

# Paths of analysis\*

Y2A

Synthia

October 10, 2022

## 1 Analysis parameters

**Analysis type:** Automatic Retrosynthesis

**Rules:** none selected

**Filters:** Exclude Diastereoselective 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:**  $\text{TUNNEL\_COEF} * \text{FGI\_COEF} * \text{STEP} * 20 + 1000 * (\text{CONFLICT} + \text{NON\_SELECTIVITY} + \text{FILTERS} + \text{PROTECT})$

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

**Min. search width:** 400

**Max. reactions per product:** 60

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**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:** 2000051.25

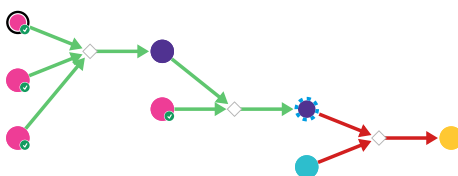
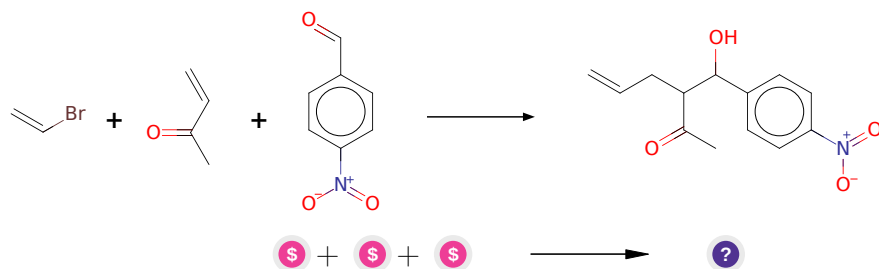


Figure 1: Outline of path 1

#### 2.1.1 Alkenylation-Aldol reaction of enones and enoate esters



**Substrates:**

1. 3-Buten-2-one - *available at Sigma-Aldrich*
2. Bromoethylene - *available at Sigma-Aldrich*
3. 4-Nitrobenzaldehyde - *available at Sigma-Aldrich*

**Products:**

1. C=CCC(C(C)=O)C(O)c1ccc([N+](=O)[O-])cc1

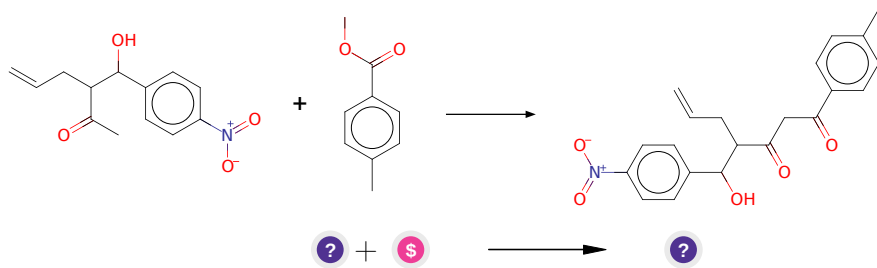
**Typical conditions:** 1.RCuLi.2.RCHO

**Protections:** none

**Reference:** [10.1016/S0040-4039\(01\)80891-1](#) AND [10.1016/S0040-4020\(01\)82115-3](#) AND [10.1021/jo2010186](#) AND [10.1021/jo101439h](#) AND [10.1021/ja906241w](#)

**Retrosynthesis ID:** 20547

### 2.1.2 Condensation of methyl ketones with esters



**Substrates:**

1. C=CCC(C(C)=O)C(O)c1ccc([N+](=O)[O-])cc1
2. Methyl p-toluate - *available at Sigma-Aldrich*

**Products:**

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1

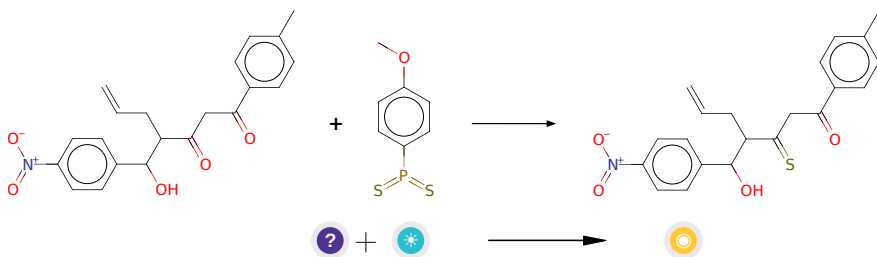
**Typical conditions:** NaOMe.MeOH

**Protections:** none

**Reference:** [10.1016/j.tetlet.2007.10.010](#) and [10.1016/j.tetlet.2013.09.025](#) and [10.1016/j.ejmech.2013.10.072](#) and [10.1002/ange.19921040631](#)

**Retrosynthesis ID:** 4792

### 2.1.3 Synthesis of Thioketones using Lawesson's Reagent



**Substrates:**

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1
2. 4-methoxyphenyl-dithiophosphonsaeureanhydrid

**Products:**

1. C=CCC(C(=S)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1

**Typical conditions:** Lawesson's Reagent.neat.microwave

**Protections:**

Functional group SMARTS	Classification	Protecting groups
<chem>[*]C([*])=O</chem>	carbonyls	1.3-Dioxanes 1.3-Dioxolanes 1.3-Dithianes 1.3-Dithiolanes Dimethyl Acetals and Ketals N,N-Dimethylhydrazones

**Reference:** DOI: [10.1021/ol990629a](https://doi.org/10.1021/ol990629a)

**Retrosynthesis ID:** 11476

## 2.2 Path 2

**Score:** 2000076.25

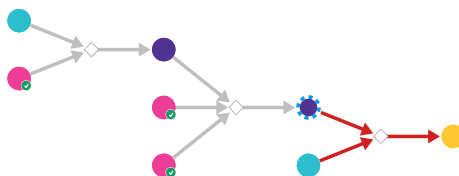
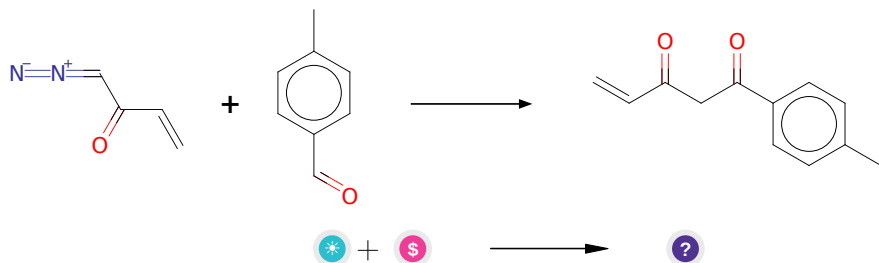


Figure 2: Outline of path 2

### 2.2.1 Homologation of aldehydes to ketones with diazoalkanes



**Substrates:**

1. 1-diazo-but-3-en-2-one
2. p-Tolualdehyde - *available at Sigma-Aldrich*

**Products:**

1. C=CC(=O)CC(=O)c1ccc(C)cc1

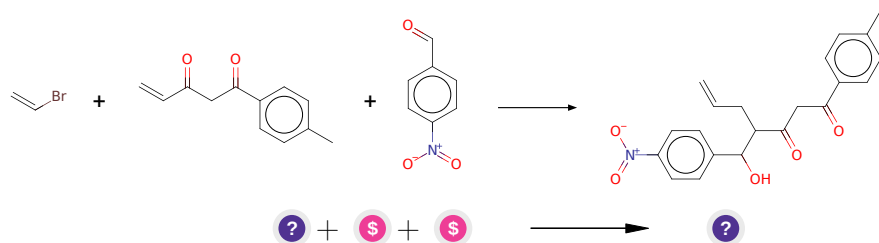
**Typical conditions:** Lewis.acid

**Protections:** none

**Reference:** [10.1021/jo00275a006](#) AND [10.1016/j.tet.2014.05.107](#) AND [10.1016/j.tet.2014.11.059](#) AND [10.1021/ol9010932](#)

**Retrosynthesis ID:** 15017

### 2.2.2 Alkenylation-Aldol reaction of enones and enoate esters



**Substrates:**

1. C=CC(=O)CC(=O)c1ccc(C)cc1
2. Bromoethylene - *available at Sigma-Aldrich*
3. 4-Nitrobenzaldehyde - *available at Sigma-Aldrich*

**Products:**

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1

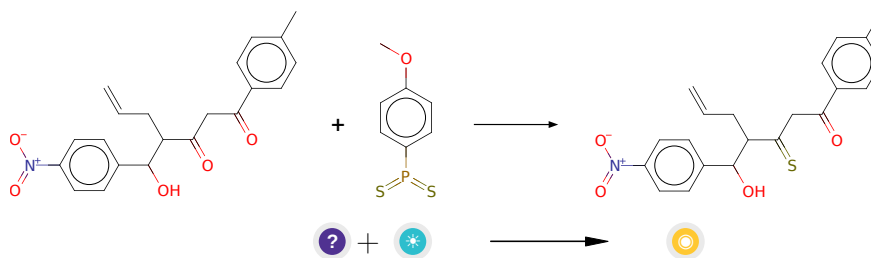
**Typical conditions:** 1.RCuLi.2.RCHO

**Protections:** none

**Reference:** [10.1021/jo2010186](#) AND [10.1021/jo101439h](#) AND [10.1021/ja906241w](#)  
AND [10.1016/S0040-4039\(01\)80891-1](#) AND [10.1016/S0040-4020\(01\)82115-3](#)

**Retrosynthesis ID:** 13048

### 2.2.3 Synthesis of Thioketones using Lawesson's Reagent



**Substrates:**

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1
2. 4-methoxyphenyl-dithiophosphonane

**Products:**

1. C=CCC(C(=S)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1

**Typical conditions:** Lawesson's Reagent.neat.microwave

**Protections:**

Functional group SMARTS	Classification	Protecting groups
[#6]C([#6])=O	carbonyls	1.3-Dioxanes
		1.3-Dioxolanes
		1.3-Dithianes
		1.3-Dithiolanes
		Dimethyl Acetals and Ketals
		N,N-Dimethylhydrazones

**Reference:** DOI: [10.1021/ol990629a](#)

Retrosynthesis ID: 11476

## 2.3 Path 3

Score: 2000076.25

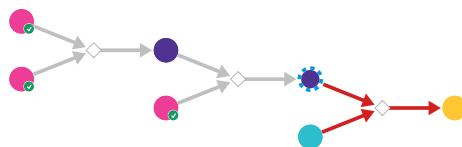
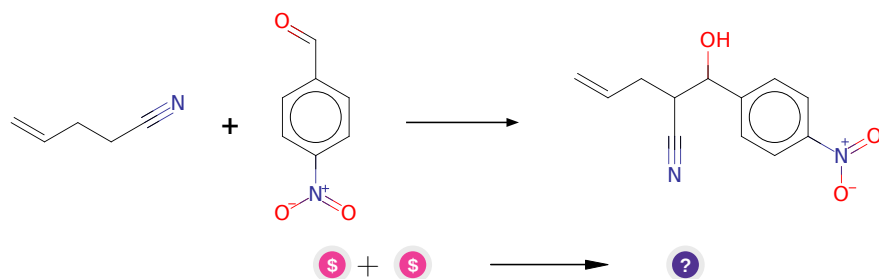


Figure 3: Outline of path 3

### 2.3.1 Aldol-like condensation with nitriles



**Substrates:**

1. 4-Nitrobenzaldehyde - *available at Sigma-Aldrich*
2. 4-Pentenitrile - *available at Sigma-Aldrich*

**Products:**

1. C=CCC(C#N)C(O)c1ccc([N+](=O)[O-])cc1

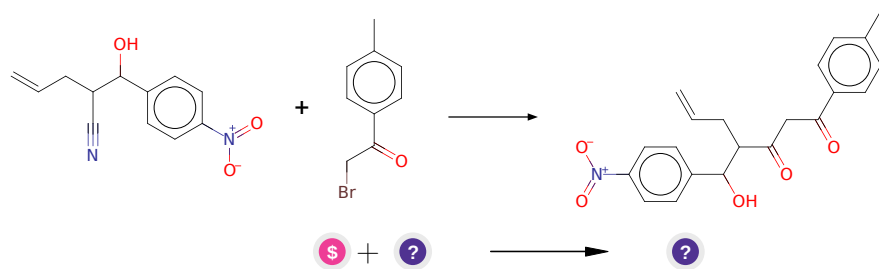
**Typical conditions:** LDA.THF.cooling

**Protections:** none

**Reference:** [10.1039/B800634B](#) and [10.1002/anie.201302613](#) and [10.1021/jm701319c](#) and [10.1016/S0040-4020\(98\)00122-7](#) and [10.1021/jo025872t](#)

Retrosynthesis ID: 23727

### 2.3.2 Blaise Reaction



#### Substrates:

1. 2-Bromo-4'-methylacetophenone - *available at Sigma-Aldrich*
2. C=CCC(C#N)C(O)c1ccc([N+](=O)[O-])cc1

#### Products:

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1

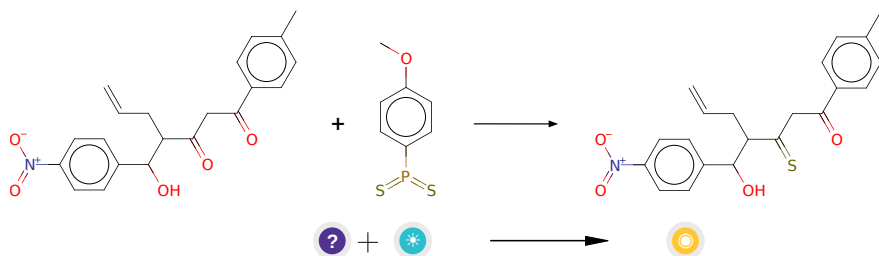
**Typical conditions:** Zn.TMSCl.THF then HCl

**Protections:** none

**Reference:** *10.1002/ejoc.201403402*

**Retrosynthesis ID:** 10000153

### 2.3.3 Synthesis of Thioketones using Lawesson's Reagent



#### Substrates:

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1
2. 4-methoxyphenyl-dithiophosphonsaeureanhydrid

#### Products:

1. C=CCC(C(=S)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1



**Typical conditions:** Lawesson's Reagent.neat.microwave

**Protections:**

Functional group SMARTS	Classification	Protecting groups
<chem>[*]C([*])=O</chem>	carbonyls	1.3-Dioxanes 1.3-Dioxolanes 1.3-Dithianes 1.3-Dithiolanes Dimethyl Acetals and Ketals N,N-Dimethylhydrazones

Reference: DOI: [10.1021/ol990629a](https://doi.org/10.1021/ol990629a)

Retrosynthesis ID: 11476

## 2.4 Path 4

Score: 2000090.31

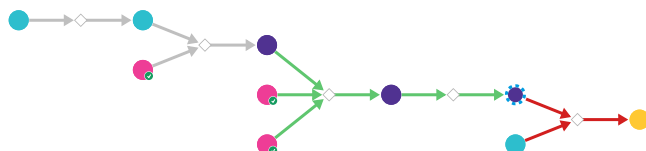
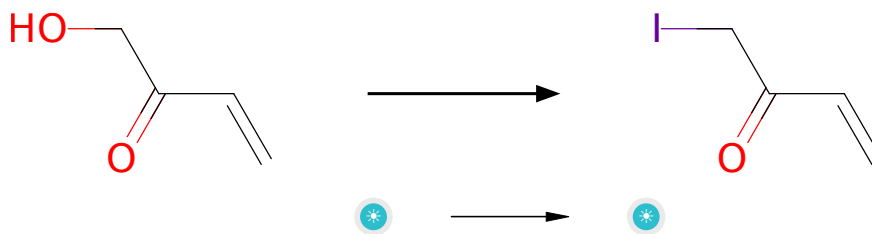


Figure 4: Outline of path 4

### 2.4.1 Synthesis Of Alkyl Iodides Via Appel Reaction



Substrates:

- 1-hydroxy-but-3-en-2-one

Products:

- 1-iodo-but-3-en-2-one

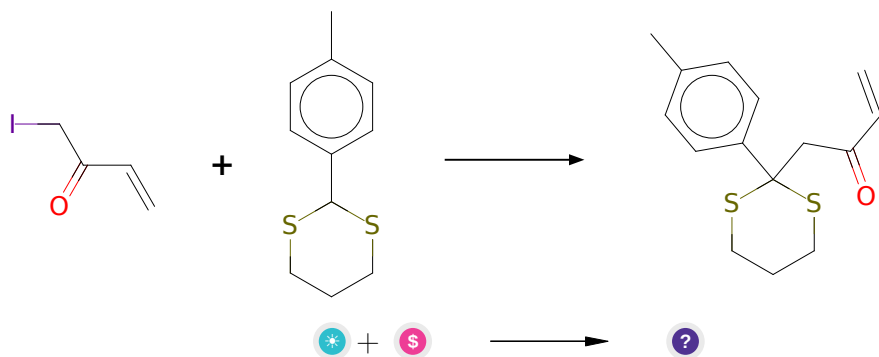
**Typical conditions:** Imidazole.PPh<sub>3</sub>.I<sub>2</sub>

**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

#### 2.4.2 Alkylation of dithianes



**Substrates:**

- 1-iodo-but-3-en-2-one
- 2-p-tolyl-[1,3]dithiane - *available at Sigma-Aldrich*

**Products:**

- C=CC(=O)CC1(c2ccc(C)cc2)SCCS1

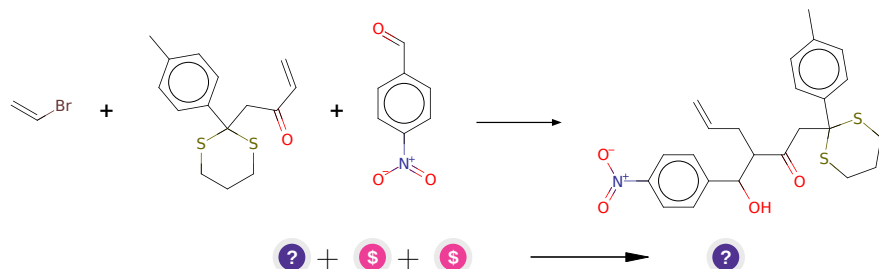
**Typical conditions:** LDA.THF

**Protections:** none

**Reference:** [10.1021/ja055740s](#) (SI) and [10.1016/S0008-6215\(99\)00275-X](#) and [10.1021/ja0618954](#)

**Retrosynthesis ID:** 34220

### 2.4.3 Alkenylation-Aldol reaction of enones and enoate esters



#### Substrates:

1. C=CC(=O)CC1(c2ccc(C)cc2)SCCCS1
2. 4-Nitrobenzaldehyde - *available at Sigma-Aldrich*
3. Bromoethylene - *available at Sigma-Aldrich*

#### Products:

1. C=CCC(C(=O)CC1(c2ccc(C)cc2)SCCCS1)C(O)c1ccc([N+](=O)[O-])cc1

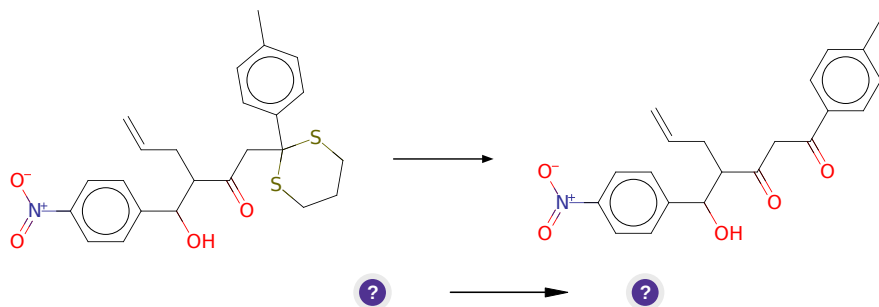
**Typical conditions:** 1.RCuLi.2.RCHO

**Protections:** none

**Reference:** [10.1016/S0040-4039\(01\)80891-1](#) AND [10.1016/S0040-4020\(01\)82115-3](#) AND [10.1021/jo2010186](#) AND [10.1021/jo101439h](#) AND [10.1021/ja906241w](#)

**Retrosynthesis ID:** 20547

### 2.4.4 Synthesis of ketones from dithianes



#### Substrates:

1. C=CCC(C(=O)CC1(c2ccc(C)cc2)SCCCS1)C(O)c1ccc([N+](=O)[O-])cc1

#### Products:

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1

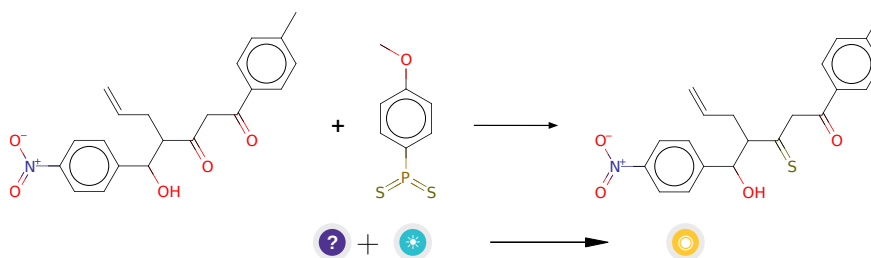
**Typical conditions:** MeI.CaCO<sub>3</sub>

**Protections:** none

**Reference:** [10.1016/j.tet.2013.09.075](https://doi.org/10.1016/j.tet.2013.09.075) and [10.1021/jo00007a015](https://doi.org/10.1021/jo00007a015) and [10.1021/jo0610412](https://doi.org/10.1021/jo0610412) and [10.1021/ol901024t](https://doi.org/10.1021/ol901024t) and [10.1021/ol500553x](https://doi.org/10.1021/ol500553x) and [10.1021/jo0626459](https://doi.org/10.1021/jo0626459)

**Retrosynthesis ID:** 31724

#### 2.4.5 Synthesis of Thioketones using Lawesson's Reagent



**Substrates:**

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1
2. 4-methoxyphenyl-dithiophosphonane

**Products:**

1. C=CCC(C(=S)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1

**Typical conditions:** Lawesson's Reagent.neat.microwave

**Protections:**

Functional group SMARTS	Classification	Protecting groups
[#6]C([#6])=O	carbonyls	1.3-Dioxanes
		1.3-Dioxolanes
		1.3-Dithianes
		1.3-Dithiolanes
		Dimethyl Acetals and Ketals
		N,N-Dimethylhydrazones

Reference: DOI: [10.1021/ol990629a](https://doi.org/10.1021/ol990629a)

Retrosynthesis ID: 11476

## 2.5 Path 5

Score: 2000115.31

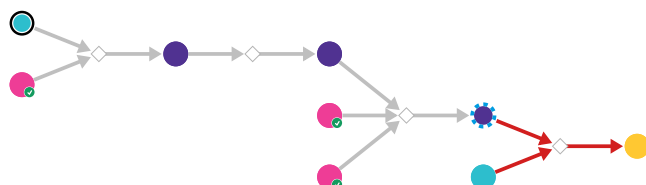
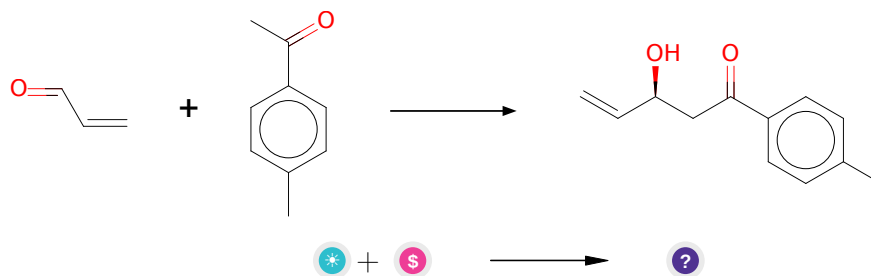


Figure 5: Outline of path 5

### 2.5.1 Asymmetric Aldol Reaction



Substrates:

1. Acrolein
2. Methyl p-tolyl ketone - [available at Sigma-Aldrich](#)

Products:

1. C=C[C@@H](O)CC(=O)c1ccc(C)cc1

Typical conditions: chiral catalyst

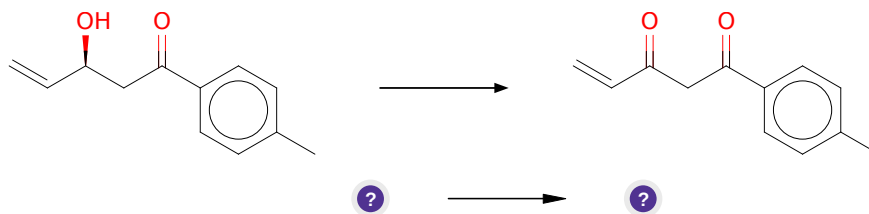
Protections: none

Reference: [10.1016/S0040-4039\(00\)95300-0](https://doi.org/10.1016/S0040-4039(00)95300-0) and [10.1016/S0040-4039\(00\)80089-1](https://doi.org/10.1016/S0040-4039(00)80089-1) and [10.1016/j.tetlet.2006.08.051](https://doi.org/10.1016/j.tetlet.2006.08.051) and [10.5012/bkcs.2010.31.03.653](https://doi.org/10.5012/bkcs.2010.31.03.653) and

10.1002/(SICI)1521-3773(19991216)38:24<3738::AID-ANIE3738>3.0.CO;2-2  
and 10.1039/B923537J

Retrosynthesis ID: 9991987

### 2.5.2 Oxidation of Chiral Alcohols



Substrates:

1. C=C[C@@H](O)CC(=O)c1ccc(C)cc1

Products:

1. C=CC(=O)CC(=O)c1ccc(C)cc1

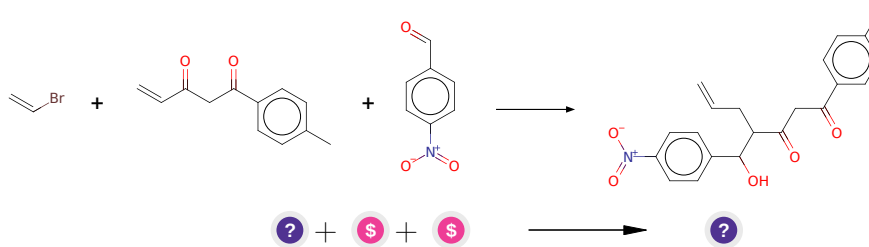
Typical conditions: CAN.NaBrO<sub>3</sub>.CH<sub>3</sub>CN or other oxidant e.g. TPAP or NaOCl

Protections: none

Reference: DOI: 10.1016/S0040-4039(00)86883-5 or 10.1021/ja00054a005 or 10.1016/S0040-4039(00)85677-4

Retrosynthesis ID: 25121

### 2.5.3 Alkenylation-Aldol reaction of enones and enoate esters



Substrates:

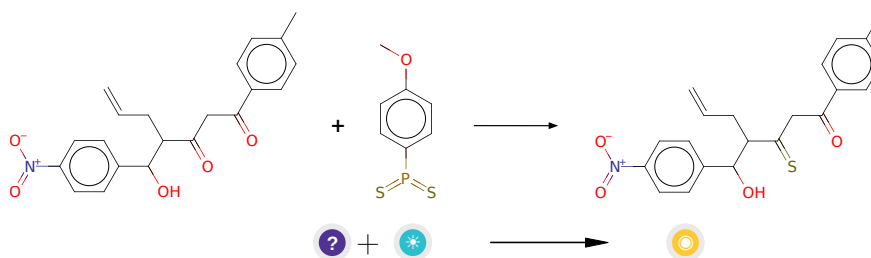
1. C=CC(=O)CC(=O)c1ccc(C)cc1
2. Bromoethylene - *available at Sigma-Aldrich*
3. 4-Nitrobenzaldehyde - *available at Sigma-Aldrich*

**Products:**

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1

**Typical conditions:** 1.RCuLi.2.RCHO**Protections:** none

**Reference:** [10.1021/jo2010186](#) AND [10.1021/jo101439h](#) AND [10.1021/ja906241w](#)  
 AND [10.1016/S0040-4039\(01\)80891-1](#) AND [10.1016/S0040-4020\(01\)82115-3](#)

**Retrosynthesis ID:** 13048**2.5.4 Synthesis of Thioketones using Lawesson's Reagent****Substrates:**

1. C=CCC(C(=O)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1
2. 4-methoxyphenyl-dithiophosphonane

**Products:**

1. C=CCC(C(=S)CC(=O)c1ccc(C)cc1)C(O)c1ccc([N+](=O)[O-])cc1

**Typical conditions:** Lawesson's Reagent.neat.microwave**Protections:**

Functional group SMARTS	Classification	Protecting groups
<chem>[*]C([*])=O</chem>	carbonyls	1.3-Dioxanes 1.3-Dioxolanes 1.3-Dithianes 1.3-Dithiolanes Dimethyl Acetals and Ketals N,N-Dimethylhydrazones



**Reference:** DOI: [10.1021/ol990629a](https://doi.org/10.1021/ol990629a)

**Retrosynthesis ID:** 11476