[766]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10}\text{-}n\text{-I})_2^-$ $n=4,9$ air-stable redox couple	S, H, B, C, IR, MS, E	[1573]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_8\text{-}4,9,12\text{-I}_3)_2^-$ air-stable redox couple	S, H, B, C, IR, MS, E	[1573]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_9\text{-}9,12\text{-Me}_2)_2^-$ air-stable redox couple	S, H, B, C, IR, MS, E	[1573]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_7\text{-}8,9,10,12\text{-Me}_4)_2^-$ air-stable redox couple	S, H, B, C, IR, MS, E	[1573]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_9)_2\text{-}\mu(4,8')\text{-}\mu(8,4')\text{-}(o\text{-C}_6\text{H}_4)_2^-$	S, X, H, B(2d), UV	[1018]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(8,8')\text{-MeC}_6\text{H}_3^-$	S, H, B, IR	[939]
$dl\text{-}3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(1,1')\text{-CH}_2\text{OCH}_2$	S, X, E	[1019]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(1,1')\text{-}(\text{CH}_2)_n^-$ $n=3\text{-}5$	X( $n=3, 4$ ), H, B, C, IR	[598]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(8,8')\text{-CH}_2\text{C}_9\text{H}_6$	S, X, H, B	[1021]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(8,8')\text{-NH}_2$	S, MS, UV	[1022]
	S, H, B, C, MS, UV, IR (diffuse reflectance)	[859]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(8,8')\text{-NH}_2(\text{CH}_2)_n$ polymers	S, H, B, C, IR (diffuse reflectance)	[859]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(8,8')\text{-NHMe}$	S, MS	[1022]
$2\text{-Co}(1,7\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(8,8')\text{-N}_2\text{C}_3\text{H}_2\text{C}(\text{O})\text{OH}$	S, X, H, B, IR, MS	[775]
$2\text{-Co}(1,7\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(8,8')\text{-N}_2\text{C}_3\text{H}_2\text{C}(\text{O})\text{OMe}^-$	S, X	[1023]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(8,8')\text{-NHCH}_2\text{C}(\text{O})\text{OMe}$	X	[1024]
$[1,2\text{-}(\text{C}_4\text{H}_8\text{NCH}_2)_2\text{C}_2\text{B}_9\text{H}_{10}]\text{Co}[1,2\text{-}(\text{C}_4\text{H}_8\text{NCH}_2)_2\text{C}_2\text{B}_9\text{H}_{10}]$ inter- and intramolecular H-bonding	S, X, H, B, C, MS	[1025]
$[1,2\text{-}(\text{C}_5\text{H}_{10}\text{NCH}_2)_2\text{C}_2\text{B}_9\text{H}_{10}]\text{Co}[1,2\text{-}(\text{C}_5\text{H}_{10}\text{NCH}_2)_2\text{C}_2\text{B}_9\text{H}_{10}]$ inter- and intramolecular H-bonding	S, X, H, B, C, MS	[1025]
$3\text{-Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}\mu(8,8')\text{-NOR}_2$ R=H, CMe <sub>2</sub> , CHPh, Me	S, H, B, MS	[1028]
$(\text{C}_2\text{B}_9\text{H}_{10}\text{-OH})\text{Co}(\text{C}_2\text{B}_9\text{H}_{10}\text{-NR}_2)$ R=H, CMe <sub>2</sub> , CHPh, Me	S, H, B, MS	[1028]

Continued



Compound	Information	References
$(C_2B_9H_{11})Co(C_2B_9H_{10})-8-PMe_2$	S, H, B, P, MS	[1029]
$(1,7-C_2B_9H_{11})Co(1,7-C_2B_9H_{10})-6-PMe_2$	S, X, H, B, P, MS	[1029]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-O^-$	X	[857,1031]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-OEt$	S, UV	[1022]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-O_2CMe$	S, H, B, IR	[776]
	X	[977]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-\{cyclo-[p-C_6H_4-NMe]_2C=O\}$ urea	S, X, H	[1033]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-O_2P(O)X$ X=Cl, OH extraction agents for $Eu^{3+}$ in 1M $HNO_3$	S, X(Cl)	[984]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-O_2P(O)NEt_2$	S, X(absolute configuration)	[984]
$2-Co(1,7-C_2B_9H_{10})_2-\mu(6,6')-X$ X= $S^-$ , SMe ( <i>d,l</i> only)	S, H, B, CD	[1034]
$2-Co(1,7-C_2B_9H_{10})_2-\mu(6,6')-S^-$	X	[1035]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-SMe$	S, H, B, MS	[1036]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-SMe_2^-$	XPS	[928]
$3-Co(C_2B_9H_{10})_2-\mu(8,8')-S_2CH$	S, H, IR	[776]
	X	[1038]
	B	[923]
$2-Co(1,7-C_2B_9H_{10})_2-\mu(8,8')-S_2CH$	S, IR, MS	[776]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-X$ (X= $S_2CH$ , SMe, $S_2Me$ , OMe)	XPS	[928]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-S_2CR-1,8-C_{10}H_5NMe_2H$ R=H, OH	S, X(R=H, OH in same crystal), H, B, IR, MS, COND	[1040]
<i>Meso</i> - $3-Co(1,2-C_2B_9H_{10})_2-\mu[1,1'-S(CH_2)_2]^-$	S, X, H, B, C, MS	[1041]
<i>Meso</i> - $3-Co(1,2-C_2B_9H_{10})_2-\mu[1,1'-SO_2(CH_2)_2]^-$	S, X, H, B, C, MS	[1041]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')SCH_2C(O)OMe$	X	[1042]
$[1,2-C_2B_9H_{10}(SH)]Co[1,2-C_2B_9H_{10}(SCHO)]^-$	S, H, IR	[776]
$Na^+ 3-Co(1,2-C_2B_9H_{10})_2-\mu(1,1')-[S(CH_2CH_2O)_3CH_2CH_2-S]^-$	S, X, H, B, C, IR	[997]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-X$ X=SEt, $SC_4H_9$ , $SCH_2CH=CH_2$	S, MS	[981]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-X$ X= $SCH_2Ph$ , $SCH_2C(O)OH$ , $SCH_2C(O)OMe$	S, MS	[981]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-EME$ E=Se, Te	S, H, B, MS, UV	[1022]
$3-LL'Co_2(1,2-C_2B_9H_{11})_2$ L,L'=CO, $CNCMe_3$ , $PMe_2Ph$	S, X[(CO) $_2$ ; CO, $PMe_2Ph$ ], H, B, C, IR	[911]
$3-Co(1,2-C_2B_9H_{11})_2^- Cu(1,10-phenanthroline)^{3+} \cdot MeCN$	X	[1436]
$3-Co(1,2-C_2B_9H_{11})_2-8,8'-[O(CH_2OCH_2)]_2R^-$ $K^+$ R= $\alpha$ - $S_2C_6H_4$ , S	X( $\alpha$ - $S_2C_6H_4$ ), H, B, C, IR, MS	[1589]
$Cs^+ 3-Co(1,2-C_2B_9H_{11})_2^-$ $Cs^+$ interactions with calix[4]arene-bis( <i>t</i> -octylbenzo)-18-crown-6	H	[1446]
$3-Co(1,2-C_2B_9H_{10}-8-X)_2^-$ X=H, I	Anions cross through synthetic lipid bilayer membranes	[1536]
$3-Co(1,2-C_2B_9H_{10})_2-\mu(8,8')-PO_3H$	Catalyst for ketimine hydrogenation to amines	[1537]
$3-Co(1,2-C_2B_9H_{11})_2^-$ in ion-selective electrodes for protonatable <i>N</i> -containing analytes dopamine, nicotinic acid, nicotinamide, histamine and metformin in aqueous solutions	H, B, C, IR, MS	[1521]

Compound	Information	References
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> X) <sub>2</sub> <sup>-</sup> X=H, Cl surfactant properties, aggregate formation	E, MS	[1525]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> (CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> NHCH <sub>2</sub> -chlorin-e6 <sup>+</sup> conjugate accumulation in cancer cells for BNCT and phototherapy	Fluorescence, cytotoxicity toward human lung adenocarcinoma	[1514]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> structure of composite with poly(ethylene oxide)	S, solid state NMR (H, B, C, Na), wide angle X-ray scattering, isothermal titration calorimetry	[1441]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> conjugates with proteins in protein data bank (PDB) and HIV for BNCT		[1453]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> Cl <sub>3</sub> ) <sub>2</sub> <sup>-</sup> conjugates with 5-ethynyl-2'-deoxyuridine	S, H, B, C, MS, cytotoxicity	[1497]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> counterion in Ba <sup>2+</sup> beauvercin complexes		[1486]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> incorporation into DABCO (diazabicyclooctane) MOFs Ag...H-B interactions	S, X	[1490,1550]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> (Ph <sub>3</sub> P) <sub>3</sub> Ag-X-Ag(Ph <sub>3</sub> P) <sub>3</sub> <sup>+</sup> X=Cl, Br	S, X, IR	[1510]
Cyclo-C <sub>6</sub> [(Me <sub>2</sub> CH-C <sub>6</sub> H <sub>3</sub> ) <sub>2</sub> (μ-O(CH <sub>2</sub> ) <sub>4</sub> O) <sub>2</sub> ] <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> derivatives.	Hexaarylbenzene-based receptor interaction of HO(CH <sub>2</sub> ) <sub>2</sub> NH <sub>3</sub> <sup>+</sup> cation	[1475]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> R] <sup>-</sup> R=NHC(O)NH <sub>3</sub> <sup>+</sup> , SC(=NH <sub>2</sub> <sup>+</sup> )NH <sub>2</sub> , NHC(=NH <sub>2</sub> <sup>+</sup> )NH <sub>2</sub> , N[(CH <sub>2</sub> ) <sub>2</sub> OH] <sub>3</sub> <sup>+</sup> , [O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>4</sub> OH, [O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>4</sub> O <sup>-</sup> , OC <sub>6</sub> H <sub>4</sub> - <i>p</i> -C(O)OMe	Light scattering/(solubility) suppression of self-assembly; biocompatibility	[1439]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I) <sup>-</sup> Pd-catalyzed B-C <sub>vinyl</sub> coupling		[1448]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> -8-IR <sup>+</sup> R=Ph, C <sub>6</sub> H <sub>4</sub> - <i>p</i> -OMe reaction with nucleophiles to form carbonitrile, pyridinium, sulfonium, thiol, acetoxy and amino derivatives	S, H, B, C	[1574]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> -8-R R=I, SCHNMe <sub>2</sub> <sup>+</sup>	S, H, B, C, MS	[1574]
(C <sub>5</sub> H <sub>4</sub> N) <sub>2</sub> [(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> O-8-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-X)] <sub>2</sub> X=H, I probe for sequential voltage tuning	S, E	[1580]
C <sub>5</sub> H <sub>4</sub> NMe <sub>2</sub> <sup>2+</sup> 2[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-X)] <sup>-</sup> X=H, I probe for sequential voltage tuning	S, E	[1580]
(C <sub>5</sub> H <sub>4</sub> NMe) <sub>2</sub> <sup>2+</sup> 2[(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11-n</sub> ) <sub>2</sub> Co <sup>-</sup> n=0,1,3 probe for sequential voltage tuning	S, E	[1580]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-CH=CHR) <sup>-</sup> R=Ph, <i>p</i> -C <sub>6</sub> H <sub>4</sub> OH, <i>n</i> -C <sub>4</sub> H <sub>9</sub> , <i>n</i> -C <sub>8</sub> H <sub>17</sub> , CN, OH	S, X(Ph), H, B, C, MS	[1448]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-CH=CH-C <sub>6</sub> H <sub>4</sub> - <i>p</i> -R) <sub>2</sub> <sup>-</sup> R=Me, F, Cl, Br, <i>m</i> -Br	S, X(Ph), H, B, C, MS	[1448]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) <sup>-</sup> adenosine and 2'-deoxyadenosine derivatives	S, H, C	[1458]
H <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> monolayer vesicles → micelles	X-ray and neutron scattering	[1450]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -8,8'-C <sub>6</sub> H <sub>4</sub> <sup>-</sup> ansa-metallacyclophane	S, H, B, C, IR, MS, UV, E	[1456]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -1.1'-P(E)Ph <sup>-</sup> E=O, S, Se ansa-metallacyclophanes	S, H, B, C, IR, MS, UV, E	[1456]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -1.1'-P(E)CMe <sub>3</sub> <sup>-</sup> E=O, S, Se ansa-metallacyclophanes	S, X, H, B, C, IR, MS, UV, E	[1456]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> -1.1'-P(E)X-8,8'-C <sub>6</sub> H <sub>4</sub> <sup>-</sup> X=PCMe <sub>3</sub> , PPh, P(E)Ph; E=O, S, Se diansa-metallacyclophanes	S, X, H, B, C, IR, MS, UV, E	[1456]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -10-I) <sub>2</sub> <sup>-</sup>	S, X, H, B, C, IR, MS	[1459]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -8,8'-CH <sub>2</sub> OMe	S, X, H, B, C, IR, MS	[1460]

Continued

Compound	Information	References
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OMe <sub>2</sub> )	S, H, B, C, IR, MS	[1460]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OMe <sub>2</sub> ) <sup>2-</sup>	S, H, B, C, IR, MS	[1460]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-CH <sub>2</sub> R)(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OMe) R=py, NH <sub>2</sub> - <i>n</i> -C <sub>6</sub> H <sub>13</sub> , NH <sub>2</sub> C <sub>2</sub> H <sub>4</sub> OH, PPh <sub>3</sub>	S, X(py, NH <sub>2</sub> C <sub>2</sub> H <sub>4</sub> OH, PPh <sub>3</sub> ), H, B, C, IR, MS	[1460]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-CH <sub>2</sub> -OC <sub>6</sub> H <sub>4</sub> CMe <sub>3</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OMe) <sup>-</sup>	S, H, B, C, IR, MS	[1460]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -8,8-CH <sub>2</sub> O <sup>-</sup>	S, X, H, B, C, IR, MS	[1460]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-CH <sub>2</sub> -O-calix[4]-arene(OCH <sub>2</sub> CN) <sub>2</sub> (OMe)(CMe <sub>3</sub> ) <sub>4</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OMe) <sup>-</sup>	S, X, H, B, C, IR, MS	[1460]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-(OCH <sub>2</sub> CH <sub>2</sub> ) <sub>n</sub> OH] <sup>-</sup> M <sup>+</sup> n=2,4,6,8,10 M=cyclo-MeNC <sub>3</sub> H <sub>3</sub> N- <i>n</i> -C <sub>4</sub> H <sub>9</sub> , cyclo-C <sub>5</sub> H <sub>5</sub> N- <i>n</i> -C <sub>4</sub> H <sub>9</sub> , cyclo-C <sub>4</sub> H <sub>8</sub> NMe- <i>n</i> -C <sub>4</sub> H <sub>9</sub> , cyclo-C <sub>5</sub> H <sub>10</sub> NMe- <i>n</i> -C <sub>4</sub> H <sub>9</sub> glycolated low-melting salts	S, H, DSC, glass transition, MP	[1463]
3-Co[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-CMPO-(CH <sub>2</sub> -CH <sub>2</sub> O) <sub>2</sub> ][1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-X] <sup>-</sup> X=Cl, Br, I CMPO=Ph <sub>2</sub> P(O)CH <sub>2</sub> C(O)( <i>t</i> -C <sub>8</sub> H <sub>17</sub> )N for extraction of lanthanides and actinides from high-level activity nuclear waste		[1464]
3-Co[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-CMPO-(CH <sub>2</sub> -CH <sub>2</sub> O) <sub>2</sub> ][1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-X] <sup>-</sup> X=Cl, Br, I CMPO=Ph <sub>2</sub> P(O)CH <sub>2</sub> C(O)( <i>t</i> -C <sub>8</sub> H <sub>17</sub> )N for extraction of Am <sup>3+</sup> from high-level activity nuclear waste		[1466]
3-Co[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-(O-CH <sub>2</sub> ) <sub>2</sub> -N(C <sub>8</sub> H <sub>17</sub> )-C(O)CH <sub>2</sub> ] <sub>2</sub> O for extraction of Am <sup>3+</sup> from high-level activity nuclear waste		[1466]
3-Co[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-CMPO-(CH <sub>2</sub> -CH <sub>2</sub> O) <sub>2</sub> ][1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-X] <sup>-</sup> X=Cl, Br, I CMPO=Ph <sub>2</sub> P(O)CH <sub>2</sub> C(O)( <i>t</i> -C <sub>8</sub> H <sub>17</sub> )N for extraction of M <sup>2+</sup> in H <sub>2</sub> O-nitrobenzene system M=Ba, Sr, Mg, Ni, Co, Zn, Cd, Mn, Cu, Pb		[1467]
3-Co[(1,2-HO(CH <sub>2</sub> ) <sub>n</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ] <sup>-</sup> Cs <sup>+</sup> n=1-3 alkylhydroxy derivatives	S, X(n=2), H, B, C, MS	[1468]
<i>syn/anti</i> -3-Co[(1,2-HO(CH <sub>2</sub> ) <sub>n</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> ] <sup>-</sup> Cs <sup>+</sup> n=1-3 alkylhydroxy derivatives	S, H, B, C, MS	[1468]
3-Co{1,2-[HO(CH <sub>2</sub> ) <sub>3</sub> ] <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> }{1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ] <sup>-</sup> Cs <sup>+</sup> alkylhydroxy derivatives	S, H, B, C, MS	[1468]
3-Co[(1,2-(HO) <sub>2</sub> (P(O)(CH <sub>2</sub> ) <sub>3</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ] <sup>-</sup> Cs <sup>+</sup> alkylhydroxy derivatives	S, X, H, B, C, MS	[1468]
[1,2-HOP(O){O(CH <sub>2</sub> ) <sub>3</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> Co[1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ]} <sub>2</sub> ] <sup>2-</sup> Cs <sup>+</sup> alkylhydroxy derivatives	S, X, H, B, C, MS	[1468]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-oligonucleotide) <sup>-</sup> electrochemical evaluation of DNA hybridization	E	[1473]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R)(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-R') <sup>-</sup> R, R'=H, OH, NH <sub>2</sub> ; RR'=μ-NH <sub>2</sub> <sup>+</sup> , μ-S, μ-O <sub>2</sub> PO <sub>2</sub> ; R=O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> R' R'=thiourea, urea, guanidine, arginine, amines, and related compounds inhibition of NO synthases activation	S, H, B, MS, fluorescence	[1474]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I)(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'- <i>n,m</i> -Me <sub>2</sub> C <sub>6</sub> H <sub>3</sub> ) <sup>-</sup> n,m=3,4; 3,5; 7,5	S, X(2,5), H, B, C, MS	[1476]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I)(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-R) <sup>-</sup> R= <i>p</i> -C <sub>6</sub> H <sub>4</sub> OMe, OPh, C <sub>6</sub> H <sub>4</sub> NMe	S, H, B, C, MS	[1476]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I)(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-PEt <sub>3</sub> )	S, X, H, B, C, MS	[1476]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-cholesteroyl) incorporation into liposomes; bidistribution in mice BNCT	S, H, B, C, IR, MS	[1485]
3,6,1,2-(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> Co <sub>2</sub> (C <sub>2</sub> B <sub>8</sub> H <sub>10</sub> )- <i>n,n'</i> -(OH) <sub>2</sub> <sup>2-</sup> n, n'=8,10; 8',10; 8',8	S, X[8,10-(OH) <sub>2</sub> ], H, B, C, IR, HIV protease inhibition	[1478]

Compound	Information	References
3,6,1,2-(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> Co <sub>2</sub> (C <sub>2</sub> B <sub>8</sub> H <sub>10</sub> )- <i>n,n'</i> -(OH) <sub>2</sub> <sup>2-</sup> <i>n, n'</i> = 8,10; 8',10; 8',8	S, H, B, C, IR, HIV protease inhibition	[1478]
3,6,1,2-(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> Co <sub>2</sub> (C <sub>2</sub> B <sub>8</sub> H <sub>9</sub> )-8-R <sup>-</sup> R=OC <sub>4</sub> H <sub>8</sub> , O(CH <sub>2</sub> ) <sub>4</sub> L L=C <sub>4</sub> H <sub>9</sub> NH <sub>2</sub> , C <sub>5</sub> H <sub>5</sub> N, Et <sub>3</sub> N, PPh <sub>3</sub> , OC <sub>6</sub> H <sub>4</sub> CMe <sub>3</sub> , OC <sub>6</sub> H <sub>4</sub> OMe	S, H, B, C, IR, HIV protease inhibition	[1478]
<i>n</i> -C <sub>4</sub> H <sub>9</sub> )HN((CH <sub>2</sub> ) <sub>4</sub> O-8-[3,6,1,2-(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> Co <sub>2</sub> ] <sub>2</sub> ) <sup>3-</sup>	S, H, B, C, IR, HIV protease inhibition	[1478]
[3-Co(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> ) <sub>2</sub> NHR <sup>-</sup> Na <sup>+</sup> R=H, <i>n</i> -C <sub>4</sub> H <sub>9</sub> , CH <sub>2</sub> CH <sub>2</sub> OH, CMe <sub>3</sub> , C(CH <sub>2</sub> OH) <sub>3</sub> , CH <sub>2</sub> Ph, (CH <sub>2</sub> ) <sub>3</sub> SO <sub>3</sub> <sup>-</sup> , 1,2-C <sub>2</sub> B <sub>10</sub> H <sub>11</sub> , <i>nido</i> -7,8-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> <sup>-</sup> metallocarborane pharmacophores	Solubility, lipophilicity (octanol-H <sub>2</sub> O partition coefficient)	[1488]
[3-Co(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> ) <sub>2</sub> NR <sub>2</sub> <sup>2-</sup> 2Na <sup>+</sup> R <sub>2</sub> =SO <sub>2</sub> Ph, 8,8'-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> Co, 2Et metallocarborane pharmacophores	Solubility, lipophilicity (octanol-H <sub>2</sub> O partition coefficient)	[1488]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-8-O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> OC <sub>6</sub> H <sub>4</sub> - <i>o</i> -CH <sub>2</sub> Ph <sup>-</sup> conjugates with poly(2-ethylloxazoline) (PEOX) and polyethylene oxide biocompatible hydrophilic polymers; hybrid nanospheres	S, H, light-scattering, ATM, isothermal titration calorimetry, SAXS, transmission electron microscopy	[1496]
HRN((CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> O-8-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )) <sub>2</sub> <sup>-</sup> R= <i>n</i> -C <sub>4</sub> H <sub>9</sub> , H, (CH <sub>2</sub> ) <sub>2</sub> OH, (CH <sub>2</sub> ) <sub>3</sub> C(O)OH, B <sub>10</sub> H <sub>9</sub> <sup>2-</sup> conjugates with poly(2-ethylloxazoline) (PEOX) and polyethylene oxide biocompatible hydrophilic polymers; hybrid nanospheres	S, H, light-scattering, ATM, isothermal titration calorimetry, SAXS, transmission electron microscopy	[1496]
3-Co(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) <sup>-</sup> R=C(O)H, Me, Et	Antimicrobial activity	[1501]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I) <sub>2</sub> <sup>-</sup> H <sup>+</sup> metallocarborane amphiphiles monolayered $\Theta$ -shaped lyotropic lamellar phases	Wide-angle X-ray scattering (SWAXS)	[1503]
[3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -)](CMPO) <sub>3</sub> - <i>t</i> -butyl-calix[4]arene] <sup><i>n</i>-</sup> derivatives CMPO= carbamoyl methyl diphenyl phosphine oxide extraction of lanthanide and actinide 3+ cations	S, H, B, extraction of M <sup>3+</sup> ions M=Eu, Am	[1506]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -8,8'-S <sub><i>n</i></sub> -(CH <sub>2</sub> ) <sub><i>m</i></sub> -estradiol <i>n</i> =0,1 <i>m</i> =1,2	S, X( <i>n</i> =1, <i>m</i> =1), activation of $\alpha$ and $\beta$ estrogen receptors	[1507]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-R) R=deoxyuridine, deoxyguanosine phosphoramidites DNA oligomers	S, H, B, C, P, UV, MS, lipophilicity, antisense activity in MCF-7 cells, resistance to degradation by snake venom	[1508]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-CH=CH <sub>2</sub> ) <sub>2</sub>	S(Pd-catalyzed cross-coupling), IR, MS	[1512]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-O(CH <sub>2</sub> ) <sub>5</sub> R] <sup>-</sup> R=N <sub>3</sub> , NC <sub>2</sub> HR'; R'=CH <sub>2</sub> OH, (CH <sub>2</sub> ) <sub>2</sub> OH, Ph, C <sub>5</sub> H <sub>4</sub> FeCp	S, H, B, C, MS	[1517]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> N <sub>3</sub> C <sub>2</sub> HR] <sup>-</sup> R=CH <sub>2</sub> OH, (CH <sub>2</sub> ) <sub>2</sub> OH, Ph, (CH <sub>2</sub> ) <sub>6</sub> Me, C <sub>5</sub> H <sub>4</sub> FeCp, hydroxyestraacetate	S, H, B, C, MS	[1517]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-O(CH <sub>2</sub> ) <sub>2</sub> -X-(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> C $\equiv$ CH] <sup>-</sup> X=O, CH <sub>2</sub>	S, H, B, C, MS	[1517]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-O(CH <sub>2</sub> ) <sub>2</sub> -X-(CH <sub>2</sub> ) <sub>2</sub> -O-(CH <sub>2</sub> ) <sub>2</sub> CHN <sub>3</sub> -CH <sub>2</sub> C(O)OMe] <sup>-</sup> X=O, CH <sub>2</sub> triazole	S, H, B, C, MS	[1517]
N <sub>3</sub> C <sub>2</sub> H-C-N-((CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> -O-8-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )-3-Co(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )) <sub>2</sub> ) <sub>2</sub> <sup>2-</sup> triazole	S, H, B, C, MS	[1517]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )( $\mu$ -H) <sub>3</sub> CuPh	S, X, H, B, P	[1518]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> Cu(PPh <sub>3</sub> ) <sub>3</sub> <sup>+</sup>	S, X, H, B, P	[1518]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-C $\equiv$ C-C <sub>6</sub> H <sub>3</sub> -3',5'-[C(O)OH] <sub>2</sub> ) <sub>2</sub> <sup>-</sup> PPh <sub>4</sub> <sup>+</sup> incorporation in Cu <sub>2</sub> O <sub>8</sub> paddlewheel MOFs with 3 topologies	S, X, H, B	[1520]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) R=O(CH <sub>2</sub> ) <sub>2</sub> -O(CH <sub>2</sub> ) <sub>2</sub> NMe <sub>2</sub> (CH <sub>2</sub> ) <sub>3</sub> NH- <i>bicyclo</i> -C <sub>5</sub> H <sub>4</sub> N <sub>4</sub> (NH <sub>2</sub> )- <i>cyclo</i> -C <sub>4</sub> H <sub>5</sub> (OH)(CH <sub>2</sub> OH)O deoxyadenosine	S, H, B, C, cytotoxicity	[1522]

Continued

Compound	Information	References
3- <b>Co</b> (1,2- $C_2B_9H_{11}$ )(1,2- $C_2B_9H_{10}$ -8-R) <sup>-</sup> R = O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> SMe <sub>2</sub> <sup>+</sup> , O(CH <sub>2</sub> ) <sub>2</sub> OH, O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> SMe, O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> SC <sub>4</sub> H <sub>4</sub> O <sup>+</sup> , O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> S(CH <sub>2</sub> ) <sub>2</sub> C[C(O)OMe][NH(O)OCMe <sub>3</sub> , O(CH <sub>2</sub> ) <sub>2</sub> O-(CH <sub>2</sub> ) <sub>2</sub> S <sup>+</sup> -9'- <i>nido</i> -7',8'- $C_2B_9H_9$ <sup>-</sup> , O(CH <sub>2</sub> ) <sub>2</sub> O(CH <sub>2</sub> ) <sub>2</sub> PPh <sub>3</sub> <sup>+</sup> , O(CH <sub>2</sub> ) <sub>2</sub> O <sub>3</sub> SMe, O(CH <sub>2</sub> ) <sub>2</sub> N <sub>3</sub> , O(CH <sub>2</sub> ) <sub>2</sub> (NH <sub>2</sub> ) <sub>2</sub> <sup>+</sup> , O(CH <sub>2</sub> ) <sub>2</sub> SH, O(CH <sub>2</sub> ) <sub>2</sub> N <sub>3</sub> C <sub>2</sub> HPh	S, H, B, C, MS	[1577]
3- <b>Co</b> (1,2- $C_2B_9H_{11}$ ) <sub>2</sub> <sup>-</sup> imaging in living cells	Raman (B-H)	[1524]
3- <b>Co</b> (1,2- $C_2B_9H_{11}$ )(1,2- $RC_2B_9H_{10}$ ) <sup>-</sup> R = C(O)OH, CH <sub>2</sub> C(O)OH, (CH <sub>2</sub> ) <sub>2</sub> C(O)OH, OCH <sub>2</sub> C(O)OH, (CH <sub>2</sub> ) <sub>n</sub> C(O)-C <sub>6</sub> H <sub>4</sub> - <i>p</i> -NO <sub>2</sub> (n = 1, 2), (CH <sub>2</sub> ) <sub>n</sub> C(O)NHR n = 1, 2; R = <i>n</i> -C <sub>4</sub> H <sub>9</sub> , CH <sub>2</sub> Ph	S, H, B, C	[1528]
3- <b>Co</b> (1,2- $C_2B_9H_{11}$ )(1,2- $RC_2B_9H_{10}$ )-1-CH <sub>2</sub> C(O)NH(CH <sub>2</sub> ) <sub>2</sub> -1'-(1',2'- $C_2B_9H_{10}$ )Co(1',2'- $C_2B_9H_{11}$ ) <sub>2</sub> <sup>2-</sup>	S, H, B, C	[1528]
3- <b>Co</b> (1,2- $C_2B_9H_{11}$ )(1,2- $C_2B_9H_{10}$ -8-C <sub>6</sub> H <sub>4</sub> -R) <sup>-</sup> R = <i>p</i> -CH=CH <sub>2</sub> , <i>m/p</i> -CHO	S(Pd-catalyzed B-C cross-coupling), X, H, B, C, IR, MS	[1531]
3- <b>Co</b> (1,2- $C_2B_9H_{10}$ -8-I)(1',2'- $C_2B_9H_{10}$ -8'-R) R = <i>cyclo</i> -S(CH <sub>2</sub> ) <sub>2</sub> (O)(CH <sub>2</sub> ) <sub>2</sub> , ONPy, <i>cyclo</i> -O(CH <sub>2</sub> ) <sub>5</sub> charge-compensated ligand	S, X(ONpy), H, B, C, MS	[1533]
3- <b>Co</b> (1,2- $C_2B_9H_{11}$ )(1,2- $C_2B_9H_{10}$ )-8-O[(CH <sub>2</sub> ) <sub>2</sub> O] <sub>n</sub> -8'-(1,2- $C_2B_9H_{10}$ ) <b>Co</b> (1,2- $C_2B_9H_{11}$ ) n = 5, 6 aqueous self-assembly and cation selectivity	S, X(n = 6), H, B, C, IR, MS, dynamic light scattering, cation exchange with H <sup>+</sup> , Li <sup>+</sup> , Rb <sup>+</sup> , Cs <sup>+</sup> , and NMe <sub>4</sub> <sup>+</sup>	[1535]
3- <b>Co</b> (1,2- $C_2B_9H_{11}$ )(1,2- $C_2B_9H_{10}$ )-8-O[(CH <sub>2</sub> ) <sub>2</sub> O](CH <sub>2</sub> ) <sub>n</sub> C <sub>6</sub> H <sub>4</sub> - <i>p</i> -(CH <sub>2</sub> ) <sub>n</sub> O[(CH <sub>2</sub> ) <sub>2</sub> O]-8'-(1,2- $C_2B_9H_{10}$ ) <b>Co</b> (1,2- $C_2B_9H_{11}$ ) n = 0, 1 aqueous self-assembly and cation selectivity	S, H, B, C, IR, MS, dynamic light scattering, cation exchange with H <sup>+</sup> , Li <sup>+</sup> , Rb <sup>+</sup> , Cs <sup>+</sup> , and NMe <sub>4</sub> <sup>+</sup>	[1535]
(fumaramide rotaxanes) <sup>2+</sup> {3- <b>Co</b> (1,2- $C_2B_9H_8$ Br <sub>3</sub> )(1,2- $C_2B_9H_7$ Br <sub>3</sub> )-8-O[(CH <sub>2</sub> ) <sub>2</sub> O] <sub>2</sub> } <sup>-</sup> pirouetting motion	S, H	[1538]
3- <b>Co</b> (1,2- $C_2B_9H_{10}$ -8- <sup>125</sup> I)(1',2'- $C_2B_9H_{10}$ )-8'-[O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> -OC(O)Ph <sup>-</sup> in vivo imaging	S, IR, MS, biodistribution in rodents	[1545]
3- <b>Co</b> (1,2- $C_2B_9H_{10}$ -8-R)(1',2'- $C_2B_9H_{10}$ -8'-R') <sup>-</sup> R = I; R' = OH, OMe, OEt, OBu, OCH <sub>2</sub> C≡CH, O(CH <sub>2</sub> ) <sub>2</sub> C≡CH, O(CH <sub>2</sub> ) <sub>2</sub> Br, O(CH <sub>2</sub> ) <sub>2</sub> OH, Br	S, X(OH), H, B, C, MS	[1548]
3- <b>Co</b> (1,2- $C_2B_9H_{10}$ -8-R)(1',2'- $C_2B_9H_{10}$ -8'-R') <sup>-</sup> R = OH; R' = OH, OBu	S, H, B, C, MS	[1548]
3- <b>Co</b> (1,2- $C_2B_9H_{10}$ -8-OBu) <sub>2</sub> <sup>-</sup>	S, H, B, C, MS	[1548]
3- <b>Co</b> (1,2- $C_2B_9H_{11}$ )(1',2'- $C_2B_9H_{10}$ ) <sup>-</sup> -8'-[O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> -X-(CH <sub>2</sub> ) <sub>3</sub> SiO <sub>3</sub> -silica X = NH <sub>2</sub> <sup>+</sup> , S stationary phases for HPLC		
Na <sup>+</sup> {3- <b>Co</b> (1,2- $C_2B_9H_{11}$ )(1',2'- $C_2B_9H_{10}$ )-8'-[O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> [O(CH <sub>2</sub> ) <sub>4</sub> ] <sub>n</sub> OH} <sup>-</sup> n = 1-22 redox-active polymers for surface modification	S, H, B, C, IR, MS, E	[1551]
Na <sup>+</sup> {3- <b>Co</b> (1,2- $C_2B_9H_{10}$ ) <sub>2</sub> -μ(8,8')-(CH <sub>2</sub> ) <sub>4</sub> [O(CH <sub>2</sub> ) <sub>2</sub> ] <sub>2</sub> O <sup>-</sup>	S, X, H, B, C, IR, MS, E	[1551]
3- <b>Co</b> (1,2- $RC_2B_9H_{10}$ )(1',2'- $C_2B_9H_{11}$ ) <sup>-</sup> R = <i>bicyclo</i> -COC(CH <sub>2</sub> ) <sub>n</sub> NC(O)-C <sub>6</sub> H <sub>4</sub> n = 2, 3 oxazolo-isoindalone	S, X, H, B, C, MS	[1564]
3- <b>Co</b> (1,2- $RC_2B_9H_{11}$ )(1',2'- $C_2B_9H_{10}$ ) <sup>-</sup> R = C(O)C <sub>6</sub> H <sub>4</sub> - <i>o</i> -C(O)N(CH <sub>2</sub> ) <sub>3</sub> azetidin	S, X, H, B, C, MS	[1564]
3- <b>Co</b> (1,2- $C_2B_9H_{10}$ ) <sub>2</sub> -1,1'- <i>bicyclo</i> -COC(O)C <sub>6</sub> H <sub>4</sub> <sup>-</sup> benzofuran	S, X, H, B, C, MS	[1564]
3- <b>Co</b> [1,2-(H <sub>2</sub> NCH <sub>2</sub> ) <sub>2</sub> R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ](1',2'- $C_2B_9H_{11}$ ) <sup>-</sup>	S, X, H, B, C, MS	[1564]
(1,2- $C_2B_9H_{11}$ ) <b>Co</b> {1,2-[XO(CH <sub>2</sub> ) <sub>n</sub> ] $C_2B_9H_{10}$ } <sup>-</sup> n = 1-3 X = SO <sub>2</sub> Me, SO <sub>2</sub> C <sub>6</sub> H <sub>4</sub> Me	S, X(n = 2, X = SO <sub>2</sub> Me), H, B(2d), C	[1587]
<b>Co</b> {1,2-[XO(CH <sub>2</sub> ) <sub>n</sub> ] $C_2B_9H_{10}$ } <sub>2</sub> <sup>-</sup> n = 1, 2 X = SO <sub>2</sub> Me, SO <sub>2</sub> C <sub>6</sub> H <sub>4</sub> Me	S, H, B(2d), C	[1587]

Compound	Information	References
$(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{Co}[1,2\text{-}[\text{RR}'\text{N}(\text{CH}_2)_n]\text{C}_2\text{B}_9\text{H}_{10}]^-$ $n=1\text{-}3$ R=H, Et; R'=CH <sub>2</sub> Ph, <i>n</i> -C <sub>4</sub> H <sub>9</sub> , Et	S, H, B(2d), C	[1587]
$\text{Co}[1,2\text{-}[\text{H}_2\text{N}(\text{CH}_2)_2\text{C}_2\text{B}_9\text{H}_{10}]_2]^-$	S, H, B(2d), C	[1587]
$\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_{10})_2\text{-}1,1'\text{-CH}_2\text{NRCH}_2^-$ R=H, <i>n</i> -C <sub>4</sub> H <sub>9</sub>	S, H, B(2d), C	[1587]
<b><i>Closo-commo-(C<sub>2</sub>B<sub>9</sub>)Co(C<sub>2</sub>B<sub>8</sub>) clusters</i></b>		
$(\text{C}_2\text{B}_9\text{H}_{11})_2\text{Co}_x(\text{C}_2\text{B}_8\text{H}_{10})_{x-1}^{x-}$ $x=2\text{-}7$	Raman	[927]
$[(2,3\text{-C}_2\text{B}_8\text{H}_{10})\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_{11})]^-$	E	[309]
$[2,3\text{-C}_2\text{B}_8\text{H}_9\text{-}7\text{-C}_5\text{H}_5\text{N}]\text{Co}[1,2\text{-C}_2\text{B}_9\text{H}_{11}]$	S, H, B, IR, UV	[456]
$[2,3\text{-C}_2\text{B}_8\text{H}_{11}\text{-}11\text{-C}_5\text{H}_5\text{N}]\text{Co}[1,2\text{-C}_2\text{B}_9\text{H}_{11}]^-$	S, H, B, IR, UV	[456]
	X	[458]
$\text{Mn}[1,10\text{-C}_{12}\text{H}_8\text{N}_2]_3^{2+}$ $(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{Co}(1,2\text{-C}_2\text{B}_8\text{H}_{10})\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_{11})^{2-}$ · MeCN	S, X	[1043]
$(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{Co}(1,2\text{-C}_2\text{B}_8\text{H}_{10})\text{Co}(1,2\text{-C}_2\text{B}_8\text{H}_{10})\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_{11})^{3-}$	S, H, IR, UV, E	[924]
	X	[1044]
	B	[1045]
$[(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{Co}(1,2\text{-C}_2\text{B}_8\text{H}_{10})\text{Co}(2,n\text{-C}_2\text{B}_8\text{H}_{10})]^{2-}$ $n=3,4$	S, H, B, IR, UV	[460]
$(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{Co}(1,2\text{-C}_2\text{B}_8\text{H}_{10})\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_{11})^{2-}$	E	[872]
	MAG	[929]
	COND, aggregation in H <sub>2</sub> O; dynamic and static light-scattering; atomic force microscopy; scanning electron microscopy	[932]
$(1,2\text{-C}_2\text{B}_9\text{H}_{11})\text{Co}(1,2\text{-C}_2\text{B}_8\text{H}_{10})\text{Co}(1,2\text{-C}_2\text{B}_9\text{H}_{11})^{n-}$	MAG	[1048]
<b><i>Closo-Co<sub>2</sub>C<sub>2</sub>B<sub>8</sub> clusters</i></b>		
$2,n,1,7\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_{10})$ $n=4, 10$	S, H, B, IR, MS, E	[447]
$2,7,1,12\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_{10})$	S, H, B, IR, MS, UV, E	[307]
$2,8,1,12\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_{10})$	S, H, B, IR, MS, E	[447]
$2,9,1,12\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_{10})$	S, H, B, IR, MS, E	[447]
$3,4,1,2\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_{10})$	S, H, B, IR, MS, UV	[308]
$3,6,1,2\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_9)\text{-}n\text{-HgCl}$ $n=8, 9$	S, X( $n=8$ ), H, B	[1053]
	E ( $n=9$ )	[873]
$3,6,1,2\text{-CpCo}_2(\text{C}_2\text{B}_8\text{H}_6(\text{HgCl})_4)$	S, H, B	[1053]
$3,6,1,2\text{-Cp}_2\text{Co}_2(\text{RC}_2\text{B}_8\text{H}_9)$ R=HOCH <sub>2</sub> , CHO	S, IR	[1054]
$3,6,1,2\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_8)\text{-}8,9\text{-}[\text{OC}(\text{O})\text{CF}_3]_2$	S, H, B	[1053]
$3,6,1,2\text{-CpLCo}_2(\text{C}_2\text{B}_8\text{H}_{10})$ L= <i>n</i> -C <sub>4</sub> H <sub>9</sub> -C <sub>5</sub> H <sub>4</sub> , PhC <sub>5</sub> H <sub>4</sub>	S, H	[1055]
$3,6,1,7\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_{10})$	S, H, B, MS	[314]
$3,9,1,7\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_{10})$	S, H, B, MS	[314,447]
$6,9,1,7\text{-Cp}_2\text{Co}_2(\text{C}_2\text{B}_8\text{H}_{10})$	S, H, B, MS	[314,447]
<b><i>Closo-CoPC<sub>2</sub>B<sub>8</sub> clusters</i></b>		
$1,2,3,6\text{-}(\text{C}_4\text{Me}_4)\text{Co}(\text{PC}_2\text{B}_8\text{H}_9\text{-}5\text{-R})$ R=H, Cl	S, H, B, P, E(Cl)	[1056]
$1,2,4,8\text{-}(\text{C}_4\text{Me}_4)\text{Co}(\text{PC}_2\text{B}_8\text{H}_9\text{-}5\text{-Cl})$	S, X(H), H, B, P, E	[1056]
$1,2,3,10\text{-}(\text{C}_4\text{Me}_4)\text{Co}(\text{PC}_2\text{B}_8\text{H}_9\text{-}5\text{-Cl})$	S, H, B, P	[1056]

Continued

Compound	Information	References
<b>Closo-CoMC<sub>2</sub>B<sub>8</sub> clusters</b>		
(C <sub>8</sub> H <sub>12</sub> )PtCpCo(C <sub>2</sub> B <sub>8</sub> H <sub>10</sub> )	H	[1057]
<b>Nido-Co<sub>2</sub>C<sub>4</sub>B<sub>6</sub> clusters</b>		
Cp* <sub>2</sub> Co <sub>2</sub> (Et <sub>4</sub> C <sub>4</sub> B <sub>6</sub> H <sub>6</sub> )	S, H, B, IR, UV, MS	[129]
Cp* <sub>2</sub> Co <sub>2</sub> (Et <sub>4</sub> C <sub>4</sub> B <sub>6</sub> H <sub>6</sub> )	S	[31,162]
(NC <sub>4</sub> Me <sub>4</sub> ) <sub>2</sub> Co <sub>2</sub> (Et <sub>4</sub> C <sub>4</sub> B <sub>6</sub> H <sub>6</sub> )	S, H, B, IR, UV, MS	[129]
Cp* <sub>2</sub> Co <sub>2</sub> (C <sub>4</sub> B <sub>6</sub> H <sub>8</sub> Cl <sub>2</sub> )	S, H, B, IR, MS	[148]
<b>Rhodium</b>		
<b>Closo-RhCB<sub>10</sub> clusters</b>		
(Ph <sub>3</sub> P)XRh(Me <sub>3</sub> CNH <sub>2</sub> -CB <sub>10</sub> H <sub>10</sub> ) X=Br, Cl	S, X(Br), H, B, C, P	[861]
(Ph <sub>3</sub> P)X(Me <sub>3</sub> CNH <sub>2</sub> )Rh(Me <sub>3</sub> CH <sub>2</sub> N-CB <sub>10</sub> H <sub>10</sub> ) X=Br, Cl	S, H, B, C, P	[861]
(Ph <sub>3</sub> P)(Me <sub>3</sub> CNH <sub>2</sub> ) <sub>2</sub> Rh(Me <sub>3</sub> CH <sub>2</sub> N-CB <sub>10</sub> H <sub>10</sub> ) X=Br, Cl	S, X, H, B, C, P	[861]
Other derivatives of preceding		[861]
2,1-(Ph <sub>3</sub> P) <sub>2</sub> HRh(CB <sub>10</sub> H <sub>10</sub> -1-NH <sub>2</sub> )	S, H, P, IR	[1065]
2,1-Br(Ph <sub>3</sub> P)Rh[(NHRR')CB <sub>10</sub> H <sub>10</sub> ] R=H, CH <sub>2</sub> CH=CHMe; R'=CH <sub>2</sub> CH=CHMe	S, X, H, P, IR	[1066]
2,1-Cl(Ph <sub>3</sub> P)Rh[(Me <sub>3</sub> N)CB <sub>10</sub> H <sub>10</sub> ]	S, H, B, P	[1067]
2,1-(MeC <sub>6</sub> H <sub>5</sub> )Rh[(Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>10</sub> ]	S, H, B, C	[1069]
2,1-(1,3,5-C <sub>6</sub> H <sub>3</sub> Me <sub>3</sub> )Rh[(Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>10</sub> ]	S, X, H, B, C	[1069]
2,1-(C <sub>16</sub> H <sub>16</sub> )Rh[(Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>10</sub> ] C <sub>16</sub> H <sub>16</sub> =paracyclophane	S, X, H, B, C	[1069]
2,1-(CO) <sub>2</sub> Rh <sub>2</sub> {[(Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>10</sub> ]} <sub>2</sub>	S, X, H, B, C	[1069]
2,1-[HC(pyrazolyl)] <sub>3</sub> Rh[(Me <sub>3</sub> CNH)CB <sub>10</sub> H <sub>10</sub> ] <sup>+</sup>	S, X, H, B, C	[1069]
Related derivatives		[1069]
2,1-(Ph <sub>3</sub> P) <sub>2</sub> HRh[(Me <sub>3</sub> N)CB <sub>10</sub> H <sub>10</sub> ]	S, H, P, IR	[1070]
2,1-(Ph <sub>3</sub> P)XRh[(CH <sub>2</sub> =CH-CH <sub>2</sub> ) <sub>2</sub> N]CB <sub>10</sub> H <sub>10</sub> X=Cl, Br	S, H, C, P	[1071]
2,1-Cl(Ph <sub>3</sub> P)Rh[(Me <sub>2</sub> C=NH)CB <sub>10</sub> H <sub>10</sub> ]	S, X, H, B, C, P	[1072]
2,1-(Ph <sub>3</sub> P)Rh[(O=CRNH)CB <sub>10</sub> H <sub>10</sub> ] (Rh-O) R=Me, CMe=CH <sub>2</sub> , exo- CH=CHMe, Ph	S, X(CMe=CH <sub>2</sub> ), H, B, C, P	[1072]
Related derivatives		[1072]
<b>Closo-RhC<sub>2</sub>B<sub>9</sub> clusters</b>		
<b>Mono(dicarbollyl) Rh complexes</b>		
<b>CpRh(C<sub>2</sub>B<sub>9</sub>) and Cp*Rh(C<sub>2</sub>B<sub>9</sub>) complexes</b>		
3,1,2-Cp*Rh(LC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) L=PPh <sub>2</sub> , P(S)Ph <sub>2</sub>	S, X, H, B, P, IR	[1424]
[3,1,2-Cp*Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-PPh <sub>2</sub> Ag(OC <sub>4</sub> H <sub>8</sub> )] <sub>2</sub> (μ-SO <sub>3</sub> CF <sub>3</sub> ) <sub>2</sub>	S, X, H, B, P, IR	[1424]
3,1,2-(η <sup>5</sup> -C <sub>5</sub> R <sub>5</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+</sup> R=H, Me	S, X(Me), H, B, P	[1077]
<b>Other (hydrocarbon)Rh(C<sub>2</sub>B<sub>9</sub>) complexes</b>		
3,1,2-(η <sup>5</sup> -EtC <sub>5</sub> H <sub>4</sub> )Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, MS	[1080]
3,1,2-(η <sup>2</sup> ,η <sup>3</sup> -H <sub>2</sub> C=CH-C <sub>5</sub> H <sub>4</sub> )Rh(RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=H, Me; R'=H, Me, Ph vinylcyclopentadienyl	S, H, B, C	[1081]

Compound	Information	References
3,1,2-(C <sub>7</sub> H <sub>8</sub> )Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> C <sub>7</sub> H <sub>8</sub> = norbornadiene	S, H	[1080]
3,1,2-(C <sub>7</sub> H <sub>7</sub> CH <sub>2</sub> )Rh(RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R = H, Me, Ph; R' = H, Me C <sub>7</sub> H <sub>7</sub> CH <sub>2</sub> = 2-methylenenorbornadienyl	S, X(Me <sub>2</sub> ), H, C	[1082]
2,1,7-(C <sub>7</sub> H <sub>7</sub> CH <sub>2</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) C <sub>7</sub> H <sub>7</sub> CH <sub>2</sub> = 2-methylenenorbornadienyl	S, H, C, IR, MS	[1082]
"Pseudocloso"-3,1,2-(C <sub>7</sub> H <sub>7</sub> CH <sub>2</sub> )Rh[(p-MeC <sub>6</sub> H <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] C <sub>7</sub> H <sub>7</sub> = norbornadienyl	S, X	[1083]
3,1,2-(C <sub>7</sub> H <sub>7</sub> CH <sub>2</sub> )Rh[(PhCH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] C <sub>7</sub> H <sub>7</sub> = norbornadienyl	S, X, H	[1085]
3,1,2-(C <sub>7</sub> H <sub>7</sub> CH <sub>2</sub> )Rh(MePhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) C <sub>7</sub> H <sub>7</sub> = norbornadienyl	S, H	[1086]
3,1,2-(R'' <sub>2</sub> C <sub>8</sub> H <sub>n</sub> )Rh(RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) n = 9, 11; R, R' = o-(CH <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>4</sub> , Me; R'' = H, Me	S, H	[1086]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Rh(RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> RR' = Me <sub>2</sub> , H <sub>2</sub> , HPh	S, H	[1089]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )	S, X, H, B, IR	[1091]
3,1,2-(η <sup>3</sup> -C <sub>8</sub> H <sub>13</sub> )Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) C <sub>8</sub> H <sub>13</sub> = xylylene	S, H, B	[1078]
3,1,2-(η <sup>3</sup> -C <sub>8</sub> H <sub>13</sub> )Rh(MePhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	UV (detailed; photometric detection)	[1000]
3,1,2-(η <sup>3</sup> -C <sub>8</sub> H <sub>13</sub> )(CO)Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) (C <sub>8</sub> H <sub>13</sub> = xylylene)	S, H, C	[1087]
3,1,2-(η <sup>3</sup> -C <sub>8</sub> H <sub>13</sub> )Rh{[o-C <sub>6</sub> H <sub>4</sub> (CH <sub>2</sub> ) <sub>2</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> } 2 isomers C—H...M C <sub>8</sub> H <sub>13</sub> = xylylene	S, H(2d, variable temp), C(HETCOR), IR	[1095]
3,1,2-(η <sup>3</sup> -C <sub>8</sub> H <sub>13</sub> )Rh[(PhS)R'C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> D] R = Ph, Me	S, H, B, C, IR	[1096]
3,1,2-(C <sub>9</sub> H <sub>7</sub> )Rh(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) C <sub>9</sub> H <sub>7</sub> = indenyl	X	[1097]
3,1,2-(C <sub>9</sub> H <sub>7</sub> )Rh[(MeOCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] C <sub>9</sub> H <sub>7</sub> = indenyl	S, H, B(2d), IR	[1097]
3,1,2-(C <sub>9</sub> Me <sub>7</sub> )Rh(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) C <sub>9</sub> Me <sub>7</sub> = heptamethylindenyl	S, X, H, B, IR	[1098]
3,1,2-(C <sub>9</sub> H <sub>2</sub> Me <sub>5</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) pentamethylindenyl	S, X, H, B	[1099]
"Pseudocloso"-3,1,2-(C <sub>9</sub> Me <sub>7</sub> )Rh(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, H, B(2d), IR	[1098]
3,1,2-(η <sup>3</sup> -dicyclopentenyl)Rh(RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup> C—H...Rh R = H, Me; R' = H, Me, CH=CH <sub>2</sub> , CMe=CH <sub>2</sub> , CH <sub>2</sub> OH	S, X(H, CH=CH <sub>2</sub> ; H, CH <sub>2</sub> OH), H, C	[1100]
3,1,2-(η <sup>4</sup> -C <sub>10</sub> H <sub>12</sub> )Rh(HOCH <sub>2</sub> -C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup>	S, X, H	[1101]
3,1,2-CpM(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-4-SMe <sub>2</sub> <sup>+</sup> M = Rh, Ir	S, X(Rh), H, B	[1592]
3,1,2-CpM(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-4-SMe M = Rh, Ir	S, X(Rh), H, B	[1592]
<i>(borole)Rh(C<sub>2</sub>B<sub>9</sub>) complexes</i>		
3,1,2-Cp*Rh(C <sub>6</sub> H <sub>5</sub> )-(H <sub>4</sub> C <sub>4</sub> B)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>2+</sup>	S, B	[1102]
3,1,2-Cp*Ir(H <sub>4</sub> C <sub>4</sub> B-Ph)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>2+</sup>	S, X, B	[1102]
3,1,2-(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )Rh(H <sub>4</sub> C <sub>4</sub> B)-(C <sub>6</sub> H <sub>5</sub> )Ir(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>2+</sup>	S, H, B	[1102]
<i>(phosphino)Rh(C<sub>2</sub>B<sub>9</sub>) complexes</i>		
3,1,2-(Et <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-Rh(CODH)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> C—C (COD = cyclooctadiene)	S, X, H, B, P	[1105]
3,1,2-(Et <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -6,10-D <sub>2</sub> )	S, B	[1104]
3,1,2-(Et <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -3,4,7-D <sub>3</sub> )	S, B	[1104]
3,1,2-(Et <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>6</sub> -4,5,6,7,11-D <sub>5</sub> )	S, B	[1104]
3,1,2-(Et <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6-Ph)	S, B	[1104]
[3,1,2-(R <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> R = Et, Me <sub>2</sub> Ph	S, H, B, P	[1105]
3,1,2-(Et <sub>3</sub> P) <sub>2</sub> HRh[(μ-CH <sub>2</sub> ) <sub>3</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B, P, IR	[1107]

Continued



Compound	Information	References
3,1,2-(Et <sub>3</sub> P) <sub>2</sub> HRh[(μ-CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B, P, IR	[1107]
[2,1,8-(Et <sub>3</sub> P) <sub>2</sub> HRh(MePhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )] <sub>2</sub>	S, H, B, P, IR	[1107]
2,1,8-(Et <sub>3</sub> P) <sub>2</sub> HRh(MePhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	X	[1108]
3,1,2-(Et <sub>3</sub> P) <sub>2</sub> HRh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, P, IR	[1107]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> DRh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	IR	[1104,1113]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh(H <sub>2</sub> C <sub>2</sub> B <sub>9</sub> D <sub>9</sub> ) deuterium exchange	S, H, IR	[1114]
<i>d</i> -3,1,2-(Ph <sub>2</sub> HRh(MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ))	X, P, OR	[1112]
3,1,2/2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, X, P	[1115]
2,1,12-(Ph <sub>3</sub> P) <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, P	[1115]
[3,1,2-(Ph <sub>3</sub> P)Rh(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )] <sub>2</sub> Rh-Rh	S, X, B, P	[1105]
2,1,7-(PhMe <sub>2</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, P	[835]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> (HSO <sub>4</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, P, IR	[1117]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> (NO <sub>3</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, IR	[1118]
	P	[1117]
3,1,2-(Ph <sub>3</sub> P)[η <sup>2</sup> -SC(H)PPh <sub>3</sub> ]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, IR	[1118]
2,1,7-[η <sup>2</sup> -S(H)C=C(PPh <sub>3</sub> )S]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, IR	[1118]
(Ph <sub>3</sub> P)(C <sub>2</sub> H <sub>4</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> isomers R=CO, C <sub>2</sub> H <sub>4</sub>		[1115]
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> HRh(MePhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, H, B, P, IR	[1119]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=H, Me, <i>n</i> -C <sub>4</sub> H <sub>9</sub> , Ph hydrogenation catalyst	S, H, P	[835]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-7-NC <sub>5</sub> H <sub>5</sub>	S, H, P	[835]
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> HRh(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=H, Me, Ph	S, H, P	[835]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-4-polystyrylMe	XPR	[1111]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6-R) R=PhCONH, NMe <sub>2</sub> hydrosilylation catalyst	S	[1120]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-1-CH <sub>2</sub> -cyc/α-N <sub>3</sub> P <sub>3</sub> (C <sub>5</sub> H <sub>10</sub> N) <sub>4</sub> Me	S, H, P, IR	[619]
[3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-1-CH <sub>2</sub> PMeNP(C <sub>5</sub> H <sub>10</sub> N) <sub>2</sub> N] <sub>n</sub>	S, P, IR	[619]
3,1,2-(Ph <sub>3</sub> P)HRh(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) enantiomers	OR	[1121]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> ClRh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S	[1123]
2,1,7-(Ph <sub>3</sub> P)ClRh(RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=H, Me; R'=Ph, Me	S, H, P	[1112]
3,1,2-(Ph <sub>3</sub> P)Cl <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, P, IR	[1124]
3,1,2-(Ph <sub>3</sub> P)(CO)ClRh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, P, IR	[1113,1123]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> BrRh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, P, IR	[1113]
3,1,2-(Ph <sub>3</sub> P)Br <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> K <sup>+</sup> (18-crown-6)	S, H, B, P, IR	[1124]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> I <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, B, P, IR	[1124]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh[(μ-CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, P(variable temp), IR	[1107]
	X	[1108]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>+</sup> BPh <sub>4</sub> <sup>-</sup>	S, H, B, IR, UV, E	[827]

Compound	Information	References
2,1,7-(CO)(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-10-CHLC <sub>6</sub> H <sub>4</sub> Me L = PMe <sub>3</sub> , PEt <sub>3</sub> , PMe <sub>2</sub> Ph	S, X(PEt <sub>3</sub> ), H, B, C, P, IR	[1125]
Related derivatives		[1125]
3,1,2-(Ph <sub>3</sub> P)(CO)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-4-NC <sub>5</sub> H <sub>5</sub>	X	[1126]
3,1,2-H(PPh <sub>3</sub> P) <sub>2</sub> Rh{[PhNHC(O)] <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> } hydrosilylation catalyst	S, IR	[1127]
{3,1,2-H(PPh <sub>3</sub> P) <sub>2</sub> Rh{[HNPhNHC(O)](CO)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> }} <sub>n</sub> hydrosilylation catalyst	S, IR	[1127]
3,1,2-(Ph <sub>3</sub> P)(CO)Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, C, P, IR	[1128]
3,1,2-(Ph <sub>3</sub> P)(CO)HRh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, C, P, IR	[1128]
3,1,2-(NO)(Ph <sub>3</sub> P)Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, P, IR	[1092]
3,1,2-(Ph <sub>3</sub> P)[η <sup>2</sup> -C(C <sub>6</sub> H <sub>5</sub> )NOC(=O)]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, P, IR	[1129,1130]
3,1,2-(Ph <sub>3</sub> P)[η <sup>2</sup> -C( <i>m</i> -C <sub>6</sub> H <sub>4</sub> F)NOC(=O)]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, P, IR	[1129,1130]
3,1,2/2,1,12-(Ph <sub>3</sub> P){-CH <sub>2</sub> CH <sub>2</sub> C[O( <i>n</i> -C <sub>4</sub> H <sub>9</sub> )]=O-}Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P	[1131]
2,1,7-(Ph <sub>3</sub> P)[η <sup>2</sup> -C( <i>m</i> -C <sub>6</sub> H <sub>4</sub> F)NOC(=O)]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, X, H, P, IR	[1129]
3,1,2-Cl <sub>2</sub> (Ph <sub>3</sub> P)Rh(RC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -8-SR'R'') R', R'' = Me, Et, Ph	S, X(H, Me, Me), H, B, P, IR	[1090]
[3,1,2-(C <sub>6</sub> H <sub>6</sub> )(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub>	S, H, B, IR, UV	[827]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh[( <i>n</i> -C <sub>4</sub> H <sub>9</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] hydrogenation catalyst	S, H, B, P, IR	[1132]
3,1,2-[μ-1,3-(CH <sub>2</sub> =CHCH <sub>2</sub> CH <sub>2</sub> )](Ph <sub>3</sub> P) <sub>2</sub> HRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) hydrogenation catalyst	S, X, H, B, P, IR	[1132]
3,1,2-(Ph <sub>3</sub> P)HRh(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )-9,12-Br <sub>2</sub>	S, X	[1133]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> HRh[(μ-CH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> CH <sub>2</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, P, IR	[1107]
[1,2-μ-(3,4-CH <sub>2</sub> CH <sub>2</sub> CMe=CHCH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> )]-3,1,2-(Ph <sub>3</sub> P)HRh-(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X	[1134]
μ-HC(C <sub>6</sub> H <sub>4</sub> Me)-3,1,2-(Ph <sub>3</sub> P)(CO)Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[1135]
μ-HC(C <sub>6</sub> H <sub>4</sub> Me)-3,1,2-(C <sub>8</sub> H <sub>12</sub> )Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[1135]
3,1,2-(S <sub>2</sub> CH)(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	X	[1136]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> (HSO <sub>4</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, B, P, IR	[1113]
3,1,2-(Ph <sub>3</sub> P)[C(Ph)-C(Ph <sub>3</sub> P)-CH-C(Ph)]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, P, IR	[1113]
[3,1,2-(Ph <sub>3</sub> P)(μ-CN)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>4</sub>	S, X, H, B, P, IR	[1113]
3,1,2-[CH <sub>2</sub> CH <sub>2</sub> C(O)OC <sub>4</sub> H <sub>9</sub> ](Ph <sub>3</sub> P)Rh(MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, X, H, B, P	[1112]
3,1,2-[CH <sub>2</sub> CH(Ph)C(O)OEt](Ph <sub>3</sub> P)Rh(MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, H, P	[1112]
3,1,2-[CH <sub>2</sub> CH(Me)C(O)OMe](Ph <sub>3</sub> P)Rh(MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R = H, Me	S, H, P	[1112]
2,1,7-LL'ClRh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L = Ph <sub>3</sub> P; L' = CO, C <sub>5</sub> H <sub>5</sub> N, MeCN, Ph <sub>3</sub> P; LL' = (Ph <sub>2</sub> PCH <sub>2</sub> ) <sub>2</sub>	S, B, P	[1112]
3,1,2-(Ph <sub>3</sub> P)Me(CO) <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P, IR	[1137]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> Me(CO) <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, IR	[1137]
2,1,7/2,1,12-(Ph <sub>3</sub> P) <sub>2</sub> Me(CO) <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P	[1137]
2,1,7-(Ph <sub>3</sub> P)[MeCHO(O)CMe]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P, IR	[1137]
3,1,2-(Ph <sub>3</sub> P)(C <sub>3</sub> H <sub>5</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X	[1137]
3,1,2/2,1,7-L(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> L = CO, C <sub>2</sub> H <sub>4</sub>	S, H, B, P, IR	[1138]
2,1,12-L(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> L = CO, Ph <sub>3</sub> P	S, H, B, P, IR	[1138]

Continued

Compound	Information	References
3,1,2-Ph <sub>2</sub> P(C <sub>6</sub> H <sub>5</sub> -)(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	X	
3,1,2/2,1,7-(Ph <sub>3</sub> P)[MeC(O)Me]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, P	[1140]
3,1,2/2,1,7-(Ph <sub>3</sub> P)(CO)[C(O)Me]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, P, IR	[1140]
Related derivatives		[1140]
3,1,2/2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> MeRh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S	[1140]
[18-crown-6] <sup>+</sup> 3,1,2-(Ph <sub>3</sub> P)(PhCH <sub>2</sub> )BrRh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, P	[1140]
3,1,2-I <sub>2</sub> (PPh <sub>3</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )	S, X, H, B, P, IR	[1141]
3,1,2-(Ph <sub>2</sub> PCH <sub>2</sub> CH <sub>2</sub> PPh <sub>2</sub> )ClRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+</sup>	S, X, H, B, P	[1077]
1, <i>n</i> -C <sub>6</sub> H <sub>4</sub> [3,1,2-H(Ph <sub>3</sub> P) <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-CH <sub>2</sub> ) <sub>2</sub> <i>n</i> =3,4	S	[1142]
<i>Exo-metallated Rh(C<sub>2</sub>B<sub>9</sub>) complexes</i>		
(CO) <sub>2</sub> CpW(μ-CC <sub>6</sub> H <sub>4</sub> Me)Au-2,1,7-(CO)(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, C, P, IR	[1143]
(CO) <sub>2</sub> CpW(μ-CC <sub>6</sub> H <sub>4</sub> Me)Au <sub>2</sub> [3,1,2-(CO) <sub>2</sub> Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub>	S, X, H, C, IR	[1144]
μ-Mn(CO) <sub>4</sub> -3,1,2-L(CO)Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L=Ph <sub>3</sub> P, CO	S, H, B, C, P, IR	[1092]
μ-(Me <sub>4</sub> C <sub>4</sub> )(CO) <sub>2</sub> Co-3,1,2-(CO)(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, P, IR	[1145]
3,1,2-[η <sup>4</sup> -(Me <sub>3</sub> C) <sub>2</sub> P <sub>2</sub> -Co(CO) <sub>2</sub> (C <sub>4</sub> Me <sub>4</sub> )]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, P, IR	[1146]
μ-(Ph <sub>3</sub> P)(CO) <sub>2</sub> (H)Ir-3,1,2-(Ph <sub>3</sub> P)(CO)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, X, H, B, C, P, IR	[1145]
μ-(Ph <sub>3</sub> P)(CO) <sub>2</sub> Ir-3,1,2-(Ph <sub>3</sub> P)(CO)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup>	S, H, B, C, P, IR	[1145]
(Et <sub>3</sub> P) <sub>2</sub> Pt(μ-H)(μ-CO)-3,1,2-(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, P, Pt, IR, MS	[1147]
μ-LL'/Pt-2,1,7-(Ph <sub>3</sub> P)(CO)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) L=PEt <sub>3</sub> , PPh <sub>3</sub> , PMe <sub>2</sub> Ph; L'=PEt <sub>3</sub> , CO, PMe <sub>2</sub> Ph	S, H, B, C, P, IR	[1148]
μ-(PhC=CPh)(Et <sub>3</sub> P) <sub>2</sub> Pt-2,1,7-(Ph <sub>3</sub> P)(CO)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, H, B, C, P, IR	[1148]
Related 3,1,2- and 2,1,7- <i>exo</i> -metallated Pt derivatives		[1148]
3,1,2-(CO)(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-μ(H)-3,7-CuPPh <sub>3</sub>	S, H, B, P, IR	[1149]
(Ph <sub>3</sub> P)Cu-3,1,2-(L)(CO)Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L=CO, Ph <sub>3</sub> P, Me <sub>3</sub> P	S, H, B, C, P, IR	[1128]
3,1,2-[η <sup>4</sup> -(Me <sub>3</sub> C) <sub>2</sub> P <sub>2</sub> -Au(Ph <sub>3</sub> P)]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, P, IR	[1146]
(Ph <sub>3</sub> P) <sub>2</sub> (Ph <sub>3</sub> PAu)-3,1,2-Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, C, P	[1150]
(Ph <sub>3</sub> P) <sub>2</sub> (Ph <sub>3</sub> PAu)-3,1,2-Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, P	[1150,1151]
(Ph <sub>3</sub> P)(H)(Ph <sub>3</sub> PAu) <sub>2</sub> -3,1,2-Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, B, C, P	[1150,1151]
(Ph <sub>3</sub> P)Au(CO)-2,1,7-(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, C, P, IR	[1143]
(CH <sub>2</sub> ) <sub>n</sub> [-Ph <sub>2</sub> P-Au-(CO)-2,1,7-(Ph <sub>3</sub> P)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <i>n</i> =2-6	S, H, C, P, IR	[1143]
<i>Other (ligand)Rh(C<sub>2</sub>B<sub>9</sub>) complexes</i>		
[3,1,2-X <sub>2</sub> Rh(Me <sub>2</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> X=Cl, Br, I	S, H, B	[1152]
3,1,2-(CO) <sub>2</sub> Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, C, IR	[1144]
3,1,2-(CO) <sub>2</sub> Rh(PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )-7-SMe <sub>2</sub>	S, X, H, B, IR	[1153]
3,1,2-(CO) <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )-NC <sub>5</sub> H <sub>5</sub>	S, H, B, IR, MS	[1154]
3,1,2-I <sub>2</sub> (CO)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )	S, X, H, B, P, IR	[1141]
3,1,2/2,1,7-HB(pyrazolyl) <sub>3</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, IR, UV, MS	[1155]
3,1,2-(μ-Br) <sub>2</sub> Rh <sub>2</sub> [C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -3-SC(NHPh)N(Ph)CH <sub>2</sub> ] <sub>2</sub>	X	[1156]

Compound	Information	References
3,1,2-(MeCN)(PPh <sub>2</sub> Me) <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>+</sup> SbF <sub>6</sub> <sup>-</sup>	S, X, H, B, IR	[1157]
3,1,2-L(SCHNPh)Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=Ph <sub>3</sub> P, Cl	S, H, B, IR	[1157]
3,1,2-[η <sup>4</sup> -(Me <sub>3</sub> C) <sub>2</sub> C <sub>2</sub> P <sub>2</sub> ]Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, C, P, IR	[1146]
[2,1,8-(HO)Rh(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>4</sub>	S, X, H, B, IR	[1088]
3,1,2-[X <sub>2</sub> Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sub>2</sub> X=Cl, Br, I	S, H, B, P	[1077]
3,1,2-L(CO)Rh(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-4-CH <sub>2</sub> R <sup>-</sup> L=Ph <sub>3</sub> P, CO; R=H, Me	S, H, B, C, P, IR	[1158]
3,1,2-Cp*Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, E	[1500]
3,1,2-(Me <sub>3</sub> C <sub>3</sub> H)Rh(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R=Me, O(CH <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>4</sub> isomers Rh...H π-allyl agostic CH <sub>3</sub> ...Rh bonding catalysis of alkene hydroformylation in supercritical CO <sub>2</sub>	S, X, H(2d), B, C, IR	[1471]
3,1,2-CpRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-SMe <sub>2</sub> ) <sup>+</sup>	S, H	[1483]
3,1,2-CpRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-SMe)	S, X, H, B	[1483]
3,1,2-Br <sub>2</sub> (Me <sub>2</sub> S)Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-SMe <sub>2</sub> )	S, X(Cl,Br), H, B	[1492]
3,1,2-(η <sup>3</sup> -C <sub>6</sub> H <sub>11</sub> )Rh(MePhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) 2 diastereomers allyl	S, X, H, B, IR	[1498]
<b>Bis(dicarbollyl) Rh complexes</b>		
3,1,2-Rh(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup>	E	[1159]
3,1,2-(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Rh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-4-SMe <sub>2</sub>	S, H, B, P	[1077]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-SMe <sub>2</sub> )	S, H, B	[1492]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Rh(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-SMe <sub>2</sub> ), {1,2,4,12-XRh[(Me <sub>3</sub> C-NH)-C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ]} <sub>2</sub> (μ-I) <sub>2</sub> , 3,1,2-(C <sub>6</sub> H <sub>6</sub> )LRh(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SMe <sub>2</sub> ) <sup>2+</sup> , and 3,1,2-(C <sub>6</sub> H <sub>6</sub> )LRh(R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>+</sup> (R=H, Me) catalysts for oxidative coupling of benzoic acid with diphenylacetylene in the presence of copper(II) diacetate to form 1,2,3,4-tetraphenylnaphthalene		[1557]
<b>Closo-RhC<sub>3</sub>B<sub>8</sub> clusters</b>		
1,2,4,12-(C <sub>8</sub> H <sub>12</sub> )Rh[(12-Me <sub>3</sub> CNH)C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ]	S, H, B	[693]
1,3'-Rh[2,4,12-(Me <sub>3</sub> C-NH)C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ](1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B	[1432]
<b>Iridium</b>		
<b>Closo-IrC<sub>2</sub>B<sub>9</sub> clusters</b>		
<b>Mono(dicarbollyl) Ir complexes</b>		
<b>CpIr(C<sub>2</sub>B<sub>9</sub>) and Cp*Ir(C<sub>2</sub>B<sub>9</sub>) complexes</b>		
3,1,2-Cp*Ir(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, E	[1500]
2,1,7-(C <sub>8</sub> H <sub>12</sub> )Ir[(7-Me <sub>2</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B	[1562]
3,1,2-CpIr(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-SMe <sub>2</sub> ) <sup>+</sup>	S, H	[1483]
3,1,2-CpIr(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-SMe)	S, X, H, B	[1483]
Pseudocloso-3,1,2-X <sub>2</sub> Ir(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-SMe <sub>2</sub> ) X=Cl, Br, I	S, X, H, B	[1492]
3,1,2-Br <sub>2</sub> (Me <sub>2</sub> S)Ir(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-SMe <sub>2</sub> )	S, H, B	[1492]
<b>Other (hydrocarbon)Ir(C<sub>2</sub>B<sub>9</sub>) complexes</b>		
3,1,2-Lir(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SMe <sub>2</sub> ) <sup>2+</sup> L=C <sub>6</sub> H <sub>6</sub> , C <sub>6</sub> H <sub>5</sub> OMe, C <sub>6</sub> H <sub>3</sub> Me <sub>3</sub> , C <sub>6</sub> H <sub>2</sub> Me <sub>4</sub> , C <sub>6</sub> Me <sub>6</sub>	S, B, E	[1079]
3,1,2-(R'' <sub>2</sub> C <sub>8</sub> H <sub>11</sub> )Ir(RR'C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) R,R'=o-(CH <sub>2</sub> ) <sub>2</sub> C <sub>6</sub> H <sub>4</sub> , Me; R''=H, Me	S, H	[1086]
3,1,2-[η <sup>3</sup> -endo-1,5-cyclo-Me <sub>2</sub> C <sub>8</sub> H <sub>9</sub> ]Ir[[o-C <sub>6</sub> H <sub>4</sub> (CH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -8-R] R=H, OEt C—H...Ir	S, X, H(actual, var. T), B, C(actual), IR	[1161]

Continued

Compound	Information	References
3,1,2-( $\eta^3$ -C <sub>8</sub> H <sub>13</sub> )Ir[cyclo-(C <sub>8</sub> H <sub>12</sub> ) <sub>2</sub> Ir <sub>2</sub> E]C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> - $\mu$ (Ir, B)-OMe E = S, Se	S, X(Se), H, B, IR	[1162]
3,1,2-( $\eta^4$ -C <sub>8</sub> H <sub>12</sub> )Ir(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -9-SMe <sub>2</sub> )	S, H, B	[1164]
	X	[1165]
3,1,2-LIr(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>+</sup> L = C <sub>6</sub> H <sub>6</sub> , MeC <sub>6</sub> H <sub>5</sub> , 1,2/1,3-Me <sub>2</sub> C <sub>6</sub> H <sub>4</sub> , 1,2,4,5-Me <sub>4</sub> C <sub>6</sub> H <sub>2</sub> , 1,3,5-Me <sub>3</sub> C <sub>6</sub> H <sub>3</sub> , [2.2]paracyclophane catalysis of oxidative coupling of benzoic acid with PhC $\equiv$ CPh (L = C <sub>6</sub> H <sub>6</sub> )	S, X(Me <sub>4</sub> C <sub>6</sub> H <sub>2</sub> ), H, B	[1584]
3,1,2-(MeCN) <sub>3</sub> Ir(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>+</sup>	S, X, H, B	[1584]
3,1,2-(RC <sub>5</sub> H <sub>4</sub> )Ir(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R = H, C(O)Me	S, H	[1584]
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> Hir(RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R = H, Ph	S, H, P	[835]
3,1,2-(R <sub>3</sub> P) <sub>2</sub> Hir(RC <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R = Ph, <i>p</i> -C <sub>6</sub> H <sub>4</sub> Me	S	[1167]
(Ph <sub>3</sub> P)[ $\eta^2$ -C( <i>p</i> -C <sub>6</sub> H <sub>4</sub> Cl)NOC(=O)]-3,1,2-Ir(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, P, IR	[1129]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )Rh(H <sub>4</sub> C <sub>4</sub> B)-(C <sub>6</sub> H <sub>5</sub> )Ir(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>2+</sup>	S, H, B	[1102]
<i>Exo-metallated Ir(C<sub>2</sub>B<sub>9</sub>) complexes</i>		
3,1,2-(CO)(Ph <sub>3</sub> P)Ir(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )- $\mu$ (H)-3,7-CuPPh <sub>3</sub>	S, X, H, B, P, IR	[1149]
(CO) <sub>2</sub> CpW( $\mu$ -CC <sub>6</sub> H <sub>4</sub> Me)Au <sub>2</sub> [3,1,2-(CO) <sub>2</sub> Ir(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )] <sub>2</sub>	S, H, B, C, IR	[1170]
(Ph <sub>3</sub> PAu)(CO)-3,1,2-Lir(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L = PPh <sub>3</sub> , CO	S, X(Ph <sub>3</sub> P), H, B, C, P, IR	[1169]
(Ph <sub>3</sub> P) <sub>4</sub> Au <sub>4</sub> (CO)Ir-3,1,2-(CO) <sub>2</sub> Ir(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) <sup>+</sup> (Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, H, B, C, P, IR	[1169]
<b>Nickel</b>		
<i>Closo-NiCB<sub>10</sub> clusters</i>		
2,1-(RNC) <sub>2</sub> Ni[(RNC)CB <sub>10</sub> H <sub>10</sub> ] R = CMe <sub>3</sub> , xylyl	S, X(xylyl), H, B, C, IR	[1173]
2,1-(CO)Ni[(OEt <sub>2</sub> )(Ph <sub>2</sub> P- <i>o</i> -C <sub>6</sub> H <sub>4</sub> )CB <sub>10</sub> H <sub>9</sub> ] Ni-P	S, X, H, B, C, IR	[1173]
2,1-Ni(Me <sub>2</sub> HNCB <sub>10</sub> H <sub>10</sub> ) <sub>2</sub>	S, UV	[593]
2,1-Ni[(C <sub>3</sub> H <sub>7</sub> NHR)CB <sub>10</sub> H <sub>10</sub> ] <sub>2</sub> R = H, Me	S	[682]
2,1-Ni(RCB <sub>10</sub> H <sub>10</sub> ) <sub>2</sub> R = PhCH <sub>2</sub> NH <sub>2</sub> , HO	S, UV	[593]
2,1-(Me <sub>3</sub> CNC) <sub>2</sub> Ni(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, IR	[1176]
2,1-(NMe <sub>2</sub> CB <sub>10</sub> H <sub>10</sub> )Ni(Me <sub>2</sub> NHCB <sub>10</sub> H <sub>10</sub> ) <sup>-</sup>		[593]
<i>Closo-NiPCB<sub>9</sub> clusters</i>		
2,1,7-Ni(MePCB <sub>9</sub> H <sub>10</sub> ) <sub>2</sub>	S	[779]
2,1,7-(Ph <sub>3</sub> P)BrNi(MePCB <sub>9</sub> H <sub>10</sub> )	S	[1073]
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> Ni(MePCB <sub>9</sub> H <sub>10</sub> )	S	[1073]
<i>Closo-NiC<sub>2</sub>B<sub>9</sub> clusters</i>		
<i>Other (hydrocarbon)Ni(C<sub>2</sub>B<sub>9</sub>) complexes</i>		
4,1,6-( $\eta^3$ -C <sub>3</sub> H <sub>5</sub> )Ni(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup>	S, H, B, C	[1181]
2,1,7-( $\eta^3$ -C <sub>3</sub> H <sub>5</sub> )Ni(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )-2,6,11-Cu(PPh <sub>3</sub> )-6,11- $\mu$ (H) <sub>2</sub>	S, X, H, B, C, P	[1181]
2,1,7-( $\eta^3$ -RC <sub>3</sub> H <sub>4</sub> )Ni(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) R = H, Ph	S, H, B, C, IR	[1182]
2,1,7-[Ni(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )] <sub>2</sub> ( $\eta^4$ , $\eta^4$ -C <sub>8</sub> H <sub>8</sub> )	S, X, H, B, C, IR	[1182]
3-Ni <sup>III</sup> [1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-C $\equiv$ C-C <sub>6</sub> H <sub>4</sub> -C $\equiv$ C-C <sub>6</sub> H <sub>4</sub> -BODIPY]-[1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-C $\equiv$ C-C <sub>6</sub> H <sub>4</sub> -C $\equiv$ C-C <sub>6</sub> H <sub>4</sub> -CH <sub>2</sub> OC(O)-[NHC-(O)OCMe <sub>3</sub> ]-CH <sub>2</sub> -tryptophan] <sup>-</sup>	S, H, B, C, F, IR, MS, UV, fluorescence	[1591]

Compound	Information	References
3-Ni <sup>IV</sup> [1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-C≡C-C <sub>6</sub> H <sub>4</sub> -C≡C-C <sub>6</sub> H <sub>4</sub> <sup>-</sup> BODIPY]-[1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8'-C≡C-C <sub>6</sub> H <sub>4</sub> -C≡C-C <sub>6</sub> H <sub>4</sub> <sup>-</sup> CH <sub>2</sub> O(CO)[NHC-(O)OCMe <sub>3</sub> ]-tryptophan]	S, H, B, C, F, IR, MS, UV, fluorescence	[1591]
<b>(phosphino)Ni(C<sub>2</sub>B<sub>9</sub>) complexes</b>		
2,1,7-(Ph <sub>3</sub> P) <sub>2</sub> Ni(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, IR	[1057]
3,1,2-(Ph <sub>3</sub> P)HNi(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-PPh <sub>3</sub> )	S, H, B, IR, MS	[1184,1185]
3,1,2-(Ph <sub>3</sub> P)ClNi(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-PPh <sub>3</sub> )	S, B	[1184,1185]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>2</sub> PPh <sub>2</sub> ] <sub>2</sub> Ni(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, P	[1185]
3,1,2-[( <i>p</i> -MeC <sub>6</sub> H <sub>4</sub> ) <sub>3</sub> P]HNi(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-P( <i>p</i> -MeC <sub>6</sub> H <sub>4</sub> ) <sub>3</sub> )	S, H, B, P	[1185]
3,1,2-(Ph <sub>3</sub> P)(C <sub>5</sub> H <sub>5</sub> N)Ni(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P	[1185]
3,1,2-LNi[(Me <sub>2</sub> NCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] N → Ni L = (Ph <sub>2</sub> P) <sub>2</sub> (CH <sub>2</sub> ) <sub>2</sub> , <i>o</i> -phenanthroline]	S, X( <i>o</i> -phenanthroline), H, B, C, P, IR	[1187]
3,1,2/4,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>2</sub> PPh <sub>2</sub> ] <sub>2</sub> Ni(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -6-Et)	S, X, H, B	[1188]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> Ni[Me( <i>n</i> -C <sub>5</sub> H <sub>11</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B, C, IR	[1189]
3,1,2-(Me <sub>3</sub> CNC)INi[(Ph <sub>3</sub> PCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] L = Et <sub>3</sub> P, Me <sub>3</sub> CNC	S, X, H, B, C, P, IR	[1187]
<b>Other (ligandNi(C<sub>2</sub>B<sub>9</sub>) complexes</b>		
2,1,7-L <sub>2</sub> Ni(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L = CO, CCMe <sub>3</sub>	S, H, B, C, IR	[1182]
(μ-CO)(CO) <sub>5</sub> Co <sub>2</sub> -3,1,2-Ni(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C	[1191]
3,1,2/2,1,7-(bipyridyl)Ni(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, IR, MS	[1192–1194]
	H, B	[1194]
	X	[1195]
3,1,2-(NC <sub>5</sub> H <sub>4</sub> ) <sub>2</sub> Ni(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, C, IR, MS	[1196]
<b>Neutral bis(dicarbollyl) Ni complexes</b>		
3-Ni(1,2-D <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub>	S, IR, MS	[1199]
3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> ·L L = naphthalene, pyrene, PhNMe <sub>2</sub>	X	[1201]
3-Ni(1,2-Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub>	S, H, MS, UV, E	[1199]
3-Ni(1,2-MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> dd/II and meso isomers	S, X, H(var. temp), B(var. temp), C	[1200]
3-Ni(1,2-MeC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> <sup>-</sup> MeNC <sub>5</sub> H <sub>5</sub> <sup>+</sup> , NMe <sub>4</sub> <sup>+</sup> salts dd/II and meso isomers	S, X, H(var. temp), B(var. temp), C	[1200]
2-Ni(1,7-Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub>	S, H, B, IR, MS, UV, E	[1199]
3-Ni[1,2-(CH <sub>2</sub> ) <sub>3</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub>	S, H, IR, E	[968,1199]
3-Ni[(1,2- <i>m/p</i> -FC <sub>6</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub>	F	[761]
3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-OMe) <sub>2</sub>	X	[1203]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Ni(1,7-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, E	[1199]
[1,2-(CH <sub>2</sub> ) <sub>3</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]Ni[1,7-(CH <sub>2</sub> ) <sub>3</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, B, IR, MS, E	[968,1199]
2-Ni(1,7-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(8,8')-N <sub>2</sub> C <sub>3</sub> H <sub>3</sub>	S, X(meso, dl), H, B, IR, MS, E	[775]
3,1,2-Ph <sub>2</sub> B(pyrazolyl) <sub>2</sub> Ni(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	S, X, H, B, IR, UV, MS	[1155]
1,2-[(Me <sub>2</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> Ni <sub>2</sub> (C <sub>5</sub> H <sub>5</sub> ) <sup>+</sup> C <sub>5</sub> H <sub>5</sub> -bridged triple-decker sandwich	S, H, B	[806]

Continued

Compound	Information	References
<b>Anionic bis(dicarbollyl) Ni complexes</b>		
3-Ni <sup>III</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> ↔ Ni <sup>IV</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>0</sup> redox couple shuttle for dye-sensitized solar cells	E, UV	[1415]
X <sup>+</sup> 3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> X=Ph <sub>4</sub> P, H(phen), Cs, NMe <sub>4</sub>	ESR: influence of temp and X	[1412]
3-Ni(1,2-D <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> <sup>-</sup>	S, IR	[1199]
2-Ni(1,7-Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> <sup>-</sup>	S, IR, UV, E, OR	[1199]
3-Ni[1,2-(CH <sub>2</sub> ) <sub>3</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> <sup>-</sup>	S, IR, E	[968,1199]
[3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> ] <sub>2</sub> Ni <sup>2+</sup> ·4L L=2,2-bipyridine, pyridine	S, IR	[1194]
3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> Na <sup>+</sup> ·2L L=2,2-bipyridine	S, IR	[1194]
M <sup>+</sup> 3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> M=Cs, Me <sub>4</sub> N, Mn( <i>o</i> -phenanthroline) <sub>3</sub>	S, X[Mn( <i>o</i> -phenanthroline) <sub>3</sub> ], ESR (variable temp)	[1209]
Cp* <sub>2</sub> Fe <sup>+</sup> 3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup>	S, IR, MS, MAG	[601]
(1,10-phenanthroline) <sub>3</sub> M <sup>2+</sup> 3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> M=Mn, Ni, Cu	MAG (variable T)	[950]
3-[Ni <sup>II</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>2-</sup> ] <sub>2</sub> Ni <sup>4+</sup> ·4L L=1,10-phenanthroline	S	[1210]
X <sup>+</sup> 3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> X=Cs, Me <sub>4</sub> N, Ph <sub>4</sub> P, 0.5 Fe(phen) <sub>3</sub>	S, IR, UV, MAG, Raman, XPS	[1212]
3-Ni[1,2-(C <sub>4</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> <sup>-</sup> benzodicarbollide complex	S, H, IR, UV	[687]
3-Ni[1,2-(C <sub>4</sub> H <sub>4</sub> )C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ] <sub>2</sub> -μ(1,1')-(CH <sub>2</sub> ) <sub>4</sub> <sup>-</sup>	S, X, H, B, C, IR	[598]
3-Ni(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-Cl) <sub>2</sub> <sup>-</sup> extraction agent for Cs <sup>+</sup> and Sr <sup>2+</sup>	S	[769]
(1,2-Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )Ni(1,7-Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sup>-</sup>	S, IR, UV, E, OR	[1199]
3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -10-I) <sub>2</sub> <sup>-</sup>	S, X, H, B, C, IR, MS	[1459]
C(NHPh) <sub>3</sub> <sup>+</sup> 3-Ni(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> triphenylguanidinium	S, X	[1495]
3-Ni <sup>III</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6'-(CH <sub>2</sub> ) <sub>5</sub> C <sub>16</sub> H <sub>9</sub> ) <sup>-</sup> C <sub>16</sub> H <sub>9</sub> =pyrene templates for electro- or photocontrolled molecular motors	S, H, B, C, MS, UV, fluorescence, E	[1541]
3-Ni <sup>IV</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6'-(CH <sub>2</sub> ) <sub>5</sub> C <sub>16</sub> H <sub>9</sub> ) C <sub>16</sub> H <sub>9</sub> =pyrene templates for electro- or photocontrolled molecular motors	S, H, B, C, MS, UV, fluorescence, E	[1541]
<b>Palladium</b>		
<b>Closo-PdCB<sub>10</sub> clusters</b>		
2,1-(Me <sub>3</sub> CNC) <sub>2</sub> Pd(CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup>	S, H, IR	[1176]
2,1-(Ph <sub>2</sub> PCH <sub>2</sub> ) <sub>2</sub> Pd[(Me <sub>3</sub> N)CB <sub>10</sub> H <sub>10</sub> ]	S	[1176]
<b>Mono(dicarbollyl) Pd complexes</b>		
3,1,2-(C <sub>4</sub> Ph <sub>4</sub> )Pd(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, IR, UV	[618]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Pd(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, H, B, IR	[1217]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Pd(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, H, B, IR	[1217]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Pd(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+</sup>	S, X, H, B, IR	[1091]
3,1,2-(η <sup>2</sup> , σ-5-OMeC <sub>8</sub> H <sub>12</sub> )Pd(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )	S, X, H	[1091]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Pd[(C <sub>4</sub> H <sub>3</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, C, P, MS	[1218]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Pd[(MeOCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	S, H, B(2d), IR	[545]
3,1,2-L <sub>2</sub> Pd(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L=1,5-C <sub>8</sub> H <sub>12</sub> , Me <sub>3</sub> CNC, NH <sub>3</sub> , C <sub>4</sub> Ph <sub>4</sub>	S, B	[1219]
3,1,2-L <sub>2</sub> Pd(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) L <sub>2</sub> =(NMe <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> H <sub>4</sub> , (Ph <sub>2</sub> PCH <sub>2</sub> ) <sub>2</sub> ; L=Me <sub>3</sub> P, (MeO) <sub>3</sub> P	S, B	[1219]

Compound	Information	References
3,1,2-(Me <sub>3</sub> CNC)Pd(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[1180]
2,1,7-(Me <sub>3</sub> CNC)Pd(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, IR	[1057]
3,1,2-[C <sub>2</sub> H <sub>4</sub> (NMe <sub>2</sub> ) <sub>2</sub> ]Pd(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X	[1220]
3,1,2-(Me <sub>2</sub> NCH <sub>2</sub> CH <sub>2</sub> NMe <sub>2</sub> )Pd(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, B	[1219]
3,1,2-(Ph <sub>3</sub> P)(I)Pd(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )	X	[1221]
3,1,2-(PhMe <sub>2</sub> P)ClPd(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )	X	[1221]
3,1,2-(PMe <sub>2</sub> Ph) <sub>2</sub> Pd(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P, IR	[1222]
3,1,2-(PMe <sub>2</sub> Ph) <sub>2</sub> Pd(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )-4,5-(CN) <sub>2</sub>	S, X, H, B, P, IR	[1222]
3,1,2-(PMe <sub>2</sub> Ph) <sub>2</sub> Pd(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, C, IR	[1180]
3,1,2-(PMe <sub>2</sub> Ph) <sub>2</sub> Pd[(C <sub>4</sub> H <sub>2</sub> RS)C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -8-PMe <sub>2</sub> Ph] <sup>+</sup> (C <sub>4</sub> H <sub>2</sub> RS)-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> <sup>-</sup> R=H, Me	S, X(H), H, B, P, MS	[1223]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub>2</sub> PPh <sub>2</sub> ]Pd[(C <sub>4</sub> H <sub>3</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, P, MS	[1218]
3,1,2-[Me <sub>3</sub> C <sub>3</sub> P] <sub>2</sub> Pd{[μ-S(CH <sub>2</sub> ) <sub>3</sub> ]C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> }	S, H, B, P	[1224]
<b>Bis(dicarbollyl) Pd complexes</b>		
3-Pd(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> ·I	S, IR	[1199]
<b>Platinum</b>		
<b>Closo-PtCB<sub>10</sub> clusters</b>		
2,1-(Et <sub>3</sub> P) <sub>2</sub> (PhHg)Pt(CB <sub>10</sub> H <sub>11</sub> )	S, X, H, B, C, P	[1226]
2,1-(Et <sub>3</sub> P) <sub>2</sub> (Ph <sub>3</sub> PCu)Pt(CB <sub>10</sub> H <sub>11</sub> )	S, X, H, B, C, P	[1226]
2,1-(Et <sub>3</sub> P) <sub>2</sub> HPT(CB <sub>10</sub> H <sub>11</sub> )	S, X, H, B, C, P	[1226]
2,1-(Et <sub>3</sub> P) <sub>2</sub> HPT(CB <sub>10</sub> H <sub>11</sub> )	S, X, H, B, C, P	[1226]
2,1-(Et <sub>3</sub> P) <sub>2</sub> Pt[(Me <sub>3</sub> N)CB <sub>10</sub> H <sub>10</sub> ]	S, H, B, P, IR	[1176]
2,1-Cl(PMe <sub>2</sub> Ph) <sub>2</sub> Pt(CB <sub>10</sub> H <sub>11</sub> )	S, X, H, B, C, P	[1227]
[2,1-(PMe <sub>2</sub> Ph) <sub>2</sub> Pt(CB <sub>10</sub> H <sub>10</sub> ) <sub>2</sub>	S, X, H, B, C, P	[1227]
2,1-(PhSe)(Et <sub>3</sub> P)Pt(CB <sub>10</sub> H <sub>10</sub> -n-SePh) n=3, 7	S, X(n=3), H, B, C, P	[1228]
2,1-(PhSe)(Et <sub>3</sub> P)Pt(CB <sub>10</sub> H <sub>10</sub> )-3-O(CH <sub>2</sub> ) <sub>4</sub> Cl	S, X, H, B, C, P	[1228]
2,1-(Et <sub>3</sub> P) <sub>2</sub> Pt(CB <sub>10</sub> H <sub>10</sub> )-7-Te(Ph)CH <sub>2</sub> Cl	S, X, H, B, C, P	[1228]
exo-(PhTe)(Et <sub>3</sub> P)Pt(μ-PhTe) <sub>2</sub> -7,1-(Et <sub>3</sub> P)Pt(CB <sub>10</sub> H <sub>11</sub> )	S, X, H, B, C, P	[1228]
<b>Closo-PtC<sub>2</sub>B<sub>9</sub> clusters</b>		
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Pt(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, H, B, IR	[1217]
3,1,11-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Pt(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, H, B, IR	[1217]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> )Pt[(C <sub>4</sub> H <sub>3</sub> S)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, C, P, MS	[1218]
3,1,2-(R <sub>3</sub> P) <sub>2</sub> Pt(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) R=Me, Ph	S, IR, H, E	[1229]
3,1,2-[(n-C <sub>3</sub> H <sub>7</sub> ) <sub>3</sub> P] <sub>2</sub> Pt(PhC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )	S, IR, H, E	[1229]
3,1,2-(Et <sub>3</sub> P) <sub>2</sub> Pt(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X	[1215]
3,1,2-(Ph <sub>3</sub> P) <sub>2</sub> Pt(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, B	[1231]
3,1,11-(Me <sub>2</sub> Ph) <sub>2</sub> Pt(PhC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> -11-R) R=H, Ph	S, X, P	[1232]
2,1,7-(PR <sub>3</sub> ) <sub>2</sub> Pt(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -7-OEt) R <sub>3</sub> =Me <sub>2</sub> Ph, Et <sub>3</sub>	S, H, B, P	[1234]
2,1,7-(PhMe <sub>2</sub> P) <sub>2</sub> Pt(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X	[1215,1235]
2,1,7-L <sub>2</sub> Pt(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L=PMe <sub>2</sub> Ph, PEt <sub>3</sub> , PMe <sub>3</sub>	S, H, IR	[1057]
3,1,2-L <sub>2</sub> Pt(PhMeC <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) L=PMe <sub>2</sub> Ph, PEt <sub>3</sub> , PPh <sub>3</sub> , P(C <sub>6</sub> H <sub>4</sub> Me) <sub>3</sub>	S, X(PEt <sub>3</sub> , PPh <sub>3</sub> ), H, B, IR	[1236]

Continued



Compound	Information	References
3,1,2/2,1,8-(PMe <sub>2</sub> Ph) <sub>2</sub> Pt(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, H, B, P	[1237]
	H(variable temp), P	[1238]
2,1,8-(PMe <sub>2</sub> Ph) <sub>2</sub> Pt(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-4-Et	S, X, H, B, P, IR	[1239]
2,1,8-(PMe <sub>2</sub> Ph) <sub>2</sub> Pt(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> )-4-F	S, X, H, B, F, P, IR	[1239]
3,1,2-[CH-CPh-CH-CPh]Pt <sub>2</sub> (Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	S, X, H, B, C	[1191]
2,1,8-(Ph <sub>2</sub> PCH <sub>2</sub> CH <sub>2</sub> PPh <sub>2</sub> )Pt(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> - <i>n</i> -I) <i>n</i> =7, 10, 12	S, X, H, B, P, IR	[1186]
3,1,2-(Ph <sub>2</sub> PCH <sub>2</sub> ) <sub>2</sub> Pt(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, B	[1219,1240]
	X	[1240]
3,1,2-(Me <sub>2</sub> PhP) <sub>2</sub> Pt[(C <sub>4</sub> H <sub>3</sub> S) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ]	S, X, H, B, P, MS	[1218]
<b>Closo-Cu<sub>2</sub>B<sub>9</sub> clusters</b>		
3,1,2-(Ph <sub>3</sub> P)Cu(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-NC <sub>5</sub> H <sub>5</sub>	S, H, B	[1246]
3,1,2-(Ph <sub>3</sub> P)Cu(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-4-NC <sub>5</sub> H <sub>4</sub> CO(O)Me	S, X, H, B, P, IR	[1245]
3,1,2-(Ph <sub>3</sub> P)Cu(C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )- <i>exo</i> -4,8-(μ-H) <sub>2</sub> Cu(PPh <sub>3</sub> )	S, X, H, B, P, IR	[1245]
3,1,2-(Ph <sub>3</sub> P)Cu(Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )-μ <sub>3</sub> -B(8,9,12)-Cu(PPh <sub>3</sub> )	S, X, H, B, P	[1247]
3,1,2-(Ph <sub>3</sub> P)Cu(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )-4-SMe <sub>2</sub>	S, X, H, B	[1248]
<i>nido</i> -3-[( <i>o</i> -toluyl) <sub>3</sub> ]P(Cu(1,2-Ph <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )-μ <sub>3</sub> -B(8,9,12)-Cu[P( <i>o</i> -toluyl) <sub>3</sub> ])	S, X, H	[1247]
<b>Gold<sup>c</sup></b>		
<b>Nido-Au<sub>2</sub>B<sub>9</sub> clusters</b>		
3,1,2-(Ph <sub>3</sub> P)Au(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -NC <sub>5</sub> H <sub>5</sub> )	S, B, IR	[1246]
3,1,2-(Et <sub>2</sub> NCS <sub>2</sub> )Au(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	S, X, H, B, IR, MS	[1231,1250]
10,7,8-(μ-PtL <sub>2</sub> R)(μ-H)(Ph <sub>3</sub> P)Au(R' <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> ) L = PEt <sub>3</sub> , PMe <sub>2</sub> Ph; R = H, Me; R' = H, Me	S, X(H, Me, PEt <sub>3</sub> ), B, C, P	[1253]
<b>Theoretical Studies</b>		
<b>Molecular and electronic structure calculations</b>		
3-Cr(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup>	MNDO	[30]
3,1,2-RHC≡C=(PH <sub>3</sub> ) <sub>2</sub> Mn(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	DFT: better electron reservoir than Cp <sup>-</sup> analogue	[1256]
CpM(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) M = Mn, Re	DFT, ab initio, stable isomers	[1394]
3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>q</sup> <i>q</i> =0, -1; 3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ); 3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -I) <sup>-</sup>	DFT: effect of <i>iodo</i> -substitution on bond lengths and hyperpolarizability; redox switchable 2nd order NLO	[1487]
3,1,2-(CO) <sub>3</sub> Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	DFT	[706]
3,1,2-(CO) <sub>2</sub> (NO)Re(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	DFT	[713]
1,2, <i>n</i> -CpFe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup> <i>n</i> =3,4,8,9,12	Stability; <i>o</i> , <i>m</i> , <i>p</i> -directing influence; DFT	[1257]
1,7,8,12-CpFe(C <sub>3</sub> B <sub>8</sub> H <sub>11</sub> )	Stability; <i>o</i> , <i>m</i> , <i>p</i> -directing influence; DFT	[1257]
1,2,3,4/1,2,3,5/1,2,3,6/1,2,4,5/1,2,4,10-CpFe(PC <sub>2</sub> B <sub>8</sub> H <sub>10</sub> )	Stability; <i>o</i> , <i>m</i> , <i>p</i> -directing influence; DFT	[1257]
	D-D transitions	[764]
	DFT( <sup>1</sup> B), orbital analysis	[1408]
(TMTSF) <sup>+</sup> 3-Fe(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> TMTSF = tetramethyltetraselenafulvalenium	S, X, COND	[936]
3-Fe(1,2-RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> <sup>-</sup> R = H, Me, Ph	Effective magnetic moments	[1262]

Compound	Information	References
3,1,2-CpFeC <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	Effective magnetic moments	[1262]
	Electron transfer	[1263]
3,1,2-(1',3',5'-C <sub>6</sub> H <sub>3</sub> Me <sub>3</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	DFT, GIAO ( <sup>11</sup> B)	[1259]
1,2,3, <i>n</i> -CpFe(C <sub>3</sub> B <sub>8</sub> H <sub>11</sub> ) <i>n</i> =4,5	DFT	[785]
1,2,4,12-CpFe[(12-Me <sub>3</sub> CNH)C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ]	DFT	[423]
3,1,2-Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SMe <sub>2</sub> ) <sub>2</sub>	Dicarbollide rotational energy	[772]
3,1,2-(C <sub>6</sub> H <sub>6</sub> )Fe(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+</sup>	DFT: Fe bonding	[1530]
1,2,4,12-(C <sub>6</sub> H <sub>6</sub> )Fe{[12-(Me <sub>3</sub> C)HN]C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> }	DFT: Fe bonding	[1530]
3,1,2-Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SMe <sub>2</sub> ) <sub>2</sub>	Dicarbollide rotational energy	[772]
3,1,2-(MeC≡CPh)(CO) <sub>2</sub> Ru(Me <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> )	ZINDO	[831]
1,2,4,12-Cp*Ru[(12-Me <sub>3</sub> CNH)C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> ]	DFT	[423]
3,1,2-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> )Ru(RC <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -8-SMe <sub>2</sub> ) <sup>+</sup>	Antipodal effects	[819]
3,1,2-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> )Ru(RC <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -8-SMe)	Antipodal effects	[819]
3,1,2-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> )Ru(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SH) C—H...S—H...H—B interactions; 2-D polymeric network	DFT	[814]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub><i>n</i></sub> PPh <sub>2</sub> ]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <i>n</i> =2,5 influence of phosphino ligands on reactivity	DFT: energies, geometry	[1532]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub><i>n</i></sub> PPhC <sub>6</sub> H <sub>4</sub> ]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <i>n</i> =2,5 C <sub>6</sub> H <sub>4</sub> —B influence of phosphino ligands on reactivity	DFT: energies, geometry	[1532]
3,1,2-[Ph <sub>2</sub> P(CH <sub>2</sub> ) <sub><i>n</i></sub> P(C <sub>6</sub> H <sub>4</sub> ) <sub>2</sub> ]ClRu(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <i>n</i> =2,5 2 C <sub>6</sub> H <sub>4</sub> —B influence of phosphino ligands on reactivity	DFT: energies, geometry	[1532]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(1,1')-SiMe <sub>2</sub> <sup>-</sup>	DFT	[978]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -μ(1,1')-SiMeH <sup>-</sup>	DFT	[978]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> -μ(1,1')-SiMeR-μ(8,8')-C <sub>6</sub> H <sub>4</sub> <sup>-</sup>	DFT	[978]
(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )Co[1,2-(Me <sub>3</sub> Si)C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]-μ(8,8')-C <sub>6</sub> H <sub>4</sub> <sup>-</sup>	DFT	[978]
3-Co[1,2-(Me <sub>3</sub> Si)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sub>2</sub> <sup>-</sup>	DFT	[978]
3-(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-R) <sup>-</sup> (R=I, Me, Et, Ph, C <sub>6</sub> H <sub>4</sub> Ph, C <sub>6</sub> H <sub>4</sub> -4- <i>n</i> -C <sub>4</sub> H <sub>9</sub> , CH <sub>2</sub> CH <sub>2</sub> Ph)	Ab initio	[1015]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -9,12-I <sub>2</sub> ) <sub>2</sub> <sup>-</sup> R=Me, Ph air-stable redox couple	DFT: redox potentials)	[1573]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4,9,12-I <sub>3</sub> ) <sub>2</sub> <sup>-</sup> air-stable redox couple	DFT: redox potentials)	[1573]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -8,9,10,12-Me <sub>4</sub> ) <sub>2</sub> <sup>-</sup> air-stable redox couple	DFT: redox potentials)	[1573]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>7</sub> -8,9,10,12-I <sub>4</sub> ) <sub>2</sub> <sup>-</sup> air-stable redox couple	DFT: redox potentials)	[1573]
3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )[(Me <sub>2</sub> HSi)C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ] <sup>-</sup>	DFT: Si—H...H—C bonding; loss of H <sub>2</sub> and formation of 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sub>2</sub> -u(1,1')-SiMe <sub>2</sub> <sup>-</sup>	[1268]
3,1,2-(η <sup>5</sup> -indenyl)Co(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )	Extended Hückel, ligand conformation	[867]
3,1,2-(NC <sub>4</sub> H <sub>4</sub> )Co[(MeS) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	Effect of electron density transfer on carborane C—C distance	[903]
3,1,2-Co(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SMe <sub>2</sub> ) <sub>2</sub> <sup>+</sup>	Dicarbollide rotational energy	[772]
3,1,2-CpCo(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SH) C—H...S—H...H—B interactions; 2-D polymeric network	DFT	[814]
Cp <sub>2</sub> Co <sub>2</sub> (Me <sub>4</sub> C <sub>4</sub> B <sub>6</sub> H <sub>6</sub> )	Polyhedral cage distortion	[1269]
Cs <sup>+</sup> 3-Co(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> <sup>-</sup> Cs <sup>+</sup> interactions with calix[4]arene-bis( <i>t</i> -octylbenzo)-18-crown-6	DFT	[1446]

Continued

Compound	Information	References
3- <b>Co</b> (1,2-(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-I) <sup>-</sup> Pd-catalyzed B-C <sub>vinyl</sub> coupling	DFT: reaction profile	[1448]
3- <b>Co</b> (1,2-(C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> )(1',2'-R'C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> R, R'=H, NH <sub>2</sub> , NO <sub>2</sub> , Me, OMe, Ph, Cl nonlinear optical properties	DFT: β <sub>tot</sub> (first-order hyperpolarizability), dipole moment, UV	[1575]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> ) <b>Rh</b> (C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )	Extended Hückel, charge distribution	[1091]
3- <b>Co</b> [1,2-(H <sub>2</sub> NCH <sub>2</sub> )R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ](1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>-</sup>	DFT: geometry; GIAO, <sup>11</sup> B shifts	[1564]
<i>Pseudocloso</i> -3,1,2-Cp* <b>Rh</b> [(PhCH <sub>2</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	DFT, solid state	[1093]
<i>Pseudocloso</i> -3,1,2-(η <sup>3</sup> -C <sub>8</sub> H <sub>13</sub> ) <b>Rh</b> [(MeC <sub>6</sub> H <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ]	DFT, solid state	[1093]
1,2,4,12-(C <sub>8</sub> H <sub>12</sub> ) <b>Ir</b> {[12-(Me <sub>3</sub> C)HN]C <sub>3</sub> B <sub>8</sub> H <sub>10</sub> }	DFT: Fe bonding	[1529]
3- <b>Ni</b> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sub>2</sub> -L L = naphthalene, pyrene, PhNMe <sub>2</sub>	Ligand orientation, charge transfer	[1201]
3- <b>Ni</b> (1,2-R <sub>2</sub> C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> R = H, Me	DFT: HOMO/LUMO vertical electronic excitation	[1202]
3- <b>Ni</b> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -8-SMe <sub>2</sub> ) <sub>2</sub>	Dicarbollide rotational energy	[772]
3- <b>Ni</b> <sup>III</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6-(CH <sub>2</sub> ) <sub>n</sub> C <sub>16</sub> H <sub>9</sub> <sup>-</sup> n = 3,5,7	DFT: conformer energies	[1541]
3- <b>Ni</b> <sup>IV</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6-(CH <sub>2</sub> ) <sub>n</sub> C <sub>16</sub> H <sub>9</sub> n = 3,5,7 C <sub>16</sub> H <sub>9</sub> = pyrene	DFT: conformer energies	[1541]
3- <b>Ni</b> <sup>III</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6'-(CH <sub>2</sub> ) <sub>5</sub> C <sub>16</sub> H <sub>9</sub> <sup>-</sup> C <sub>16</sub> H <sub>9</sub> = pyrene	DFT: conformer energies	[1541]
3- <b>Ni</b> <sup>IV</sup> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1',2'-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -6'-(CH <sub>2</sub> ) <sub>5</sub> C <sub>16</sub> H <sub>9</sub> C <sub>16</sub> H <sub>9</sub> = pyrene	DFT: conformer energies	[1541]
2,1-(Me <sub>3</sub> CNC) <sub>2</sub> <b>Pd</b> [(Me <sub>3</sub> N)CB <sub>10</sub> H <sub>10</sub> ]	CB <sub>10</sub> ligand slip-distortion	[1215]
3,1,2-(η <sup>4</sup> -C <sub>8</sub> H <sub>12</sub> ) <b>Pd</b> (C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+</sup>	Extended Hückel, charge distribution	[1091]
2,1-(PH <sub>3</sub> ) <sub>2</sub> <b>Pt</b> (CB <sub>10</sub> H <sub>11</sub> ) <sup>-</sup>	Extended Hückel, d orbital contributions	[1272]
2,1-(PhSe)(Et <sub>3</sub> P) <b>Pt</b> (CB <sub>10</sub> H <sub>10</sub> -n-SePh) n = 3, 7	Extended Hückel	[1228]
<i>Other Calculations</i>		
3,1,2-(η <sup>6</sup> -C <sub>6</sub> H <sub>6</sub> ) <b>Fe</b> (C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-L <sup>+</sup> R = H, Me; L = SMe <sub>2</sub> , NMe <sub>3</sub> )	DFT, redox potentials	[743]
3,1,2-(η <sup>5</sup> -C <sub>6</sub> H <sub>7</sub> ) <b>Fe</b> (C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> -4-L R = H, Me; L = SMe <sub>2</sub> , NMe <sub>3</sub> )	DFT, redox potentials	[743]
3,1,2-L <sub>3</sub> <b>Fe</b> (C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sup>+</sup> L = Me <sub>3</sub> CNC, P(OMe <sub>3</sub> ) <sub>3</sub>	DFT, redox potentials	[743]
3- <b>Fe</b> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> ) <sub>2</sub>	DFT, redox potentials	[743]
3- <b>Fe</b> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-SMe <sub>2</sub> )(1,2-C <sub>2</sub> B <sub>9</sub> H <sub>10</sub> -4-NMe <sub>2</sub> )	DFT, redox potentials	[743]
(C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <b>Co</b> (C <sub>2</sub> B <sub>8</sub> H <sub>10</sub> ) <b>Co</b> (C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>n-</sup>	Magnetic properties of <i>nido</i> -C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> <sup>2-</sup> and <i>arachno</i> -C <sub>2</sub> B <sub>8</sub> H <sub>10</sub> <sup>4-</sup> ligands	[1048]
3- <b>Co</b> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> Cl <sub>3</sub> ) <sub>2</sub> <sup>-</sup> separation of Am and Pu from products of irradiated Be		[1567]
3- <b>Co</b> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>8</sub> Cl <sub>3</sub> ) <sub>2</sub> <sup>-</sup> solvent system with nitrophenyloctyl ether for extraction of radioactive Cs from acidic wastes	Diffusion coefficient	[1544]
3- <b>Co</b> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>9</sub> ) <sub>2</sub> -1.1'-P(E)X-8,8'-C <sub>6</sub> H <sub>4</sub> <sup>-</sup> X = PCMe <sub>3</sub> , PPh, P(E)Ph; E = O, S, Se diansa-metallacyclophanes	<sup>11</sup> B, NMR <sup>31</sup> P NMR	[1456]
3- <b>Co</b> (1,2-C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> )(1,2-RC <sub>2</sub> B <sub>9</sub> H <sub>10</sub> ) <sup>-</sup> R = C(O)OH, CH <sub>2</sub> C(O)OH, (CH <sub>2</sub> ) <sub>2</sub> C(O)OH, OCH <sub>2</sub> C(O)OH	DFT: <sup>11</sup> B shifts, geometry	[1528]
3,1,2-(C <sub>6</sub> H <sub>6</sub> ) <b>Ir</b> (C <sub>2</sub> B <sub>9</sub> H <sub>11</sub> ) <sup>+</sup>	DFT, Ir-catalyzed decarboxylation of benzoic acid	[1584]

<sup>a</sup>Transition metals and other heteroatoms (other than carbon) incorporated into the cluster framework are in **boldface**.

<sup>b</sup>S, synthesis; X, X-ray diffraction; H, <sup>1</sup>H NMR; B, <sup>11</sup>B NMR; C, <sup>13</sup>C NMR; F, <sup>19</sup>F NMR; P, <sup>31</sup>P NMR; Li, <sup>7</sup>Li NMR; Pt, <sup>195</sup>Pt NMR; Si, <sup>29</sup>Si NMR; 2d, two-dimensional (COSY) NMR; IR, infrared data; MS, mass spectroscopic data; UV, UV-visible data; E, electrochemical data; CD, circular dichroism; ESR, electron spin resonance data; NLO, nonlinear optical data; COND, electrical conductivity; MAG, magnetic susceptibility; NQR, nuclear quadrupole resonance; OR, optical rotation; XPS, X-ray photoelectron spectra; TGA, thermogravimetric analysis; DSC, differential scanning calorimetry.

<sup>c</sup>For complexes containing exo-polyhedral metal-ligand groups bound to *nido*-C<sub>2</sub>B<sub>9</sub> clusters, see Tables 7-2 and 7-3.