

Paths of analysis*

C128

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

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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: 106.37

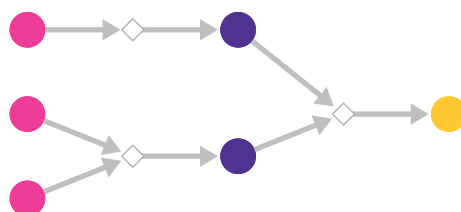
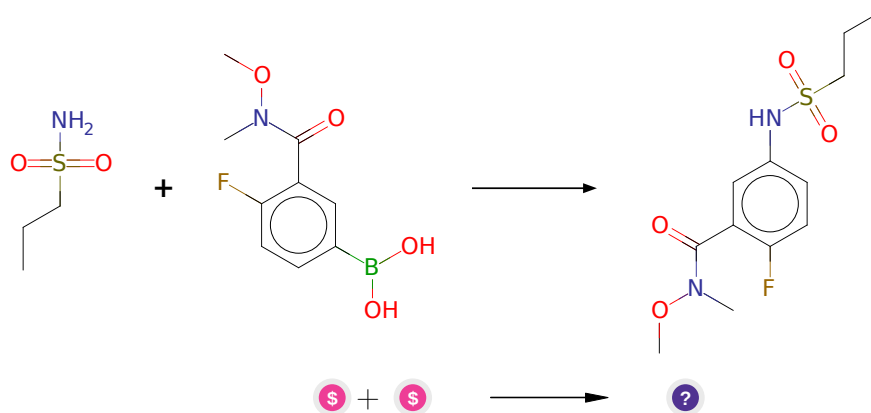


Figure 1: Outline of path 1

2.1.1 Chan-Lam Coupling



Substrates:

1. Propane-1-sulfonamide - *Combi-Blocks*
2. 4-Fluoro-3-[(N-methoxy-N-methyl)carbamoyl]phenylboronic acid - *AOBChem*

Products:

1. CCCS(=O)(=O)Nc1ccc(F)c(C(=O)N(C)OC)c1

Typical conditions: Cu(OAc)₂.K₂CO₃.H₂O or Cu(OAc)₂.pyridine.DCM.MS 4A

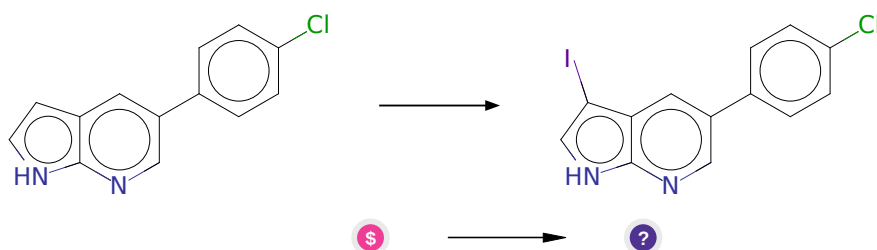
Protections: none

Yield: good

Reference: [10.1016/j.molcata.2014.02.017](https://doi.org/10.1016/j.molcata.2014.02.017) and [10.1039/C4RA08137D](https://doi.org/10.1039/C4RA08137D) and WO2008073956 p.88

Retrosynthesis ID: 31015970

2.1.2 Iodination of aromatic compounds



Substrates:

1. 5-(4-Chlorophenyl)-1H-pyrrolo[2,3-b]pyridine - *Combi-Blocks*

Products:

1. Clc1ccc(-c2cnc3[nH]cc(I)c3c2)cc1

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

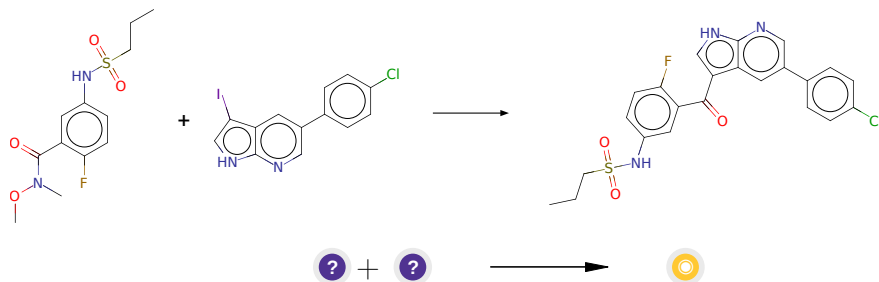
Protections: none

Yield: good

Reference: DOI: [10.1039/C5SC00964B](https://doi.org/10.1039/C5SC00964B) and [10.1016/j.tetlet.2005.05.117](https://doi.org/10.1016/j.tetlet.2005.05.117) and [10.1007/s11178-005-0256-1](https://doi.org/10.1007/s11178-005-0256-1)

Retrosynthesis ID: 10697

2.1.3 Synthesis of ketones from Weinreb amides



Substrates:

1. CCCS(=O)(=O)Nc1ccc(F)c(C(=O)N(C)OC)c1
2. Clc1ccc(-c2ncnc3[nH]cc(I)c3c2)cc1

Products:

1. CCCS(=O)(=O)Nc1ccc(F)c(C(=O)c2c[nH]c3ncc(-c4ccc(Cl)cc4)cc23)c1

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

Protections: none

Yield: good

Reference: [10.1021/jm051185t](#) and [10.1021/ol101021v](#) (supporting info)

Retrosynthesis ID: 5060

2.2 Path 2

Score: 169.66

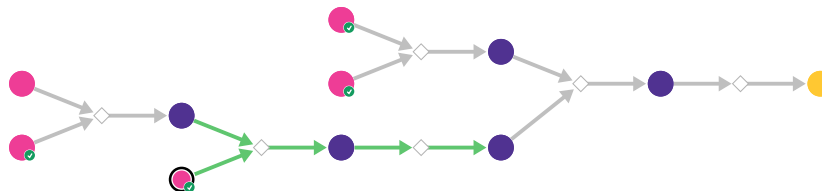
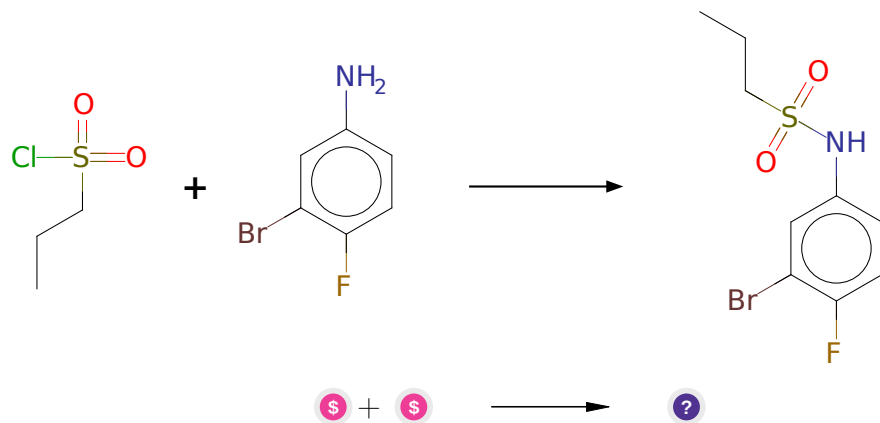


Figure 2: Outline of path 2

2.2.1 N-Sulfonylation



Substrates:

1. 3-Bromo-4-fluoroaniline - *Combi-Blocks*
2. 1-Propanesulfonyl chloride - *available at Sigma-Aldrich*

Products:

1. CCCS(=O)(=O)Nc1ccc(F)c(Br)c1

Typical conditions: THF, rt

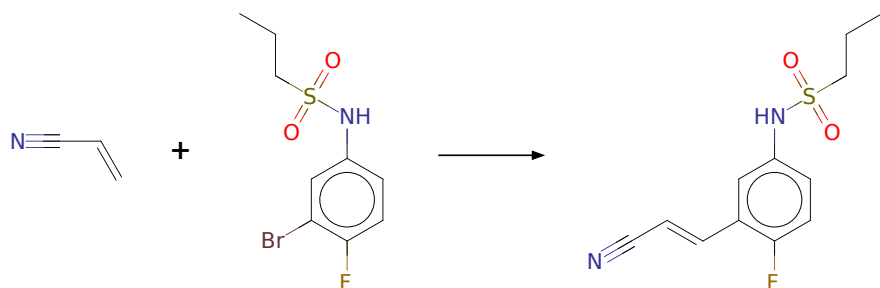
Protections: none

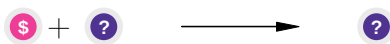
Yield: good

Reference: *10.1055/s-0029-1217565* and *10.1002/(SICI)1099-0690(199806)1998:6<945::AID-EJOC945>3.0.CO;2-3* and *10.1055/s-2001-14567* and *10.1016/j.bmc.2014.07.022*

Retrosynthesis ID: 14717

2.2.2 Heck Reaction





Substrates:

1. Acrylonitrile - *available at Sigma-Aldrich*
2. CCCC(=O)(=O)Nc1ccc(F)c(Br)c1

Products:

1. CCCC(=O)(=O)Nc1ccc(F)c(/C=C/C#N)c1

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

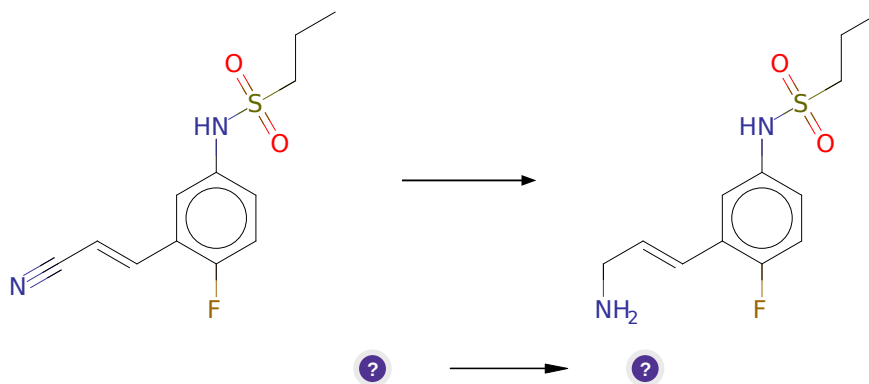
Protections: none

Yield: moderate

Reference: DOI: [10.1039/C3GC40493E](https://doi.org/10.1039/C3GC40493E) DOI: [10.1021/ol0360288](https://doi.org/10.1021/ol0360288) or DOI: [10.1021/ol702755g](https://doi.org/10.1021/ol702755g) or DOI: [10.1055/s-0033-1340319](https://doi.org/10.1055/s-0033-1340319) or DOI: [10.1016/j.tet.2004.10.049](https://doi.org/10.1016/j.tet.2004.10.049)

Retrosynthesis ID: 9180

2.2.3 Reduction of Nitriles to Amines



Substrates:

1. CCCC(=O)(=O)Nc1ccc(F)c(/C=C/C#N)c1

Products:

1. CCCC(=O)(=O)Nc1ccc(F)c(/C=C/CN)c1

Typical conditions: LAH.THF.-78C

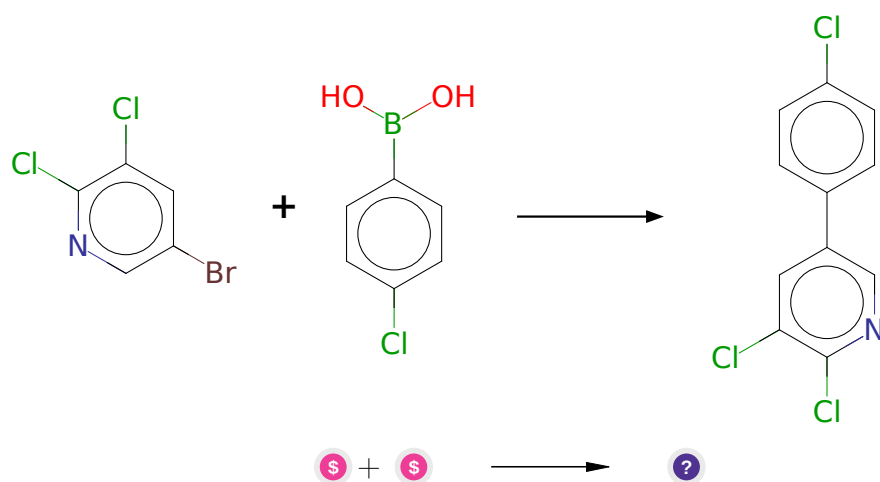
Protections: none

Yield: moderate

Reference: DOI:[10.1021/ja01145a082](https://doi.org/10.1021/ja01145a082)

Retrosynthesis ID: 11142

2.2.4 Suzuki coupling of arylboronic acids with aryl bromides



Substrates:

1. 5-Bromo-2,3-dichloropyridine - [available at Sigma-Aldrich](#)
2. 4-Chlorophenylboronic acid - [available at Sigma-Aldrich](#)

Products:

1. Clc1ccc(-c2nc(Cl)c(Cl)c2)cc1

Typical conditions: Pd catalyst.base.solvent

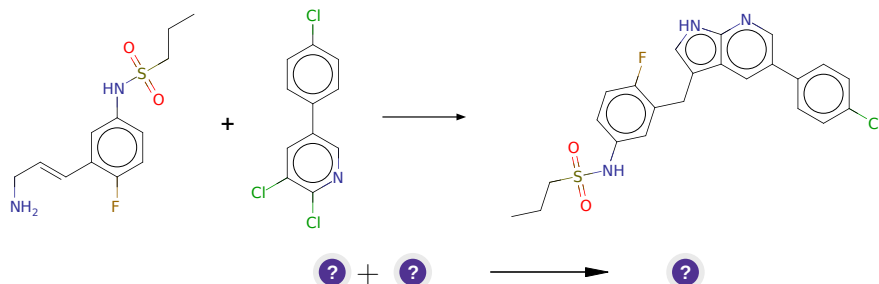
Protections: none

Yield: good

Reference: [10.1021/cr00039a007](https://doi.org/10.1021/cr00039a007) and [10.1007/3418_2012_32](https://doi.org/10.1007/3418_2012_32) and [10.1021/cr0505268](https://doi.org/10.1021/cr0505268) and [10.1016/j.jfluchem.2016.01.018](https://doi.org/10.1016/j.jfluchem.2016.01.018) and [10.1039/C3CS60197H](https://doi.org/10.1039/C3CS60197H) and [10.1016/j.ejmech.2018.08.092](https://doi.org/10.1016/j.ejmech.2018.08.092) and [10.1038/s41929-020-00564-z](https://doi.org/10.1038/s41929-020-00564-z) (metal-free coupling)

Retrosynthesis ID: 25150

2.2.5 Heck-type synthesis of indoles



Substrates:

1. CCCCS(=O)(=O)Nc1ccc(F)c(/C=C/CN)c1
2. Clc1ccc(-c2cnc(Cl)c(Cl)c2)cc1

Products:

1. CCCCS(=O)(=O)Nc1ccc(F)c(Cc2c[nH]c3ncc(-c4ccc(Cl)cc4)cc23)c1

Typical conditions: Pd2dba3.dppf.NaOtBu.PhMe.140C

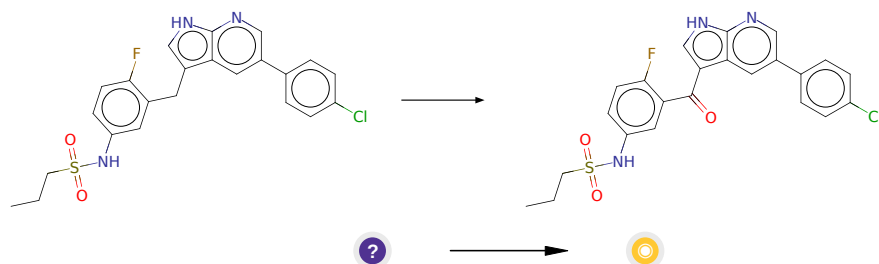
Protections: none

Yield: good

Reference: [10.1002/anie.200703763](#)

Retrosynthesis ID: 28942

2.2.6 Benzylic oxidation to ketone



Substrates:

1. CCCCS(=O)(=O)Nc1ccc(F)c(Cc2c[nH]c3ncc(-c4ccc(Cl)cc4)cc23)c1

Products:

1. CCCCS(=O)(=O)Nc1ccc(F)c(C(=O)c2c[nH]c3ncc(-c4ccc(Cl)cc4)cc23)c1

Typical conditions: oxidant eg. Oxone or O₂ or K₂S₂O₈

Protections: none

Yield: moderate

Reference: [10.1021/ol300853z](#) and [10.1021/ol0520020](#) and [10.1055/s-0037-1610678](#) and [10.1021/acs.orglett.6b02914](#)

Retrosynthesis ID: 7201