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

Reaction scoring formula: TUNNEL_COEF*FGI_COEF*STEP*20+1000 000*(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.

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

FGI Coeff: 0

Tunnels Coeff: 0

JSON Parameters: {}

2 Paths

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

2.1 Path 1

Score: 20.00

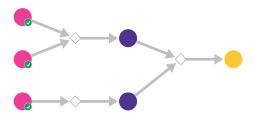
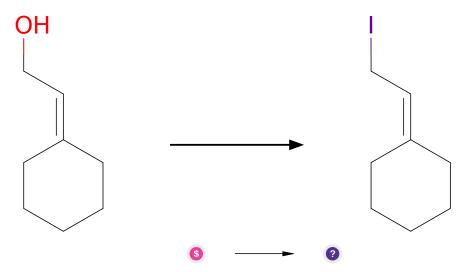


Figure 1: Outline of path 1

2.1.1 Synthesis Of Alkyl Iodides Via Appel Reaction



1. 2-cyclohexylideneethan-1-ol - available at Sigma-Aldrich

Products:

1. ICC=C1CCCCC1

Typical conditions: Imidazole.PPh3.I2

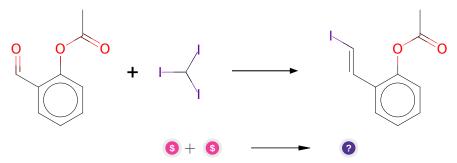
Protections: none

Reference: 10.1002/1099-0690(200102)2001:3<493::AID-EJOC493>3.0.CO2-B

(compound 20) and 10.1016/j.tet.2014.09.030

Retrosynthesis ID: 9990040

2.1.2 Takai olefination



Substrates:

1. Iodoform - available at Sigma-Aldrich

 $2. \ \ 2\text{-formylphenyl acetate} \ \ - \ \ \ \textit{available at Sigma-Aldrich}$

Products:

1. CC(=O)Oc1cccc1/C=C/I

Typical conditions: CrCl2.THF

Protections: none

Reference: 10.1021/ja00283a046 and 10.1021/ja00237a081

2.1.3 Palladium catalysed alkylation of vinyl iodides

Substrates:

- 1. CC(=O)Oc1cccc1/C=C/I
- 2. ICC=C1CCCCC1

Products:

1. CC(=O)Oc1ccccc1/C=C/CC=C1CCCCC1

 $\textbf{Typical conditions:} \ [Pd]. catalyst$

Protections: none

Reference: 10.1016/j.bmcl.2005.12.066 and 10.1021/ol052070m and 10.1021/ol5023195 and 10.1002/anie.200703134 and 10.1016/j.bmcl.2005.09.084 and 10.1021/ol0344873

Retrosynthesis ID: 25162

2.2 Path 2

Score: 20.00



Figure 2: Outline of path 2

2.2.1 Ozonolysis

Substrates:

1. 3-(2-Acetoxyphenyl)-1-propene - available at Sigma-Aldrich

Products:

1. CC(=O)Oc1cccc1CC=O

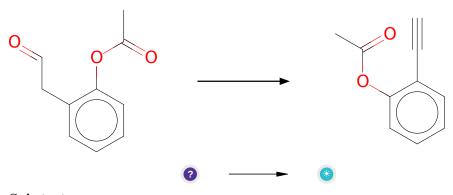
 $\textbf{Typical conditions:} \ \ O3. MeOH. CH2Cl2. PPh3 \ or \ Me2S. low \ temperature$

Protections: none

Reference: 10.1016/j.tet.2017.03.039

Retrosynthesis ID: 5074

2.2.2 Synthesis of alkynes from aldehydes



Substrates:

1. CC(=O)Oc1cccc1CC=O

${\bf Products:}$

$1. \ \ 2\text{-}acetoxyphenylacetylene}$

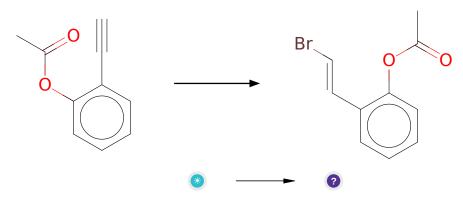
Typical conditions: P1-base.DMF

Protections: none

Reference: 10.1055/s-0028-1087919

Retrosynthesis ID: 15028

2.2.3 Bromination of vinylalanes



Substrates:

1. 2-acetoxyphenylacetylene

Products:

1. CC(=O)Oc1ccccc1/C=C/Br

 ${\bf Typical\ conditions:}\ {\bf Schwartz's\ reagent.then.Br2}$

Protections: none

Reference: DOI: 10.1039/C2CC36604E (SI, page S18) AND DOI:

10.1080/00397910008087318

Retrosynthesis ID: 7405

2.2.4 Suzuki coupling of alkyl-9-BBNs with vinyl bromides



- 1. CC(=O)Oc1ccccc1/C=C/Br
- 2. 9-(3,3-pentamethyleneallyl)-9-borabicyclo3.3.1nonane

Products:

1. CC(=O)Oc1ccccc1/C=C/CC=C1CCCCC1

Typical conditions: Pd catalyst.base.solvent

Protections: none

Reference: 10.1021/ja00183a048 and 10.1039/b707338k and 10.1016/j.tet.2015.05.039 and 10.1021/jo991064z and 10.1021/ol060290+ and 10.1246/bcsj.65.2863

Retrosynthesis ID: 25174

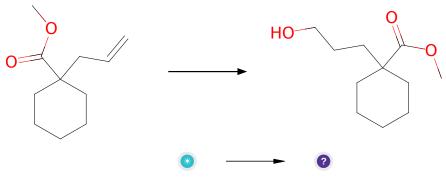
2.3 Path 3

Score: 25.00



Figure 3: Outline of path 3

2.3.1 Brown Hydroboration of Alkenes



Substrates:

1. 1-allyl-cyclohexanecarboxylic acid methyl ester

Products:

1. COC(=O)C1(CCCO)CCCCC1

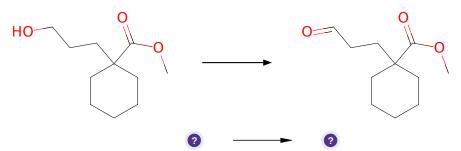
 $\textbf{Typical conditions:} \ B2H6.H2O2.THF.NaOH$

Protections: none

Reference: 10.1002/9780470638859.conrr118

Retrosynthesis ID: 4772

2.3.2 Oxidation of primary alcohols with DMP



Substrates:

1. COC(=O)C1(CCCO)CCCCC1

Products:

 $1. \ \, COC(=O)C1(CCC=O)CCCCC1$

Typical conditions: DMP.DCM.0-25 $\rm C$

Protections: none

Reference: 10.1016/j.bmc.2020.115469 p. 3, 9 and

10.1021/acs.jmedchem.8b01878 SI p. S43

Retrosynthesis ID: 50426

2.3.3 Tebbe Olefination



1. COC(=O)C1(CCC=O)CCCCC1

Products:

1. 1-but-3-enyl-cyclohexanecarboxylic acid methyl ester

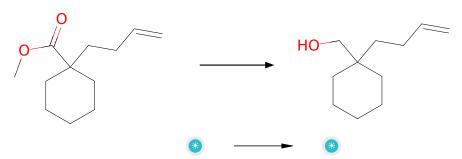
Typical conditions: Cp2TiCl2.AlMe3.toluene

Protections: none

Reference: 10.1016/j.tet.2007.03.015 and 10.1002/9780470638859.conrr617

Retrosynthesis ID: 11714

2.3.4 Esters reduction with LAH



Substrates:

1. 1-but-3-enyl-cyclohexanecarboxylic acid methyl ester

Products:

1. (1-but-3-enyl-cyclohexyl)-methanol

Typical conditions: LiAlH4.THF.0-20 $\rm C$

Protections: none

Reference: 10.1016/j.ejmech.2019.112011 p. 5, 10 and

10.1016/j.ejmech.2020.112910 p. 3, 7

2.3.5 Jones Oxidation

Substrates:

1. (1-but-3-enyl-cyclohexyl)-methanol

Products:

 $1. \ \, acide \ \, 1\text{-}(but\text{-}3\text{-}enyl) cyclohexane carboxylique}$

Typical conditions: cromate.sulfate.H2O.acetone

Protections: none

Reference: 10.1002/9780470638859.conrr349 and 10.1021/jm00270a004

Retrosynthesis ID: 11160

2.3.6 Heck Reaction

Substrates:

- $1. \ \, \hbox{o-bromophenyl acetate}$
- 2. acide 1-(but-3-enyl)cyclohexanecarboxylique

Products:

 $1. \ \mathrm{CC}(=\mathrm{O})\mathrm{Oc1ccccc1/C} = \mathrm{C/CCC1}(\mathrm{C}(=\mathrm{O})\mathrm{O})\mathrm{CCCCC1}$

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

Protections: none

Reference: DOI: 10.1039/C3GC40493E DOI: 10.1021/ol0360288 or DOI: 10.1021/ol702755g or DOI: 10.1055/s-0033-1340319 or DOI: 10.1016/j.tet.2004.10.049

Retrosynthesis ID: 9180

2.3.7 Catalytic dehydrogenative decarboxyolefination of carboxylic acids

Substrates:

1. CC(=O)Oc1ccccc1/C=C/CCC1(C(=O)O)CCCCC1

Products:

1. CC(=O)Oc1ccccc1/C=C/CC=C1CCCCC1

Typical conditions: [Ir]-photocatalyst.[Co]-catalyst.Cs2CO3.DME/H2O.blue.light.rt

Protections: none

Reference: 10.1038/s41557-018-0142-4 and 10.1021/acscatal.8b03282 and 10.1021/acs.joc.9b00167

Retrosynthesis ID: 10032330

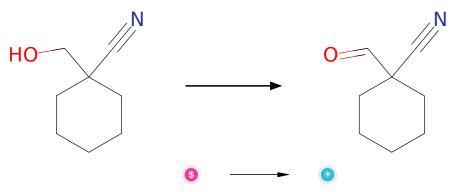
2.4 Path 4

Score: 25.00



Figure 4: Outline of path 4

2.4.1 Oxidation of primary alcohols with DMP



Substrates:

1. 1-(hydroxymethyl)cyclohexane-1-carbonitrile - available at Sigma-Aldrich

Products:

 $1. \ 1 \hbox{-formylcyclohexane} carbonitrile \\$

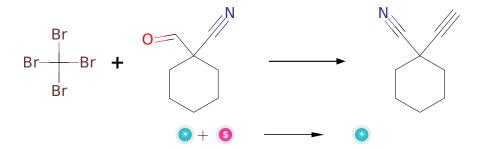
Typical conditions: DMP.DCM.0-25 $\rm C$

Protections: none

Reference: 10.1016/j.bmc.2020.115469 p. 3, 9 and 10.1021/acs.jmedchem.8b01878 SI p. S43

Retrosynthesis ID: 50426

2.4.2 Corey-Fuchs reaction



1. 1-formylcyclohexanecarbonitrile

2. Tetrabromomethane - available at Sigma-Aldrich

Products:

1. 1-ethinylcyclohexylcyanid

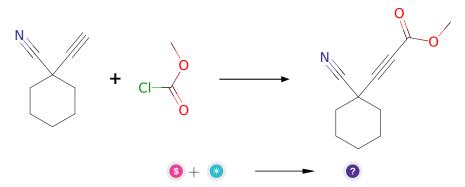
Typical conditions: PPh3.BuLi.CBr4

Protections: none

Reference: 10.1002/ejoc.200601137 and 10.1016/S0040-4039(01)94157-7

Retrosynthesis ID: 10912

2.4.3 Chloroformate Addtion To Terminal Alkynes



Substrates:

1. Methyl chloroformate - available at Sigma-Aldrich

2. 1-ethinylcyclohexylcyanid

Products:

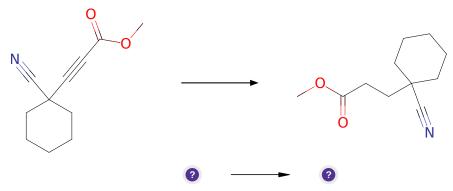
1. COC(=O)C#CC1(C#N)CCCCC1

Typical conditions: 1)n-BuLi.solvent.2) chloroformate

Protections: none

Reference: DOI: 10.1021/jo00092a015

2.4.4 Reduction of alkyne to alkane



Substrates:

1. COC(=O)C#CC1(C#N)CCCCC1

Products:

1. COC(=O)CCC1(C#N)CCCCC1

Typical conditions: H2.Pd/C

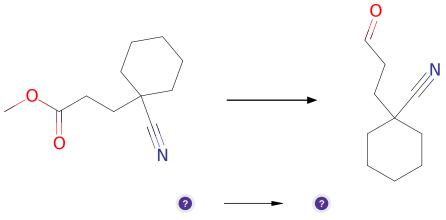
Protections: none

Reference: 10.1016/j.bmc.2011.05.030 AND 10.1021/ol048591b AND

10.1021/jo020486x

Retrosynthesis ID: 14626

2.4.5 Aldehyde Formation



Substrates:

1. COC(=O)CCC1(C#N)CCCCC1

Products:

1. N#CC1(CCC=O)CCCCC1

Typical conditions: DIBAL.solvent e.g. DCM

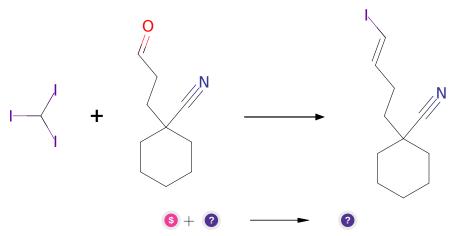
Protections: none

Reference: 10.1039/C39940000483 and 10.1039/C3CC47867J and

10.1021/jo00222a054 and 10.1021/ja9934908 and 10.1021/jo902426z

Retrosynthesis ID: 28551

2.4.6 Takai olefination



Substrates:

1. Iodoform - available at Sigma-Aldrich

 $2.\ \, \text{N}\#\text{CC1}(\text{CCC=O})\text{CCCCC1}$

Products:

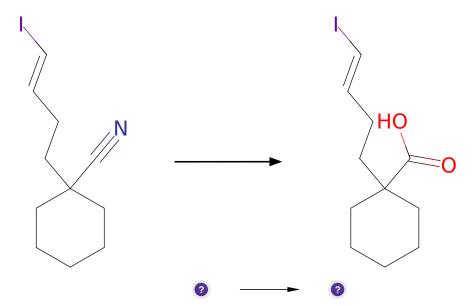
1. N#CC1(CC/C=C/I)CCCCC1

Typical conditions: CrCl2.THF

Protections: none

Reference: 10.1021/ja00283a046 and 10.1021/ja00237a081

2.4.7 Base hydrolysis of nitriles to carboxylic acids



Substrates:

1. N#CC1(CC/C=C/I)CCCCC1

Products:

1. O=C(O)C1(CC/C=C/I)CCCCC1

Typical conditions: NaOH.heating.H2O

Protections: none

Reference: 10.1002/1099-0690(200111)2001:22<4207::AID-EJOC4207>3.0.CO;2-

3 and 10.1021/acs.jmedchem.5b00702 and 10.1016/j.bmc.2011.07.045

Retrosynthesis ID: 15107

2.4.8 Suzuki coupling of arylboronic pinacol esters with vinyl iodides



1. 2-Acetoxyphenylboronic acid pinacol ester - available at Sigma-Aldrich

 $2. \ O{=}C(O)C1(CC/C{=}C/I)CCCCC1$

Products:

1. CC(=O)Oc1ccccc1/C=C/CCC1(C(=O)O)CCCCC1

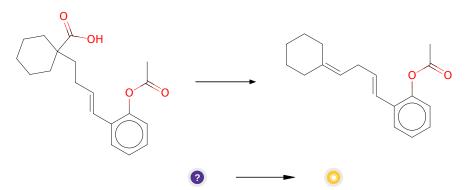
Typical conditions: Pd catalyst.base.solvent

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: 5045

2.4.9 Catalytic dehydrogenative decarboxyolefination of carboxylic acids



Substrates:

1. CC(=O)Oc1ccccc1/C=C/CCC1(C(=O)O)CCCCC1

Products:

1. CC(=O)Oc1ccccc1/C=C/CC=C1CCCCC1

Typical conditions: [Ir]-photocatalyst.[Co]-catalyst.Cs2CO3.DME/H2O.blue.light.rt

Protections: none

Reference: 10.1038/s41557-018-0142-4 and 10.1021/acscatal.8b03282 and 10.1021/acs.joc.9b00167

Retrosynthesis ID: 10032330

2.5 Path 5

Score: 45.00

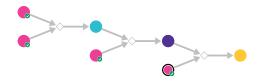
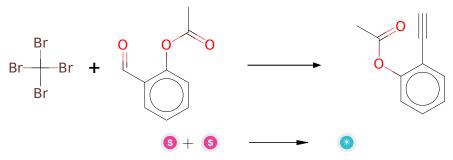


Figure 5: Outline of path 5

2.5.1 Corey-Fuchs reaction



Substrates:

1. 2-formylphenyl acetate - available at Sigma-Aldrich

2. Tetrabromomethane - available at Sigma-Aldrich

Products:

1. 2-acetoxyphenylacetylene

 $\textbf{Typical conditions:} \ PPh 3. Bu Li. CBr 4$

Protections: none

Reference: 10.1002/ejoc.200601137 and 10.1016/S0040-4039(01)94157-7

2.5.2 Hydroalkylation of alkynes

Substrates:

- 1. 2-acetoxyphenylacetylene
- 2. 2-Bromoethyl trifluoromethanesulfonate available at Sigma-Aldrich

Products:

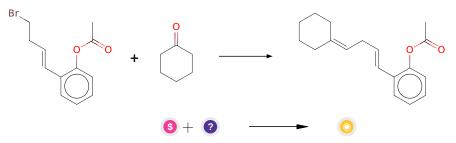
1. CC(=O)Oc1ccccc1/C=C/CCBr

Typical conditions: CsF.(Me2HSi)2O.SIPrCuOTf.dioxane

Protections: none

Reference: 10.1021/ja5124368 Retrosynthesis ID: 33512

2.5.3 HWE/Wittig Olefination



Substrates:

- 1. Cyclohexanone available at Sigma-Aldrich
- 2. CC(=O)Oc1cccc1/C=C/CCBr

Products:

 $1. \ CC(=O)Oc1ccccc1/C=C/CC=C1CCCCC1$

 ${\bf Typical\ conditions:}\ 1. PPh3\ or\ trialkylphosphite. 2. base. aldehyde$

Protections: none

Reference: 10.1002/anie.200705005 and 10.1021/ol052106a and

10.1021/jo00075a064 and 10.1021/ol3027297