



>>Entry Name: Bis(η^6 -benzene)di- μ -chlorodichlorodiruthenium, 9Cl

Synonym(s): Bis(η^6 -benzene)tetrachlorodiruthenium

CAS Registry Number: 35796-48-6

Extra CAS Registry Numbers(s): 37366-09-9

As Bis(η^6 -benzene)di- μ -bromodibromodiruthenium, [FLM15-P](#) with
X = Cl

Molecular Formula: C₁₂H₁₂Cl₄Ru₂

Molecular Weight: 500.178

Accurate Mass: 499.778004

Percentage Composition: C 28.82%; H 2.42%; Cl 28.35%; Ru 40.41%

Source/Synthesis: Synth. from RuCl₃.xH₂O and 1,3-cyclohexadiene in aq. EtOH

Use/Importance: Alkene hydrogenation catalyst; with BINAP, catalyst for asymmetric hydrogenation of carbonyls; catalyst for coupling, alkylation of thiophenes, furans with ROH; catalyses dehydrogenation of hydroxy esters → α -keto esters and cyanohydrins → α -keto nitriles. Starting material for (C₆H₆)Ru chemistry

Physical Description: Air-stable bright red solid

Melting Point: Mp 242 dec.°

Solubility: Sol. DMSO; insol. most solvs.

Other Data: Halide bridges cleaved by donor solvs., eg. MeCN, DMSO. Ir ν_{RuCl} 295, 259 cm⁻¹

>>Derivative: DMSO complex

Synonym(s): (η^6 -Benzene)dichloro[sulfinylbis[methane]-O]ruthenium, 10Cl. (Benzene)dichloro(methyl sulfoxide)ruthenium, 8Cl

CAS Registry Number: 69058-50-0

Molecular Formula: C₈H₁₂Cl₂ORuS

Molecular Weight: 328.224

Accurate Mass: 327.902937

Percentage Composition: C 29.28%; H 3.68%; Cl 21.60%; O 4.87%; Ru 30.79%; S 9.77%

Source/Synthesis: Formed when the parent compd. is dissolved in DMSO

Use/Importance: Hydrogenation catalyst precursor

Physical Description: Red needles (EtOH aq.)

Melting Point: Mp 180 dec.°

>>Derivative: Polymer

CAS Registry Number: 12628-24-9

Physical Description: Brown solid

Melting Point: Mp 242°

>>Derivative: Monomer

Synonym(s): (η^6 -Benzene)dichlororuthenium, 10Cl, 9Cl

CAS Registry Number: 51831-98-2

Molecular Formula: C₆H₆Cl₂Ru

Molecular Weight: 250.089

Accurate Mass: 249.889002

Percentage Composition: C 28.82%; H 2.42%; Cl 28.35%; Ru 40.41%

References:

- Ogata, I. *et al.*, *Tet. Lett.*, 1970, 3011, (*use*)
Zelonka, R.A. *et al.*, *Can. J. Chem.*, 1972, **50**, 3063, (*synth, ir, pmr*)
Bennett, M.A. *et al.*, *J.C.S. Dalton*, 1974, 233, (*synth, ir, pmr*)
Games, D.E. *et al.*, *J. Organomet. Chem.*, 1980, **193**, 229, (*ms*)
Tanaka, M. *et al.*, *Angew. Chem., Int. Ed.*, 1984, **23**, 518, (*use*)
Jaouhari, R. *et al.*, *Chem. Comm.*, 1986, 1255, (*use, coupling of thiophenes, furans with ROH*)
Organomet. Synth., 1986, **3**, 99, (*synth*)
Baghurst, R. *et al.*, *J. Organomet. Chem.*, 1990, **384**, C57, (*synth*)
Kitamura, M. *et al.*, *Tet. Lett.*, 1991, **32**, 4163, (*use, asymmetric hydrogenation of carbonyls*)