Paths of analysis*

O5

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

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

2.1 Path 1

Score: `45.00`

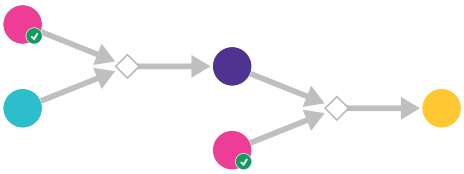
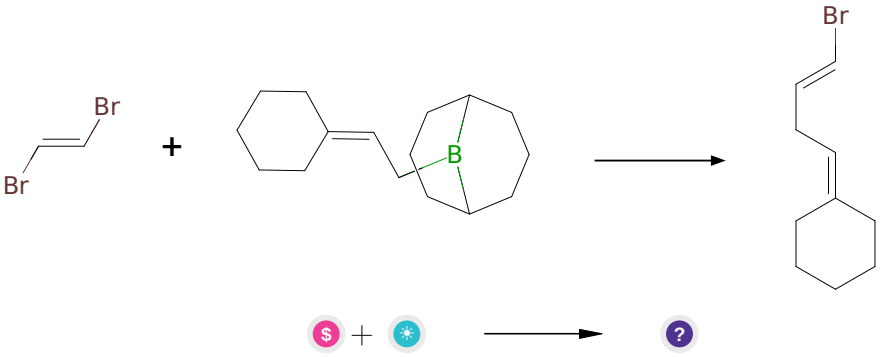


Figure 1: Outline of path 1

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



Substrates:

1. 1,2-Dibromoethylene - *available at Sigma-Aldrich*
2. 9-(3,3-pentamethyleneallyl)-9-borabicyclo3.3.1nonane

Products:

1. Br/C=C/CC=C1CCCCC1

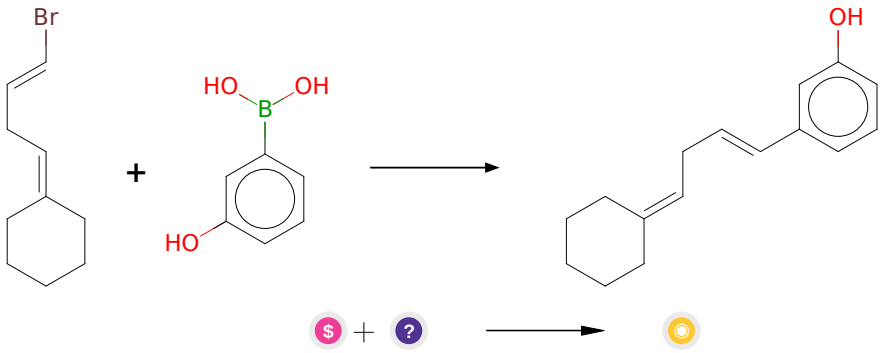
Typical conditions: Pd catalyst.base.solvent

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

2.1.2 Suzuki coupling of arylboronic acids with vinyl Bromides



Substrates:

1. 3-Hydroxyphenylboronic acid - *available at Sigma-Aldrich*
2. Br/C=C/CC=C1CCCCC1

Products:

1. Oc1cccc(/C=C/CC=C2CCCCC2)c1

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: 24862

2.2 Path 2

Score: 45.00

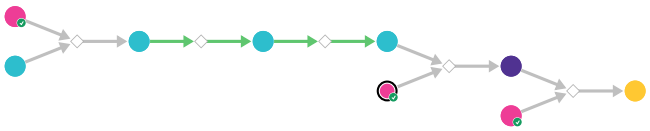
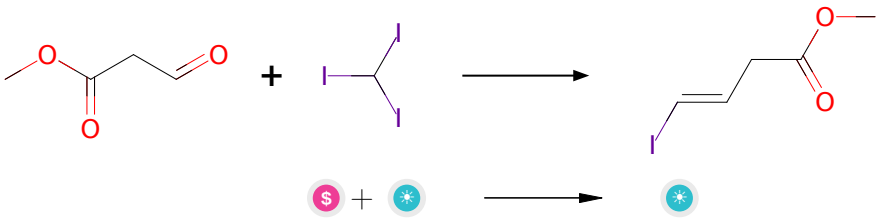


Figure 2: Outline of path 2

2.2.1 Takai olefination



Substrates:

1. Iodoform - *available at Sigma-Aldrich*
2. 3-oxo-propionsaeure-methylester

Products:

1. methyl 4-iodo-3(e)-butenoate

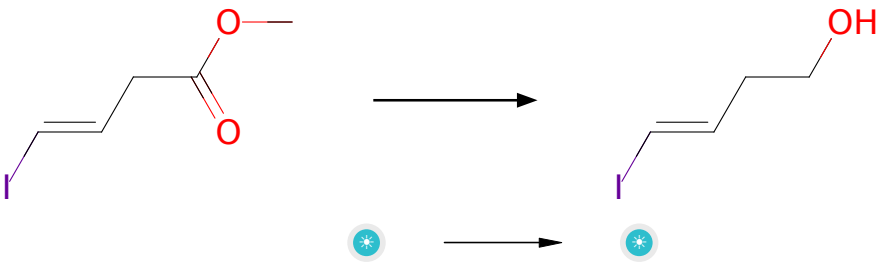
Typical conditions: CrCl2.THF

Protections: none

Reference: *10.1021/ja00283a046* and *10.1021/ja00237a081*

Retrosynthesis ID: 10497

2.2.2 Esters reduction with LAH



Substrates:

1. methyl 4-iodo-3(e)-butenoate

Products:

1. (e)-4-iodo-3-buten-1-ol

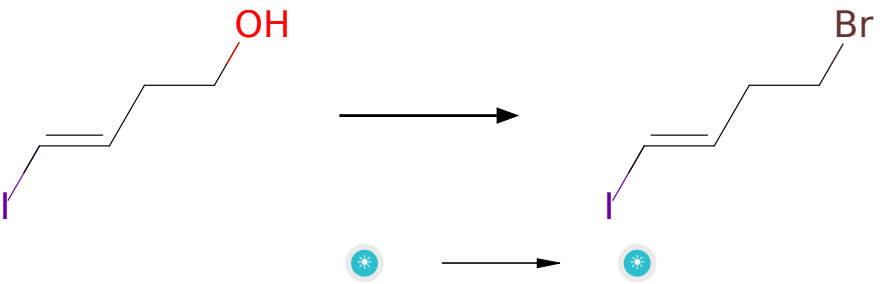
Typical conditions: LiAlH4.THF.0-20 C

Protections: none

Reference: *10.1016/j.ejmech.2019.112011* p. 5, 10 and *10.1016/j.ejmech.2020.112910* p. 3, 7

Retrosynthesis ID: 9910006

2.2.3 Appel Reaction



Substrates:

1. (e)-4-iodo-3-buten-1-ol

Products:

1. C4H6BrI

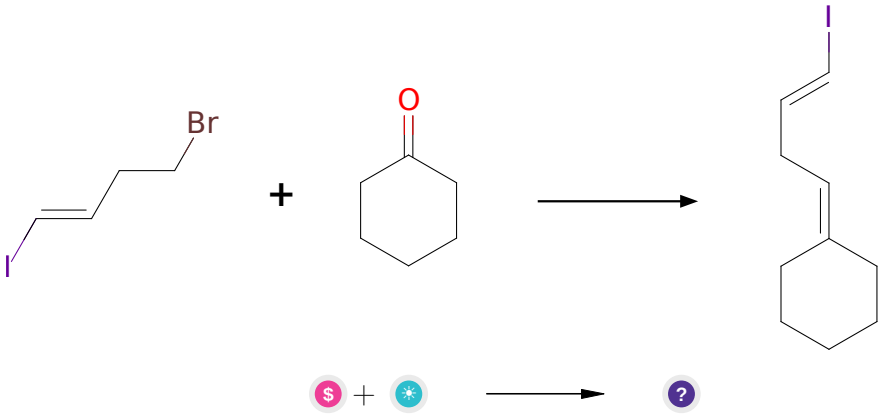
Typical conditions: PPh3.CBr4

Protections: none

Reference: *10.1021/ja800574m* and *10.1016/j.tet.2012.05.010* and *10.1016/j.tet.2004.09.021* (experimantal)

Retrosynthesis ID: 9990037

2.2.4 HWE/Wittig Olefination



Substrates:

1. Cyclohexanone - *available at Sigma-Aldrich*
2. C4H6BrI

Products:

1. I/C=C/CC=C1CCCCC1

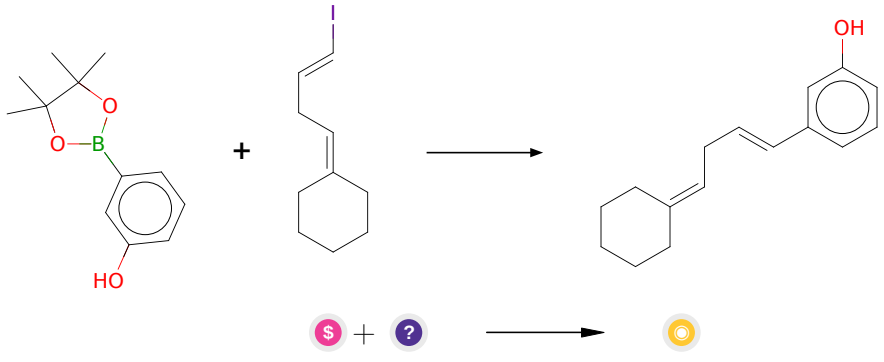
Typical conditions: 1.PPh3 or trialkylphosphite.2.base.aldehyde

Protections: none

Reference: *10.1002/anie.200705005* and *10.1021/ol052106a* and *10.1021/jo00075a064* and *10.1021/ol3027297*

Retrosynthesis ID: 24425

2.2.5 Suzuki coupling of arylboronic pinacol esters with vinyl iodides



Substrates:

1. Pinacol cyclic ester - *available at Sigma-Aldrich*
2. I/C=C/CC=C1CCCCC1

Products:

1. Oc1cccc(/C=C/CC=C2CCCCC2)c1

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

2.3 Path 3

Score: 51.25

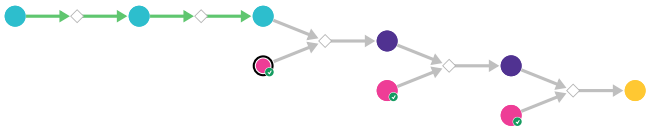
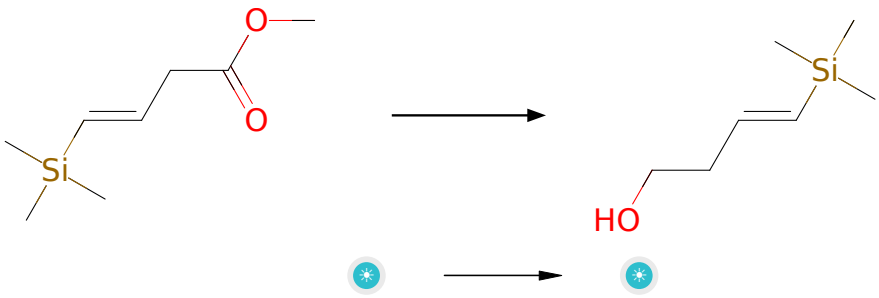


Figure 3: Outline of path 3

2.3.1 Esters reduction with LAH



Substrates:

1. C8H16O2Si

Products:

1. (e)-4-trimethylsilyl-3-buten-1-ol

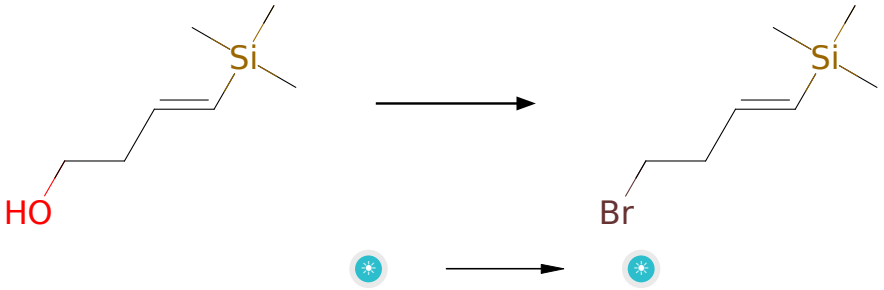
Typical conditions: LiAlH4.THF.0-20 C

Protections: none

Reference: *10.1016/j.ejmech.2019.112011* p. 5, 10 and *10.1016/j.ejmech.2020.112910* p. 3, 7

Retrosynthesis ID: 9910006

2.3.2 Appel Reaction



Substrates:

1. (e)-4-trimethylsilyl-3-buten-1-ol

Products:

1. (4-bromo-but-1-enyl)-trimethyl-silane

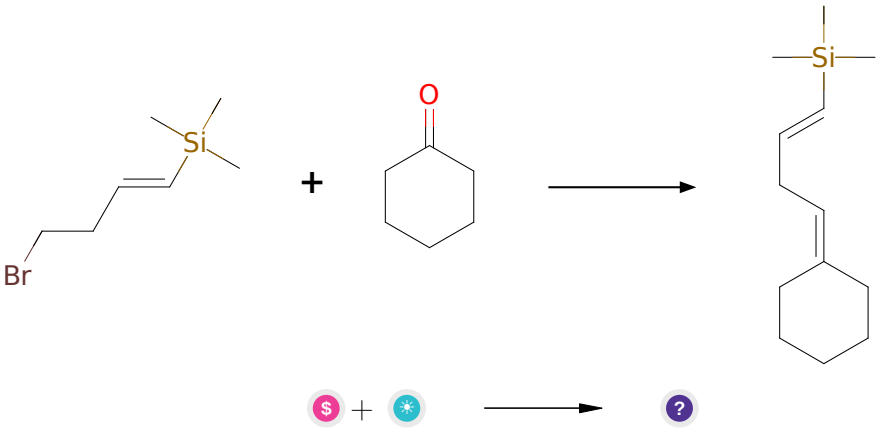
Typical conditions: PPh3.CBr4

Protections: none

Reference: *10.1021/ja800574m* and *10.1016/j.tet.2012.05.010* and *10.1016/j.tet.2004.09.021* (experimantal)

Retrosynthesis ID: 9990037

2.3.3 HWE/Wittig Olefination



Substrates:

1. Cyclohexanone - *available at Sigma-Aldrich*
2. (4-bromo-but-1-enyl)-trimethyl-silane

Products:

1. C[Si](C)(C)/C=C/CC=C1CCCCC1

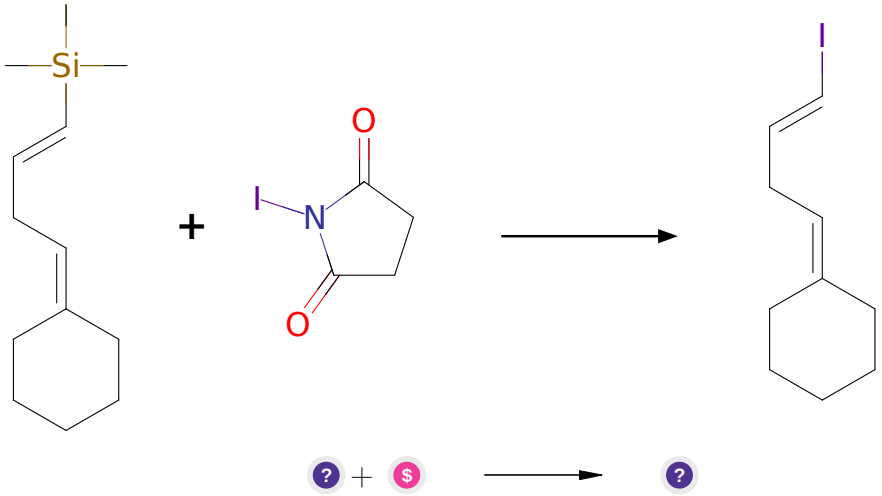
Typical conditions: 1.PPh3 or trialkylphosphite.2.base.aldehyde

Protections: none

Reference: *10.1002/anie.200705005* and *10.1021/ol052106a* and *10.1021/jo00075a064* and *10.1021/ol3027297*

Retrosynthesis ID: 24425

2.3.4 Iodination of Silyl Derivatives



Substrates:

1. C[Si](C)(C)/C=C/CC=C1CCCCC1
2. N-Iodosuccinimide - *available at Sigma-Aldrich*

Products:

1. I/C=C/CC=C1CCCCC1

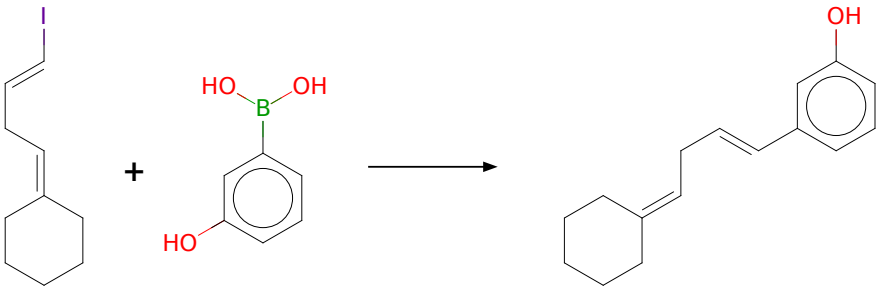
Typical conditions: NIS. 50C. MeCN

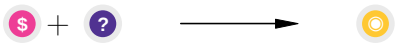
Protections: none

Reference: DOI:*10.1016/j.tetlet.2011.02.057* or DOI: *10.1016/S0040-4039(96)02000-X* or DOI: *10.1016/S0040-4020(02)00334-4*

Retrosynthesis ID: 9211

2.3.5 Suzuki coupling of arylboronic acids with vinyl iodides





Substrates:

1. 3-Hydroxyphenylboronic acid - *available at Sigma-Aldrich*
2. I/C=C/CC=C1CCCCC1

Products:

1. Oc1cccc(/C=C/CC=C2CCCCC2)c1

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: 11208

2.4 Path 4

Score: `51.25`

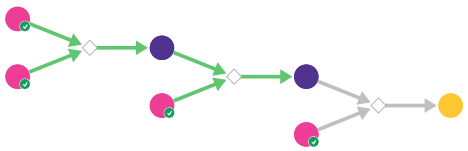
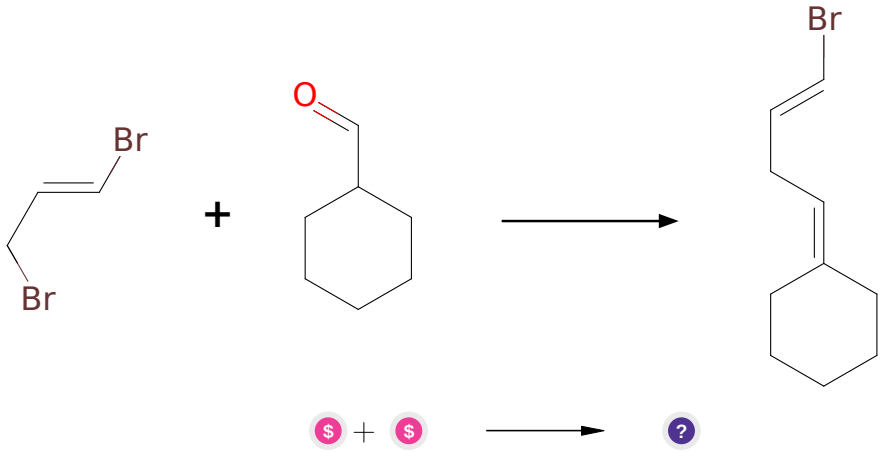


Figure 4: Outline of path 4

2.4.1 Shapiro reaction followed by alkyl bromide addition



Substrates:

1. Hexahydrobenzaldehyde - *available at Sigma-Aldrich*
2. 1,3-Dibromo-1-propene - *available at Sigma-Aldrich*

Products:

1. Br/C=C/CC=C1CCCCC1

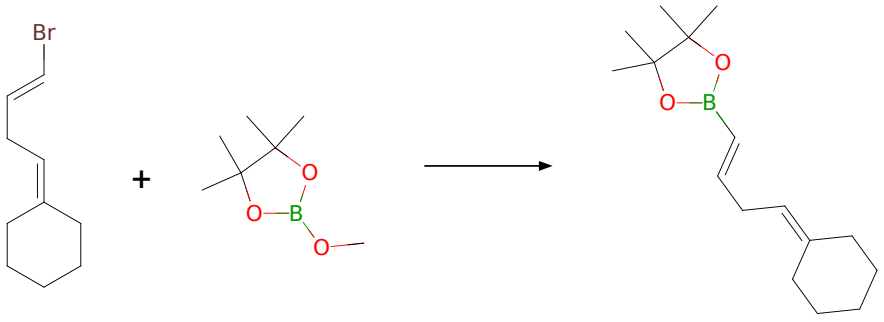
Typical conditions: 1.TsNH2NH2.2.Mes2Mg.LiCl.THF.heating then alkyl bromide.cooling

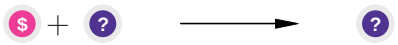
Protections: none

Reference: *10.1016/S0040-4039(00)75263-4* and *10.1021/ol300652k* and *10.1021/jo015699l*

Retrosynthesis ID: 9990458

2.4.2 Synthesis of boronic acid esters





Substrates:

1. Methoxyboronic acid pinacol ester - *available at Sigma-Aldrich*
2. Br/C=C/CC=C1CCCCC1

Products:

1. CC1(C)OB(/C=C/CC=C2CCCCC2)OC1(C)C

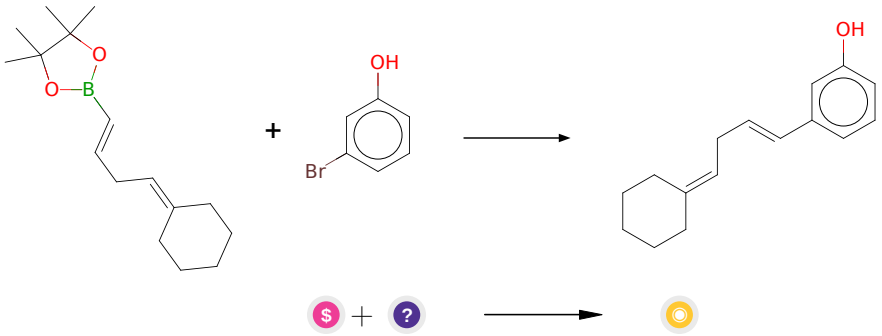
Typical conditions: Mg.THF.-78 C

Protections: none

Reference: *10.1039/B507900D* p. 3167, 3170 and *10.1021/acs.orglett.5b01434* SI p. S6

Retrosynthesis ID: 245548

2.4.3 Suzuki coupling of aryl bromides with alkenyl boronic acids pinacol esters



Substrates:

1. 3-Bromophenol - *available at Sigma-Aldrich*
2. CC1(C)OB(/C=C/CC=C2CCCCC2)OC1(C)C

Products:

1. Oc1cccc(/C=C/CC=C2CCCCC2)c1

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: 11057

2.5 Path 5

Score: 56.25

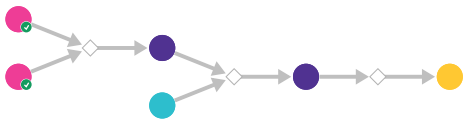
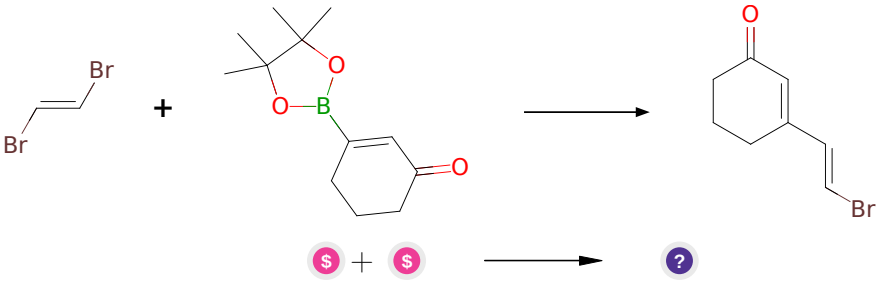


Figure 5: Outline of path 5

2.5.1 Suzuki coupling of vinyl bromides with alkenyl boronic acids pinacol esters



Substrates:

1. 3-(tetramethyl-1,3,2-dioxaborolan-2-yl)cyclohex-2-en-1-one - *available at Sigma-Aldrich*
2. 1,2-Dibromoethylene - *available at Sigma-Aldrich*

Products:

1. O=C1C=C(/C=C/Br)CCC1

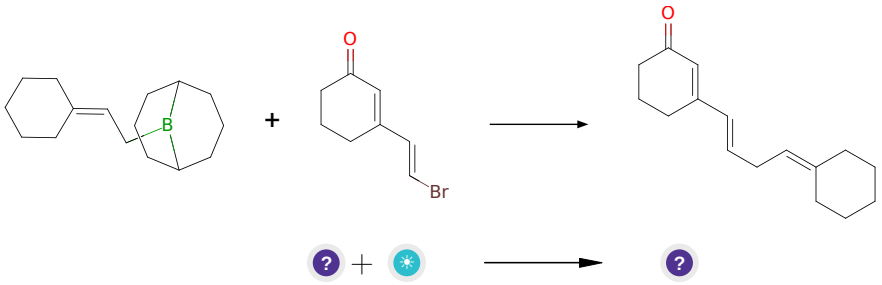
Typical conditions: Pd catalyst.base.solvent

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: 25047

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



Substrates:

1. O=C1C=C(/C=C/Br)CCC1
2. 9-(3,3-pentamethyleneallyl)-9-borabicyclo3.3.1nonane

Products:

1. O=C1C=C(/C=C/CC=C2CCCCC2)CCC1

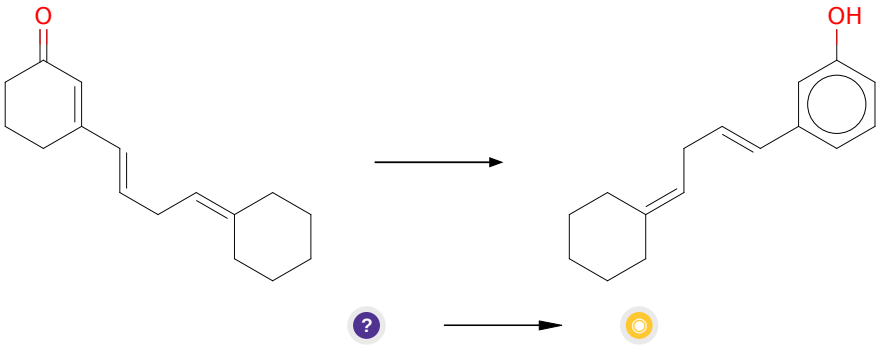
Typical conditions: Pd catalyst.base.solvent

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

2.5.3 DDQ mediated aromatization



Substrates:

1. O=C1C=C(/C=C/CC=C2CCCCC2)CCC1

Products:

1. Oc1cccc(/C=C/CC=C2CCCCC2)c1

Typical conditions: DDQ

Protections: none

Reference: *10.1021/ja054872i* and *10.1021/ja00311a085* and *10.1021/ja00122a011*

Retrosynthesis ID: 9999983