

Paths of analysis*

Analysis 3

Synthia

March 3, 2022

1 Analysis parameters

Analysis type: Automatic Retrosynthesis

Rules: none selected

Filters: FGI, FGI with protections

Max. paths returned: 5

Max. iterations: 300

Commercial:

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

Published:

1. Max. molecular weight - 1000 g/mol
2. Popularity - 10

My Stockroom:

1. Max. molecular weight - 1000 g/mol

Reaction scoring formula: $\text{TUNNEL_COEF} * \text{FGI_COEF} * \text{STEP} * 20 + 1000000 * (\text{CONFLICT} + \text{NON_SELECTIVITY} + \text{FILTERS} + \text{PROTECT})$

Chemical scoring formula: $\text{SMALLER}^3, \text{SMALLER}^{1.5}$

Min. search width: 400

Max. reactions per product: 60

Strategies: none selected

*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.

FGI 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: 106.17

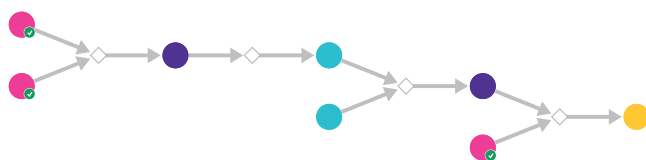
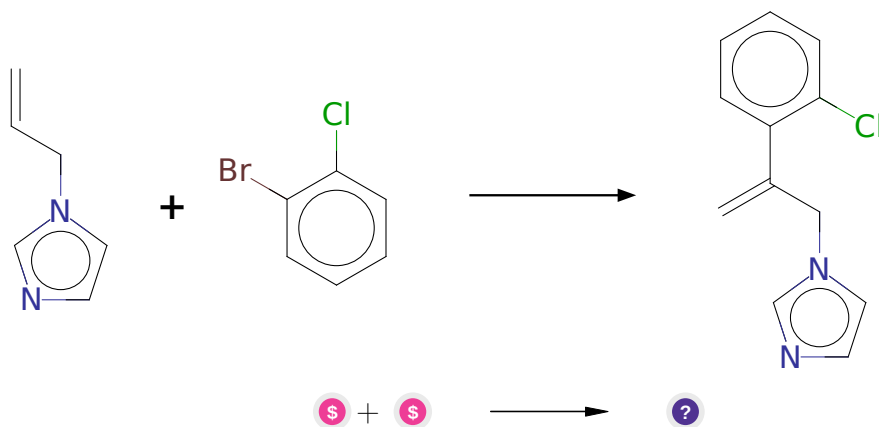


Figure 1: Outline of path 1

2.1.1 Heck Reaction



Substrates:

- 1-Bromo-2-chlorobenzene - *available at Sigma-Aldrich*
- 1-(prop-2-en-1-yl)-1H-imidazole - *available at Sigma-Aldrich*

Products:

- C=C(Cn1ccnc1)c1ccccc1Cl

Typical conditions: Pd (cat). Ligand e.g. TXPTS. Base. Temp

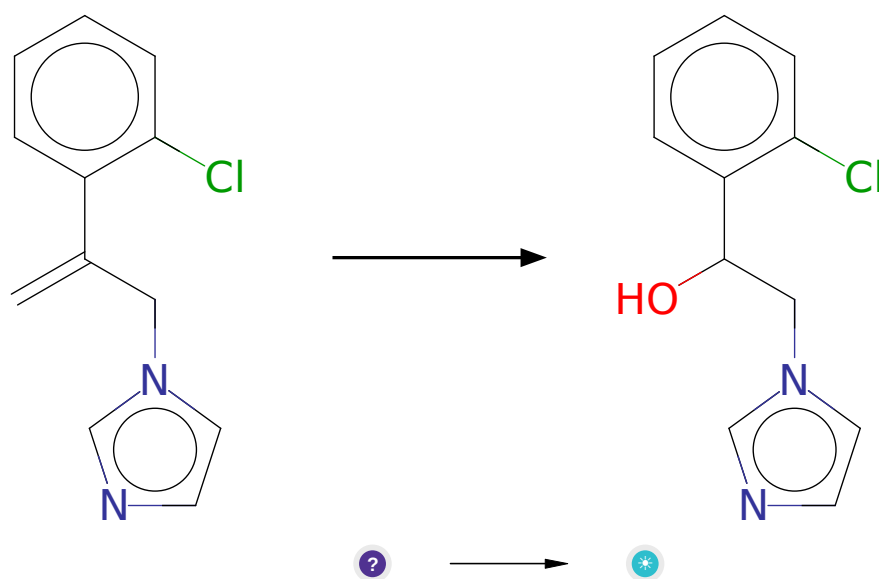
Protections: none

Yield: moderate

Reference: [10.1039/C3CC45911J](#) or [10.1021/ar00049a001](#) or
[10.1002/anie.201201806](#) or [10.1002/9780470716076](#)

Retrosynthesis ID: 9266

2.1.2 Ozonolysis followed by reduction



Substrates:

1. C=C(Cn1ccnc1)c1ccccc1Cl

Products:

1. 1-(2-chloro-phenyl)-2-imidazol-1-yl-ethanol

Typical conditions: O3.MeOH.CH2Cl2.NaBH4.low temperature

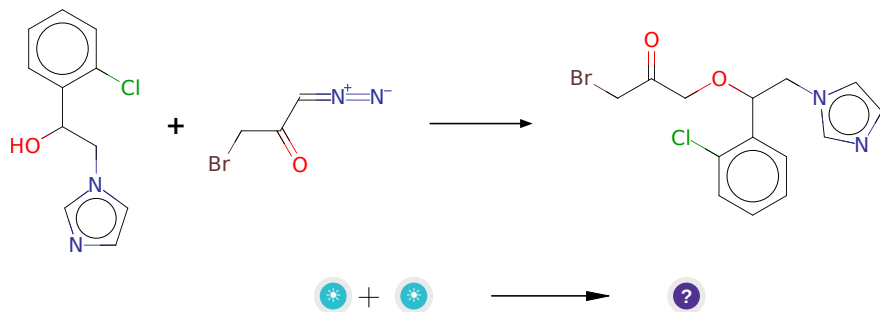
Protections: none

Yield: good

Reference: [10.1021/ja043506g](#)(SI,page S2) and [10.1016/j.jfluchem.2011.05.031](#)
and [10.1021/ja304872j](#) and [10.1021/jo026004z](#)

Retrosynthesis ID: 28553

2.1.3 Alcoholysis of alpha-diazo compounds



Substrates:

1. 1-(2-chloro-phenyl)-2-imidazol-1-yl-ethanol
2. 1-bromo-3-diazo-propan-2-one

Products:

1. O=C(CBr)COC(Cn1ccn1)c1ccccc1Cl

Typical conditions: Rh2(OAc)4

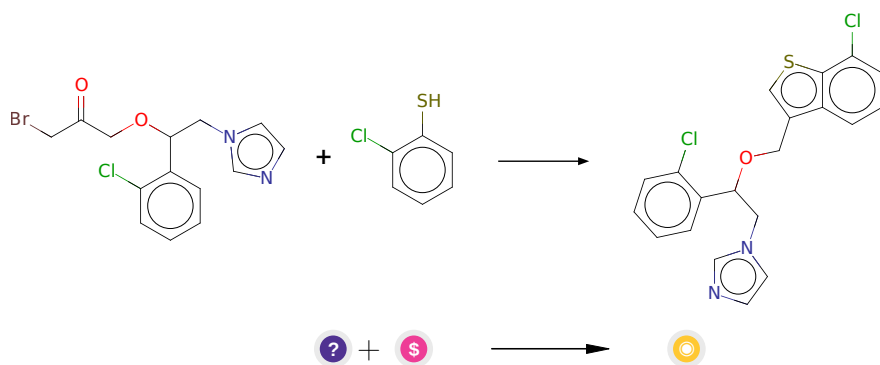
Protections: none

Yield: moderate

Reference: [10.1016/j.tetlet.2014.06.024](https://doi.org/10.1016/j.tetlet.2014.06.024) AND [10.1021/ja074729k](https://doi.org/10.1021/ja074729k) AND [10.1021/ja0607739](https://doi.org/10.1021/ja0607739) AND [10.1039/c4cc06395c](https://doi.org/10.1039/c4cc06395c)

Retrosynthesis ID: 15014

2.1.4 Synthesis of benzothiophenes from thiophenols



Substrates:

1. O=C(CBr)COC(Cn1ccnc1)c1ccccc1Cl
2. 2-Chlorothiophenol - *available at Sigma-Aldrich*

Products:

1. Clc1cccc1C(Cn1ccnc1)OCc1csc2c(Cl)cccc12

Typical conditions: Na₂CO₃.SiO₂.PPA.PhCl.135C

Protections: none

Yield: moderate

Reference: DOI: [10.1055/s-2005-918928](https://doi.org/10.1055/s-2005-918928)

Retrosynthesis ID: 295032

2.2 Path 2

Score: 123.11

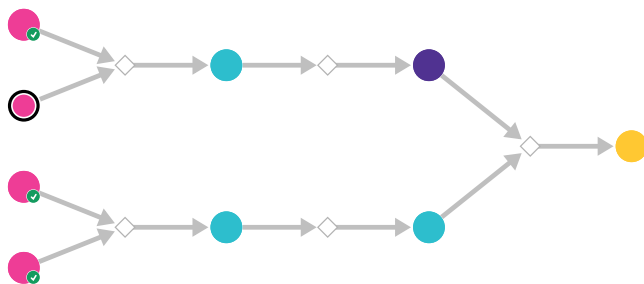
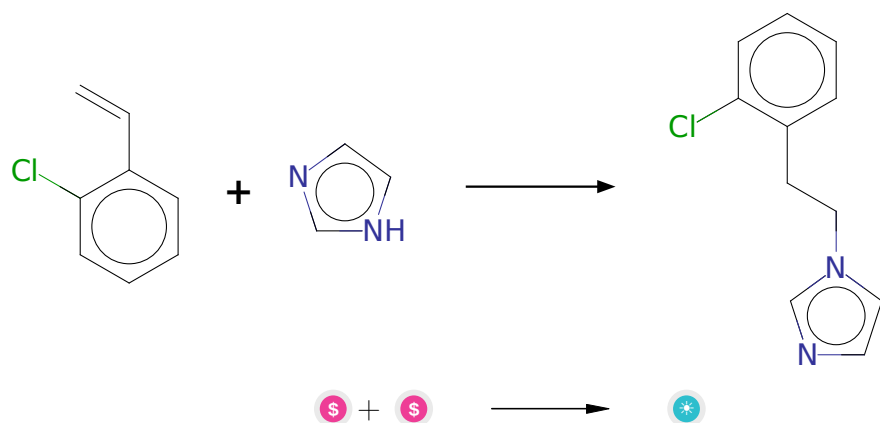


Figure 2: Outline of path 2

2.2.1 Catalytic Photoredox Anti-Markovnikov Alkene Hydroamination



Substrates:

1. 2-Chlorostyrene - *available at Sigma-Aldrich*
2. Imidazole - *available at Sigma-Aldrich*

Products:

1. 1-(2-chloro-phenethyl)-1h-imidazole

Typical conditions: 9-mesityl-10-methylacridinium-perchlorate. Phenyl. Disulfide. 2,6-Lutidine. DCM. Blue. Light. RT

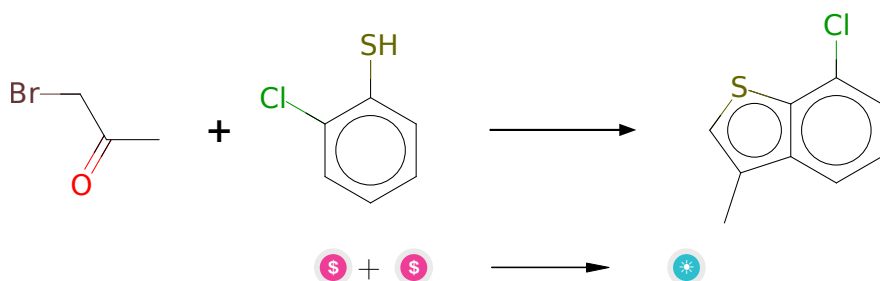
Protections: none

Yield: good

Reference: *10.1002/anie.201402443*

Retrosynthesis ID: 10032751

2.2.2 Synthesis of benzothiophenes from thiophenols



Substrates:

1. 2-Chlorothiophenol - *available at Sigma-Aldrich*
2. brom-aceton - *AstaTech*

Products:

1. 7-chloro-3-methyl-benzo[b]thiophene

Typical conditions: Na₂CO₃.SiO₂.PPA.PhCl.135C

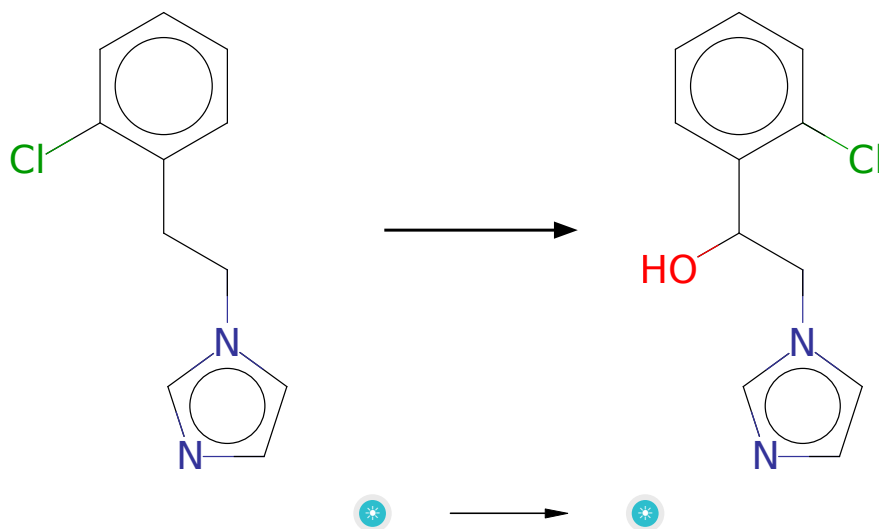
Protections: none

Yield: good

Reference: DOI: *10.1055/s-2005-918928*

Retrosynthesis ID: 295032

2.2.3 Hydroxylation of benzylic position



Substrates:

1. 1-(2-chloro-phenethyl)-1h-imidazole

Products:

1. 1-(2-chloro-phenyl)-2-imidazol-1-yl-ethanol

Typical conditions: 1.Ce(OTf)₄.MeCN.2.NaBH₄

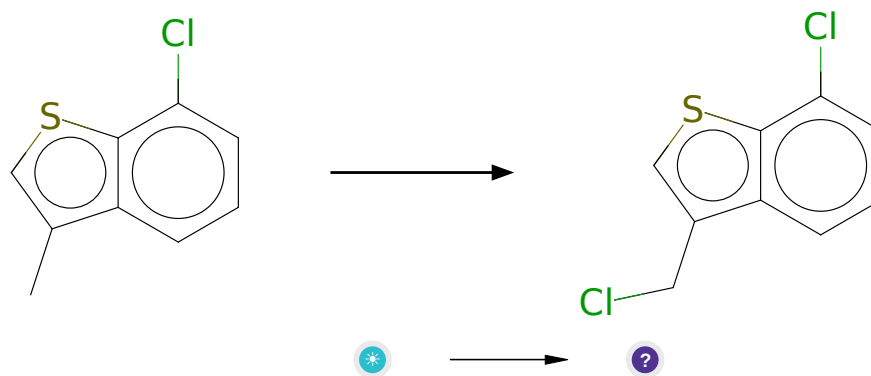
Protections: none

Yield: moderate

Reference: [10.1039/B008843I](#) and WO2012137047 p.12

Retrosynthesis ID: 27140

2.2.4 Chlorination of benzylic position



Substrates:

1. 7-chloro-3-methyl-benzo[b]thiophene

Products:

1. ClCc1csc2c(Cl)cccc12

Typical conditions: SOCl₂.AIBN or NCS/SiCl₄ or [BnNMe₃]ICl₄.AIBN

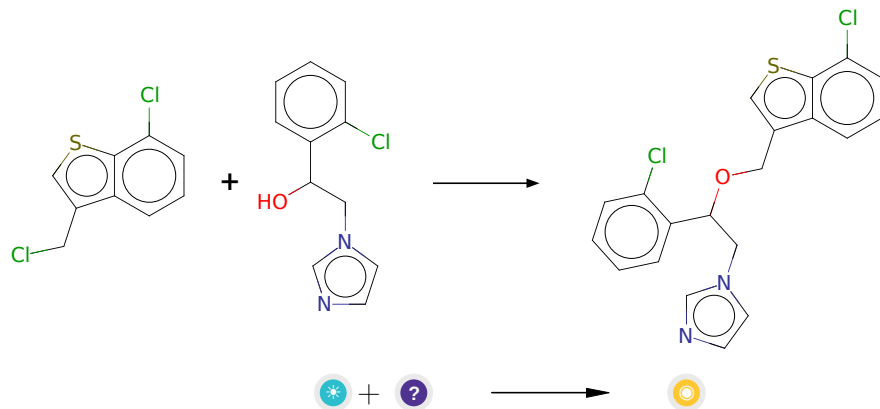
Protections: none

Yield: good

Reference: [10.1039/B803741H](#) and [10.1016/S0040-4039\(00\)82191-7](#) and [10.1016/j.tetlet.2011.05.135](#)

Retrosynthesis ID: 10001786

2.2.5 Alkylation of secondary alcohols



Substrates:

1. 1-(2-chloro-phenyl)-2-imidazol-1-yl-ethanol
2. ClCc1csc2c(Cl)cccc12

Products:

1. Clc1cccc1C(Cn1ccnc1)OCc1csc2c(Cl)cccc12

Typical conditions: K₂CO₃.acetone.heat

Protections: none

Yield: moderate

Reference: [10.1016/S0022-1139\(00\)85021-6](#) and

Retrosynthesis ID: 31011106