Paths of analysis*

L9

Synthia

October 10, 2022

1 Analysis parameters

Analysis type: Automatic Retrosynthesis

Rules: none selected

Filters: Exclude Diastereoselecitve reactions, Tunnels, FGI, FGI with protections

Max. paths returned: 50

Max. iterations: 2000

Commercial:

1. Max. molecular weight - 1000 g/mol
2. Max. price - 1500 $/g

Published:

1. Max. molecular weight - 1000 g/mol
2. Popularity - 5

My Stockroom:

1. Max. molecular weight - 1000 g/mol

Reaction scoring formula: TUNNEL_COEF*FGI_COEF*STEP*20+1000000*(CONFLICT+NON_SELECTIVITY+FILTERS+PROTECT)

Chemical scoring formula: SMALLER^ 3,SMALLER^ 1.5

Min. search width: 400

Max. reactions per product: 60

*The results stated herein were generated using the proprietary platform owned and maintained by Grzybowski Scientific Inventions, Inc., a subsidiary of Merck KGaA, Darmstadt Germany. The results are provided on an as is basis, and shall be used solely in connection with the rights afforded in the license agreement and for no other purpose.

Strategies: none selected

FGI Coeff: 0

Tunnels Coeff: 0

JSON Parameters: {}

2 Paths

2 paths found. *Paths are sorted by score. Reactions are sorted in appearance order for each path.*

2.1 Path 1

Score: `193.93`

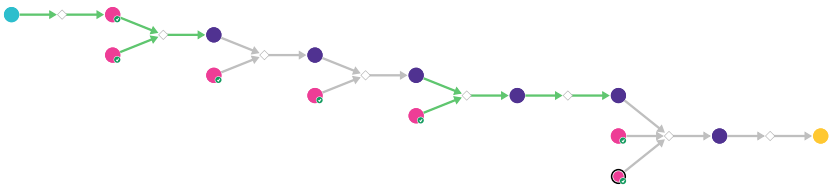
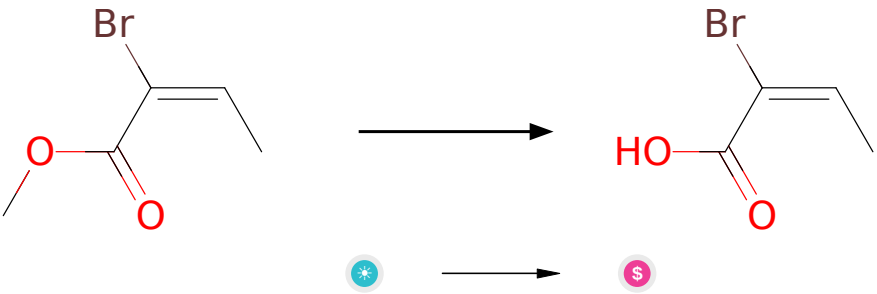


Figure 1: Outline of path 1

2.1.1 Synthesis of Carboxylic Acids via Ester Hydrolysis



Substrates:

1. e-3-methyl-2-bromacrylsaeure-methylester

Products:

1. 2-brom-cis-crotonsaeure - *available at Sigma-Aldrich*

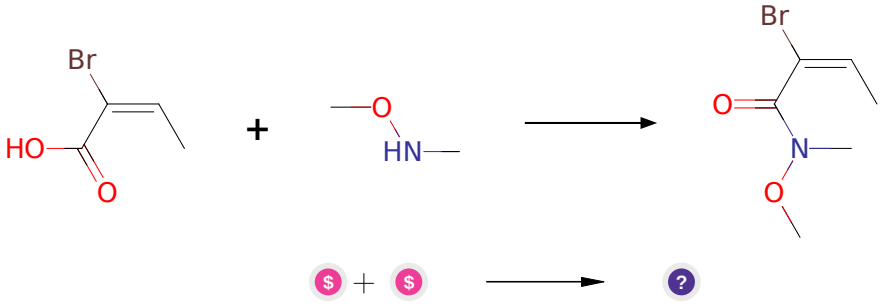
Typical conditions: water.base

Protections: none

Reference: DOI: *10.1016/j.phytochem.2012.08.001* and *10.1021/jm900803q* and *10.1002/anie.201303108* (SI page S14) and *10.1016/j.ejmech.2010.09.003*

Retrosynthesis ID: 9224

2.1.2 Synthesis of O-substituted N-substituted hydroxamic acids



Substrates:

1. 2-brom-cis-crotonsaeure - *available at Sigma-Aldrich*
2. n-methoxymethylamine - *available at Sigma-Aldrich*

Products:

1. C/C=C(/Br)C(=O)N(C)OC

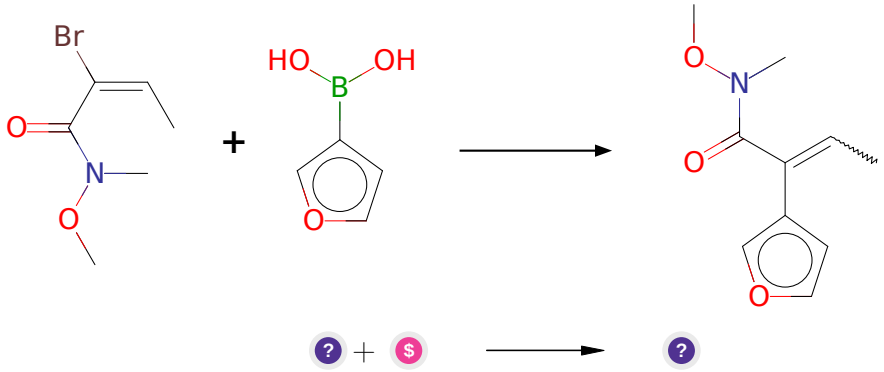
Typical conditions: DCC.DMAP or CDI.TEA.DCM

Protections: none

Reference: Patent: WO2007/67333A2, 2007 & *10.1016/j.bmcl.2008.09.100*

Retrosynthesis ID: 1152

2.1.3 Suzuki coupling of arylboronic acids with vinyl Bromides



Substrates:

1. C/C=C(/Br)C(=O)N(C)OC
2. 3-Furanboronic acid - *available at Sigma-Aldrich*

Products:

1. CC=C(C(=O)N(C)OC)c1ccoc1

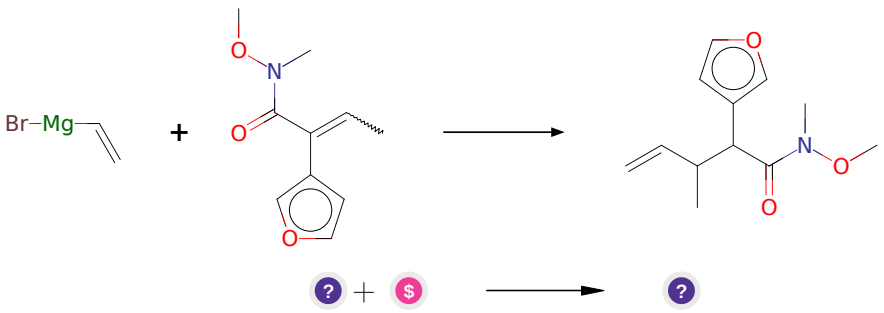
Typical conditions: Pd catalyst.base.solvent

Protections: none

Reference: *10.1021/cr00039a007* and *10.1007/3418_2012_32* and *10.1021/cr0505268* and *10.1016/j.jfluchem.2016.01.018* and *10.1039/C3CS60197H*

Retrosynthesis ID: 24863

2.1.4 Conjugate addition of organocuprate



Substrates:

1. CC=C(C(=O)N(C)OC)c1ccoc1

2. Vinylmagnesium bromide solution - *available at Sigma-Aldrich*

Products:

1. C=CC(C)C(C(=O)N(C)OC)c1ccoc1

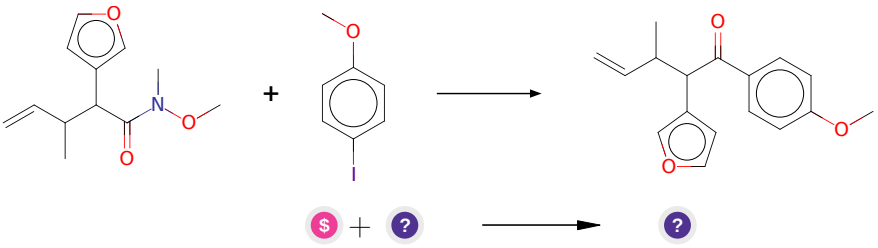
Typical conditions: 1.CuCN.LiCl.2.Eletrophile.3.NH4Cl

Protections: none

Reference: *10.1021/ol036071v* AND *10.1016/j.tet.2011.12.046* AND *10.1002/anie.201007644* AND *10.1002/anie.201007644* AND *10.1055/s-1997-1371*

Retrosynthesis ID: 10003577

2.1.5 Synthesis of ketones from Weinreb amides



Substrates:

1. 4-Iodoanisole - *available at Sigma-Aldrich*
2. C=CC(C)C(C(=O)N(C)OC)c1ccoc1

Products:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)c1ccoc1

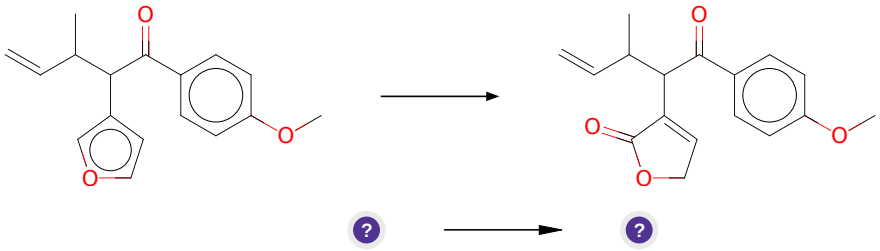
Typical conditions: 1.RmgBr.THF 2.TFA.DCM

Protections: none

Reference: *10.1021/jm051185t* and *10.1021/ol101021v* (supporting info)

Retrosynthesis ID: 5060

2.1.6 Oxidation furans to 2-(5H)-furanones



Substrates:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)c1ccoc1

Products:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)C1=CCOC1=O

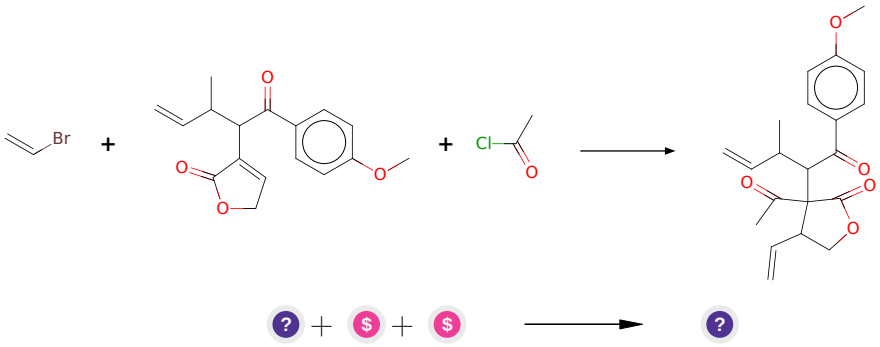
Typical conditions: 1. NBS.CHCl3.EtOH.rt 2. HCl.acetone.H2O.rt

Protections: none

Reference: DOI: *10.1055/s-2005-869865*

Retrosynthesis ID: 50717

2.1.7 Alkenylation-Acylation of enones and enoate esters



Substrates:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)C1=CCOC1=O
2. Bromoethylene - *available at Sigma-Aldrich*
3. Acetyl chloride - *available at Sigma-Aldrich*

Products:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)C1(C(C)=O)C(=O)OCC1C=C

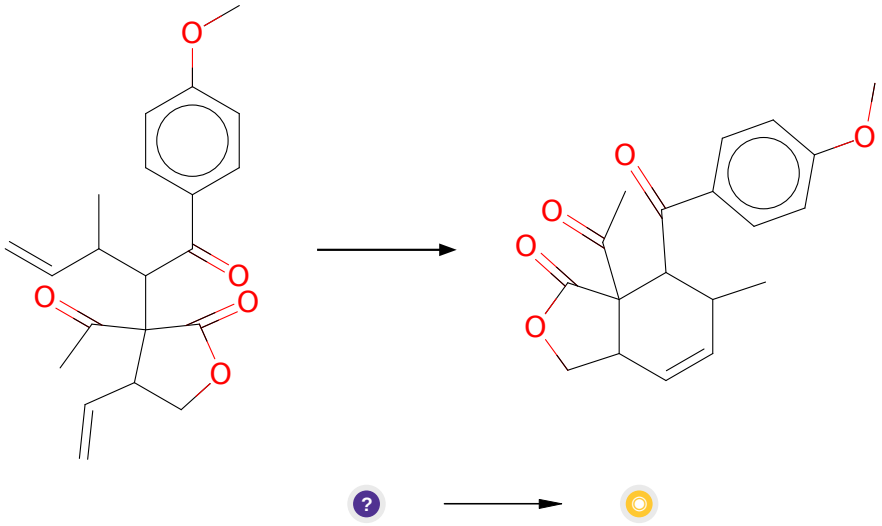
Typical conditions: 1.RCuLi.2.AcCl.HMPA

Protections: none

Reference: *10.1246/cl.1989.1063* AND *10.1248/cpb.33.1815* AND *10.1021/ja0320018* AND *10.1016/S0040-4039(01)80891-1* AND *10.1016/S0040-4020(01)82115-3*

Retrosynthesis ID: 13032

2.1.8 Ring-Closing Metathesis



Substrates:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)C1(C(C)=O)C(=O)OCC1C=C

Products:

1. COc1ccc(C(=O)C2C(C)C=CC3COC(=O)C32C(C)=O)cc1

Typical conditions: catalyst e.g. Hoveyda-Grubbs . solvent e.g. CH2Cl2

Protections: none

Reference: DOI: *10.1002/anie.200800693* and *10.1021/acs.orglett.8b04003* and *10.1021/jo0264729* and *10.1021/ja072334v* and *10.1002/ejoc.201001102*

Retrosynthesis ID: 31014187

2.2 Path 2

Score: 193.93

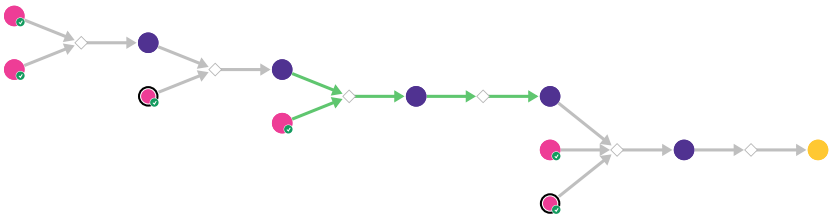
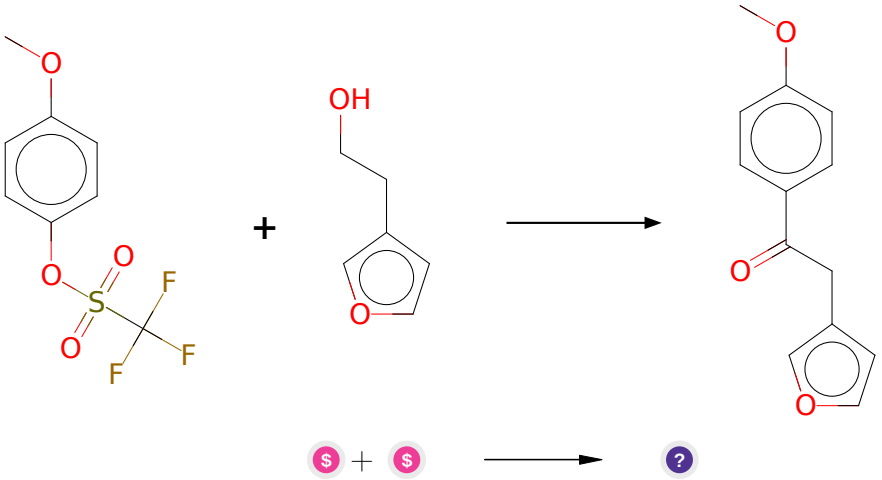


Figure 2: Outline of path 2

2.2.1 Acylation of aryl triflates with primary alcohols



Substrates:

1. 2-(furan-3-yl)ethan-1-ol - *available at Sigma-Aldrich*
2. 4-Methoxyphenyl trifluoromethanesulfonate - *available at Sigma-Aldrich*

Products:

1. COc1ccc(C(=O)Cc2ccoc2)cc1

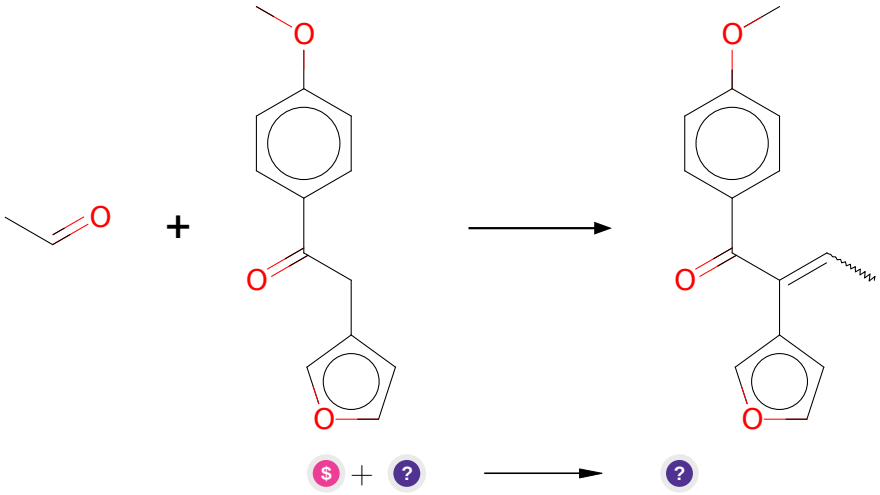
Typical conditions: Ni(cod)2.triphos.TMP.toluene.heat

Protections: none

Reference: *10.1021/jacs.9b03280*

Retrosynthesis ID: 10031467

2.2.2 Aldol Condensation



Substrates:

1. Ethanal - *available at Sigma-Aldrich*
2. COc1ccc(C(=O)Cc2ccoc2)cc1

Products:

1. CC=C(C(=O)c1ccc(OC)cc1)c1ccoc1

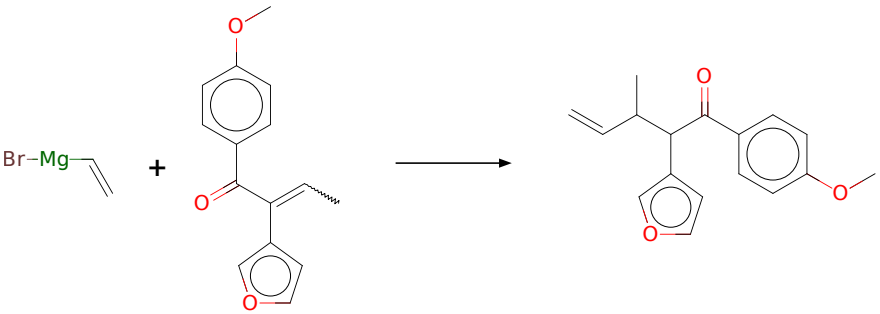
Typical conditions: NaOEt.base

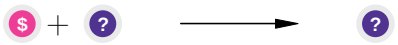
Protections: none

Reference: *10.1080/00397911.2016.1206938*

Retrosynthesis ID: 10049

2.2.3 Conjugate addition of organocuprate





Substrates:

1. Vinylmagnesium bromide solution - *available at Sigma-Aldrich*
2. CC=C(C(=O)c1ccc(OC)cc1)c1ccoc1

Products:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)c1ccoc1

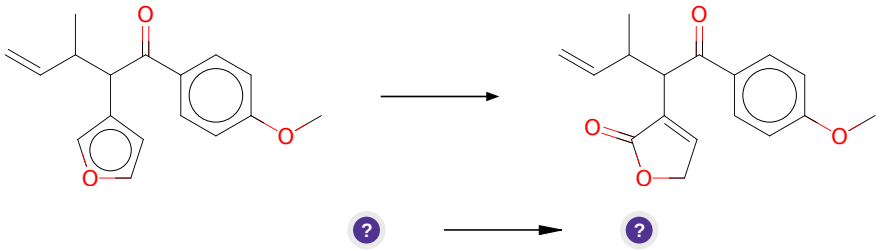
Typical conditions: 1.CuCN.LiCl.2.Eletrophile.3.NH4Cl

Protections: none

Reference: *10.1021/ol036071v* AND *10.1016/j.tet.2011.12.046* AND *10.1002/anie.201007644* AND *10.1002/anie.201007644* AND *10.1055/s-1997-1371*

Retrosynthesis ID: 10003577

2.2.4 Oxidation furans to 2-(5H)-furanones



Substrates:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)c1ccoc1

Products:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)C1=CCOC1=O

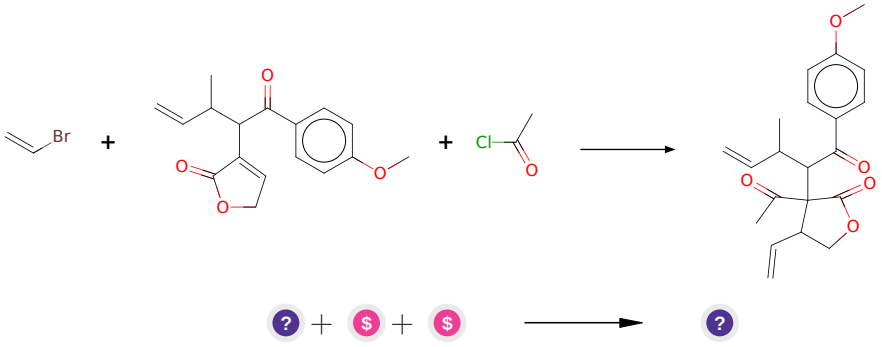
Typical conditions: 1. NBS.CHCl3.EtOH.rt 2. HCl.acetone.H2O.rt

Protections: none

Reference: DOI: *10.1055/s-2005-869865*

Retrosynthesis ID: 50717

2.2.5 Alkenylation-Acylation of enones and enoate esters



Substrates:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)C1=CCOC1=O
2. Bromoethylene - *available at Sigma-Aldrich*
3. Acetyl chloride - *available at Sigma-Aldrich*

Products:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)C1(C(C)=O)C(=O)OCC1C=C

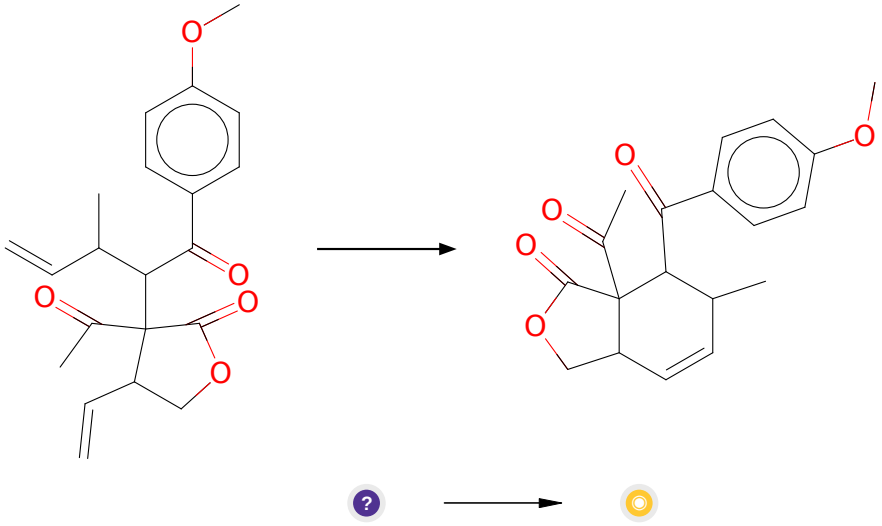
Typical conditions: 1.RCuLi.2.AcCl.HMPA

Protections: none

Reference: *10.1246/cl.1989.1063* AND *10.1248/cpb.33.1815* AND *10.1021/ja0320018* AND *10.1016/S0040-4039(01)80891-1* AND *10.1016/S0040-4020(01)82115-3*

Retrosynthesis ID: 13032

2.2.6 Ring-Closing Metathesis



Substrates:

1. C=CC(C)C(C(=O)c1ccc(OC)cc1)C1(C(C)=O)C(=O)OCC1C=C

Products:

1. COc1ccc(C(=O)C2C(C)C=CC3COC(=O)C32C(C)=O)cc1

Typical conditions: catalyst e.g. Hoveyda-Grubbs . solvent e.g. CH2Cl2

Protections: none

Reference: DOI: *10.1002/anie.200800693* and *10.1021/acs.orglett.8b04003* and *10.1021/jo0264729* and *10.1021/ja072334v* and *10.1002/ejoc.201001102*

Retrosynthesis ID: 31014187