Paths of analysis*

Synthia

October 10, 2022

1 Analysis parameters

Analysis type: Automatic Retrosynthesis

Rules: none selected

Filters: Exclude Diastereoselecitve reactions, Tunnels, FGI, FGI with protec-

tions

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

 $\begin{tabular}{ll} \textbf{Reaction scoring formula:} & TUNNEL_COEF*FGI_COEF*STEP*20+1000\\ 0000*(CONFLICT+NON_SELECTIVITY+FILTERS+PROTECT)\\ \end{tabular}$

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.

 ${f Strategies:}\ {f none}\ {f selected}$

FGI Coeff: 0

Tunnels Coeff: 0

JSON Parameters: {}

2 Paths

 $1~\mathrm{path}$ found. Paths are sorted by score. Reactions are sorted in appearance order for each path.

2.1 Path 1

Score: 326.52

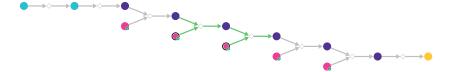
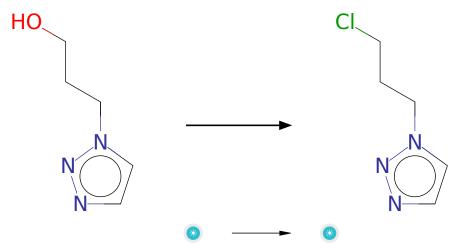


Figure 1: Outline of path 1

2.1.1 Synthesis of alkyl chlorides from alcohols



Substrates:

1. 3-[1,2,3]triazol-1-yl-propan-1-ol

Products:

1. C5H8ClN3

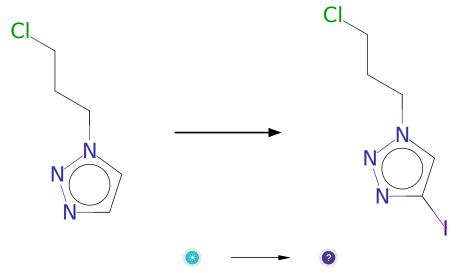
Typical conditions: cyanuric chloride.DMF.DCM.RT

Protections: none

Reference: DOI: 10.1021/ol017168p

Retrosynthesis ID: 11617

2.1.2 Iodination of aromatic compounds



Substrates:

1. C5H8ClN3

Products:

1. ClCCCn1cc(I)nn1

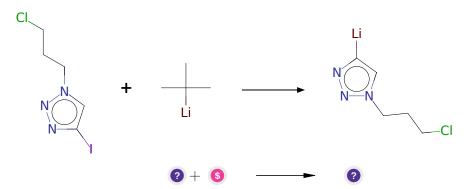
Typical conditions: I2 or other iodinating agent e.g. NIS

Protections: none

Reference: DOI: 10.1039/C5SC00964B and 10.1016/j.tetlet.2005.05.117 and

10.1007/s11178-005-0256-1

2.1.3 I/Li exchange



Substrates:

- $1. \ \, ClCCCn1cc(I)nn1$
- 2. t-BuLi available at Sigma-Aldrich

Products:

1. [Li]c1cn(CCCCl)nn1

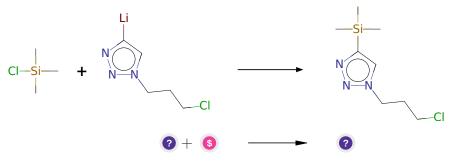
Typical conditions: nBuLi.or.tBuLi.THF.-78C

Protections: none

Reference: 10.1016/j.tet.2004.09.111 and 10.1039/c3ob41082j And 10.1016/j.bmc.2012.03.056 And 10.1002/chem.201300292

Retrosynthesis ID: 30673

2.1.4 Addition of electrophiles to lithiated arenes/heteroarenes



Substrates:

- 1. [Li]c1cn(CCCCl)nn1
- 2. TMSCl available at Sigma-Aldrich

Products:

1. C[Si](C)(C)c1cn(CCCCl)nn1

Typical conditions: THF

Protections: none

Reference: 10.1002/ejoc.200600589 and 10.1055/s-0036-1588863 and 10.1002/1099-0690(200107)2001:14<2771::AID-EJOC2771>3.0.CO;2-Y and 10.1021/ol202873d (SI)

Retrosynthesis ID: 10019541

2.1.5 Blanc bromomethylation

Substrates:

1. C[Si](C)(C)c1cn(CCCCl)nn1

2. Formalin - available at Sigma-Aldrich

Products:

1. C[Si](C)(C)c1nnn(CCCCl)c1CBr

Typical conditions: HBr.heat

Protections: none

Reference: 10.1021/ja011493q and 10.1021/ma012195g and 10.1016/S0040-ma012195g

4039(02)01769-0 and 10.1021/ja002069c

2.1.6 Synthesis of alkyl Grignard reagents

Substrates:

- $1. \ C[Si](C)(C)c1nnn(CCCCl)c1CBr \\$
- 2. Magnesium available at Sigma-Aldrich

Products:

 $1. \ C[Si](C)(C)c1nnn(CCCCl)c1C[Mg]Br \\$

 $\textbf{Typical conditions:} \ \mathrm{Mg.THF} \ \mathrm{or} \ \mathrm{iPrMgBr}$

Protections: none

Reference: DOI: 10.1021/jo00002a039 and 10.1021/jo047877r and

10.1021/ol006618v

2.1.7 Grignard-Type Reaction

Substrates:

1. C[Si](C)(C)c1nnn(CCCCl)c1C[Mg]Br

2. 2-Cyclohexen-1-one - available at Sigma-Aldrich

Products:

1. C[Si](C)(C)c1nnn(CCCCl)c1CC1(O)C=CCCC1

Typical conditions: Mg or Li.ether

Protections: none

Reference: 10.1021/jo010494y or 10.1016/j.steroids.2015.09.009 or 10.1021/jo061349t or 10.1021/ja056165v (SI page 19)

Retrosynthesis ID: 25134

2.1.8 Alkylation of tertiary alcohols

Substrates:

 $1. \ C[Si](C)(C)c1nnn(CCCCl)c1CC1(O)C=CCCC1$

Products:

 $1. \ C[Si](C)(C)c1nnn2c1CC1(C=CCCC1)OCCC2$

Typical conditions: K2CO3.acetone.heat

Protections: none

Reference: 10.1016/S0040-4020(01)90106-1 and 10.1021/acs.analchem.5b04461

and 10.3390/molecules 24091643