

# Paths of analysis\*

L5\_DIA

Synthia

October 11, 2022

## 1 Analysis parameters

**Analysis type:** Automatic Retrosynthesis

**Rules:** none selected

**Filters:** 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

**Strategies:** none selected

---

\*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.

**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:** 76.25

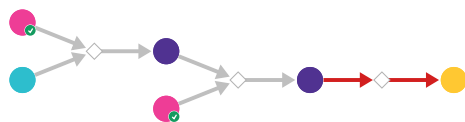
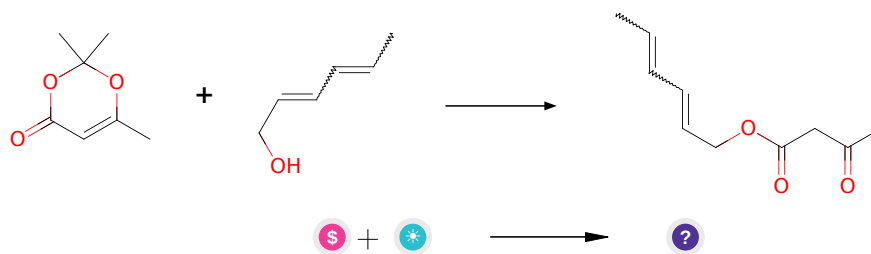


Figure 1: Outline of path 1

#### 2.1.1 Synthesis of 1,3-dicarbonyl compounds from 1,3-dioxinones



**Substrates:**

1. Diketene acetone adduct - *available at Sigma-Aldrich*
2. sorbic alcohol

**Products:**

1. CC=CC=CCOC(=O)CC(C)=O

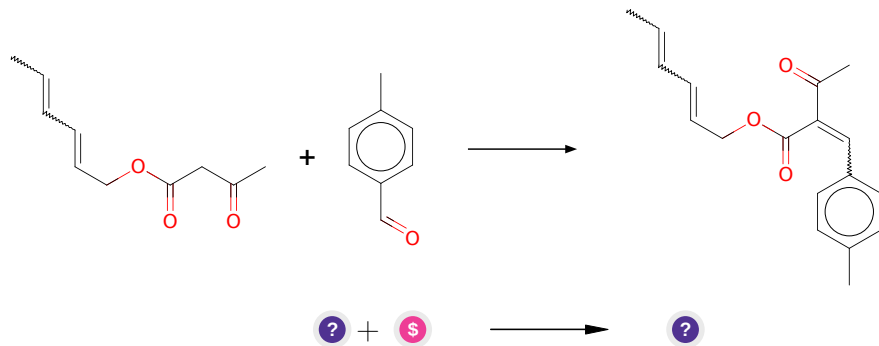
**Typical conditions:** alcohol

**Protections:** none

**Reference:** DOI: [10.1021/ja00154a049](https://doi.org/10.1021/ja00154a049)

**Retrosynthesis ID:** 12398

### 2.1.2 Knoevenagel Condensation



#### Substrates:

1. CC=CC=CCOC(=O)CC(C)=O
2. p-Tolualdehyde - *available at Sigma-Aldrich*

#### Products:

1. CC=CC=CCOC(=O)C(=Cc1ccc(C)cc1)C(C)=O

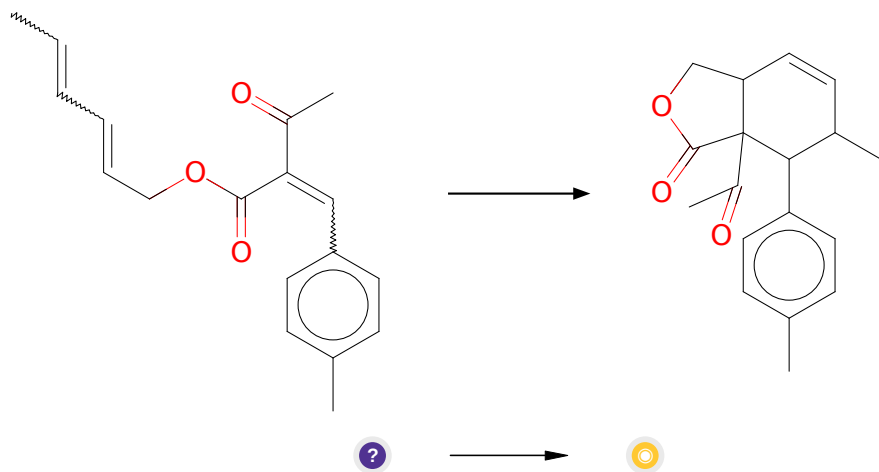
**Typical conditions:** base e.g.piperidine. solvent

**Protections:** none

**Reference:** *10.1002/0471264180.or015.02* and *10.13005/ojc/350154*

**Retrosynthesis ID:** 252

### 2.1.3 Diels-Alder



#### Substrates:

1. CC=CC=CCOC(=O)C(=Cc1ccc(C)cc1)C(C)=O

**Products:**

1. CC(=O)C12C(=O)OCC1C=CC(C)C2c1ccc(C)cc1

**Typical conditions:** Lewis acid or chiral Lewis acid. Solvent.

**Protections:** none

**Reference:** DOI: [10.1002/1521-3773\(20020517\)41:10<1668::AID-ANIE1668>3.0.CO;2-Z](https://doi.org/10.1002/1521-3773(20020517)41:10<1668::AID-ANIE1668>3.0.CO;2-Z) AND [10.1021/ja062508t](https://doi.org/10.1021/ja062508t)

**Retrosynthesis ID:** 18116

## 2.2 Path 2

Score: 76.25

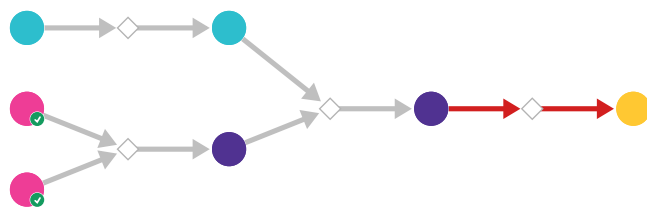
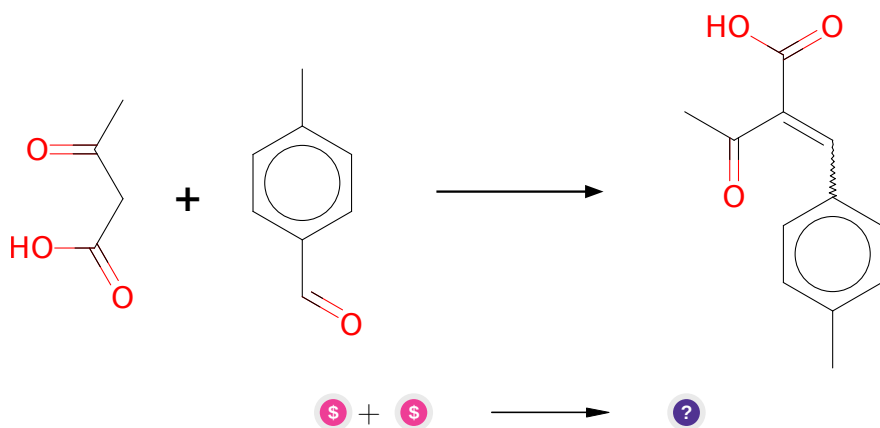


Figure 2: Outline of path 2

### 2.2.1 Knoevenagel Condensation



**Substrates:**

1. p-Tolualdehyde - *available at Sigma-Aldrich*
2. Lithium acetoacetate - *available at Sigma-Aldrich*

**Products:**

1. CC(=O)C(=Cc1ccc(C)cc1)C(=O)O

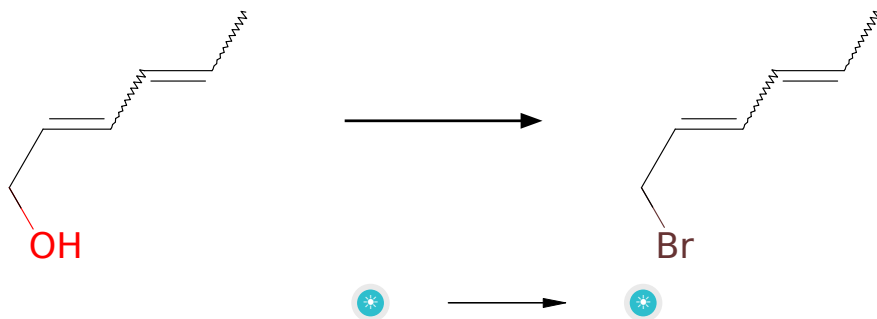
**Typical conditions:** base e.g.piperidine. solvent

**Protections:** none

**Reference:** *10.1002/0471264180.or015.02* and *10.13005/ojc/350154*

**Retrosynthesis ID:** 252

**2.2.2 Appel Reaction**



**Substrates:**

1. sorbic alcohol

**Products:**

1. 1-brom-hexa-2,4-dien

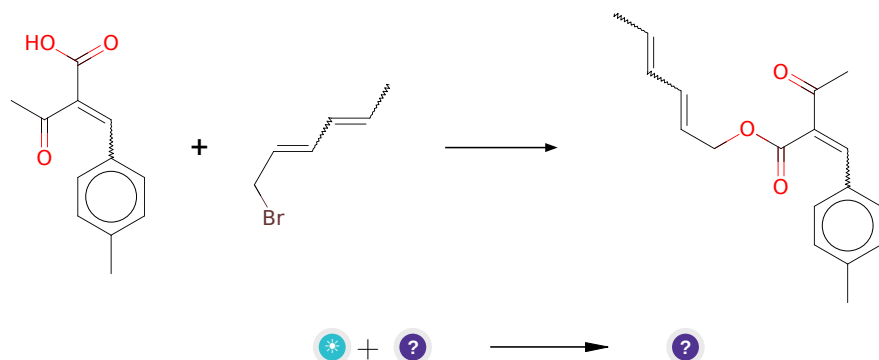
**Typical conditions:** PPh<sub>3</sub>.CBr<sub>4</sub>

**Protections:** none

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

**Retrosynthesis ID:** 9990037

### 2.2.3 Synthesis of esters from alkyl chlorides and carboxylic acids or thioacids



#### Substrates:

- 1-brom-hexa-2,4-dien
- CC(=O)C(=Cc1ccc(C)cc1)C(=O)O

#### Products:

- CC=CC=CCOC(=O)C(=Cc1ccc(C)cc1)C(C)=O

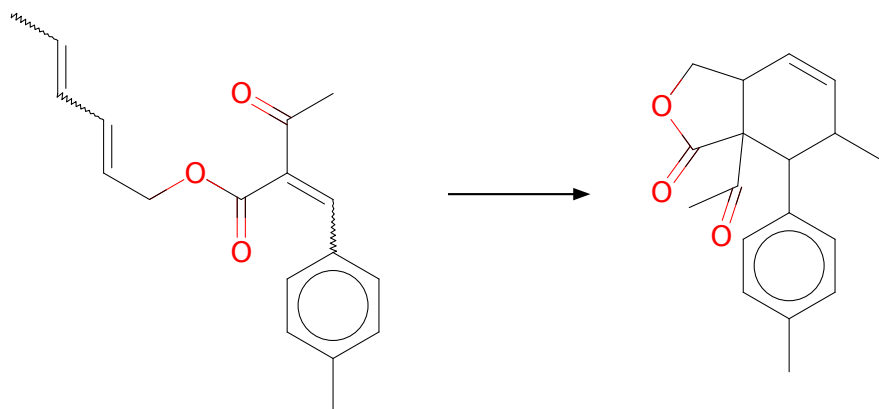
Typical conditions: K<sub>2</sub>CO<sub>3</sub>.DMF

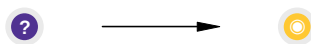
Protections: none

Reference: [10.1016/j.bmcl.2005.08.026](#) AND [10.1021/ol034655r](#) (SI) AND [10.1039/C3RA41967C](#) AND [10.1016/j.bmcl.2012.03.093](#)

Retrosynthesis ID: 14685

### 2.2.4 Diels-Alder





**Substrates:**

1. CC=CC=CCOC(=O)C(=Cc1ccc(C)cc1)C(C)=O

**Products:**

1. CC(=O)C12C(=O)OCC1C=CC(C)C2c1ccc(C)cc1

**Typical conditions:** Lewis acid or chiral Lewis acid. Solvent.

**Protections:** none

**Reference:** DOI: [10.1002/1521-3773\(20020517\)41:10<1668::AID-ANIE1668>3.0.CO;2-Z](https://doi.org/10.1002/1521-3773(20020517)41:10<1668::AID-ANIE1668>3.0.CO;2-Z) AND [10.1021/ja062508t](https://doi.org/10.1021/ja062508t)

**Retrosynthesis ID:** 18116

## 2.3 Path 3

**Score:** 76.25

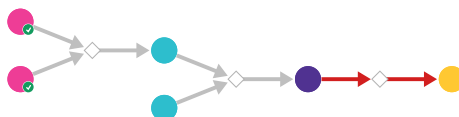
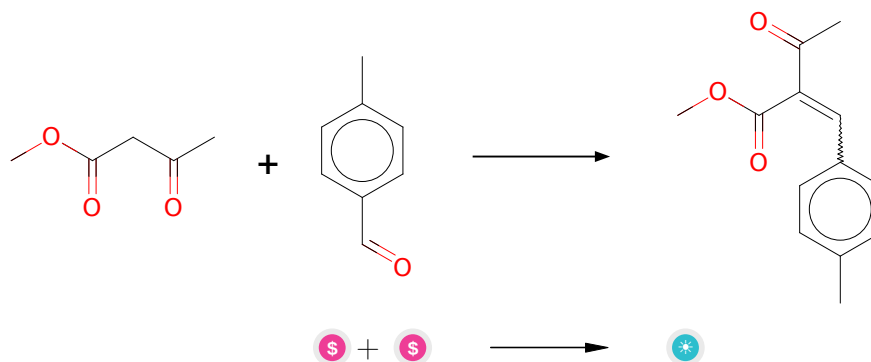


Figure 3: Outline of path 3

### 2.3.1 Knoevenagel Condensation



**Substrates:**

1. Methyl acetoacetate - *available at Sigma-Aldrich*
2. p-Tolualdehyde - *available at Sigma-Aldrich*

**Products:**

1. 2-acetyl-3-p-tolyl-acrylic acid methyl ester

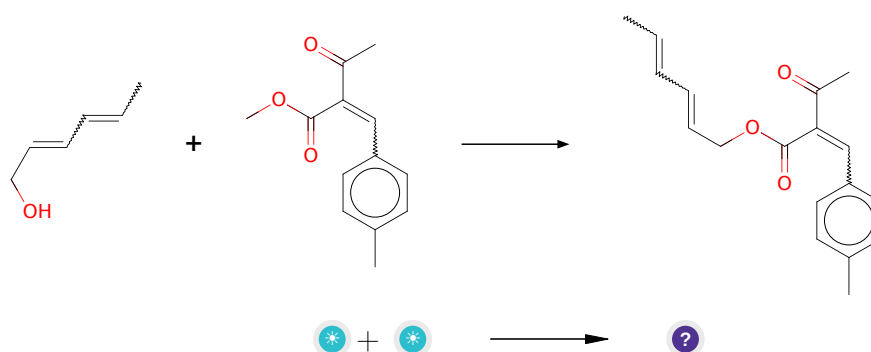
**Typical conditions:** base e.g.piperidine. solvent

**Protections:** none

**Reference:** *10.1002/0471264180.or015.02* and *10.13005/ojc/350154*

**Retrosynthesis ID:** 252

### 2.3.2 Acid catalyzed transesterification



**Substrates:**

1. sorbic alcohol
2. 2-acetyl-3-p-tolyl-acrylic acid methyl ester

**Products:**

1. CC=CC=CCOC(=O)C(=Cc1ccc(C)cc1)C(C)=O

**Typical conditions:** H+

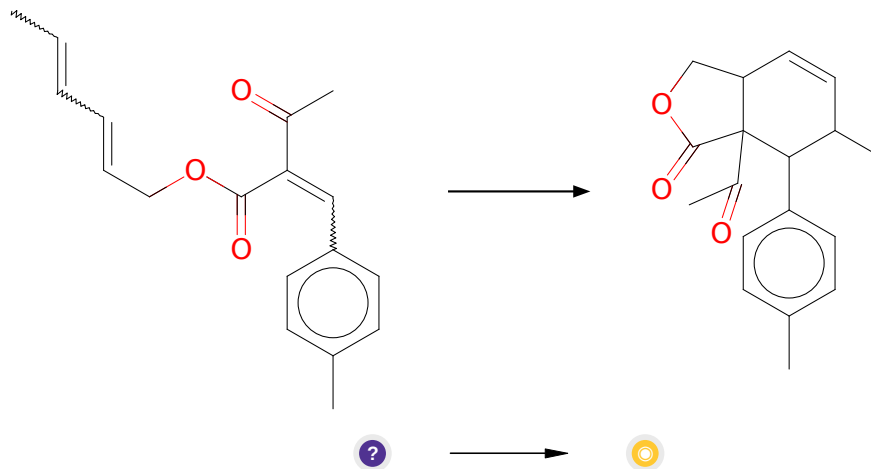
**Protections:** none

**Reference:** *10.1021/cr00020a004*

**Retrosynthesis ID:** 50438



### 2.3.3 Diels-Alder



**Substrates:**

1. CC=CC=CCOC(=O)C(=Cc1ccc(C)cc1)C(C)=O

**Products:**

1. CC(=O)C12C(=O)OCC1C=CC(C)C2c1ccc(C)cc1

**Typical conditions:** Lewis acid or chiral Lewis acid. Solvent.

**Protections:** none

**Reference:** DOI: [10.1002/1521-3773\(20020517\)41:10<1668::AID-ANIE1668>3.0.CO;2-Z](https://doi.org/10.1002/1521-3773(20020517)41:10<1668::AID-ANIE1668>3.0.CO;2-Z) AND [10.1021/ja062508t](https://doi.org/10.1021/ja062508t)

**Retrosynthesis ID:** 18116

### 2.4 Path 4

**Score:** 76.25

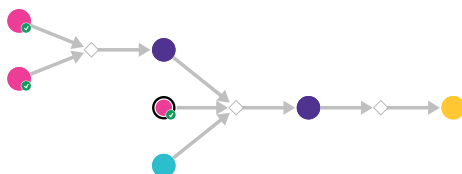
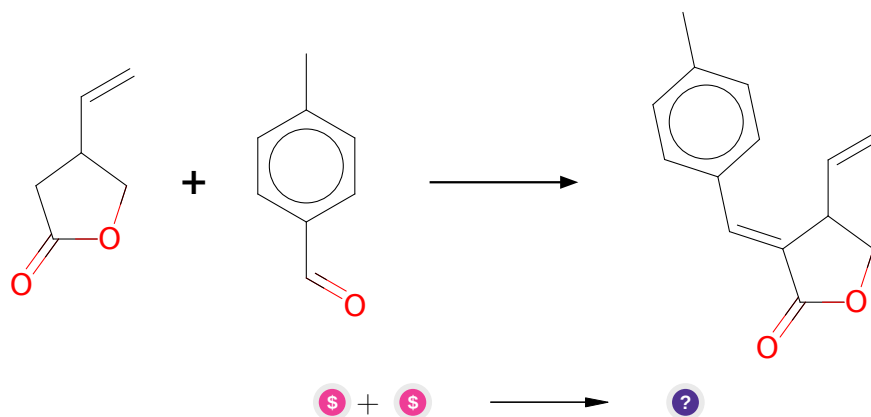


Figure 4: Outline of path 4

#### 2.4.1 Condensation of esters with aldehydes



##### Substrates:

1. 4-ethenyloxolan-2-one - *available at Sigma-Aldrich*
2. p-Tolualdehyde - *available at Sigma-Aldrich*

##### Products:

1. C=CC1COC(=O)/C1=C/c1ccc(C)cc1

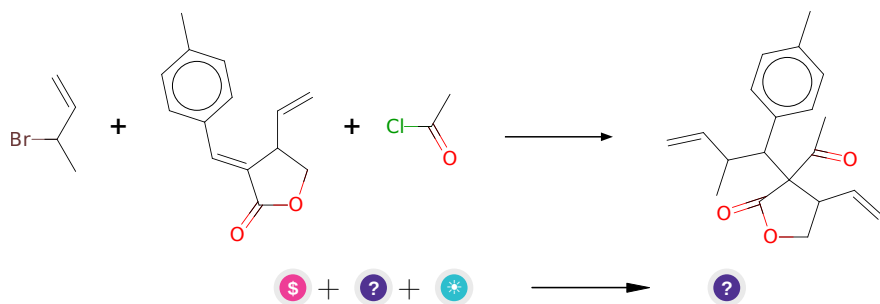
**Typical conditions:** 1.LDA.2RCHO

**Protections:** none

**Reference:** [10.1021/jo970387x](#) AND [10.1021/jo00076a051](#) AND [10.1016/S0040-4039\(97\)10827-9](#) AND [10.1055/s-2002-25767](#) AND [10.1039/P19920003277](#)

**Retrosynthesis ID:** 14981

## 2.4.2 Conjugated addition of organocuprate-acylation of enones and enoate esters



### Substrates:

1. Acetyl chloride - *available at Sigma-Aldrich*
2. C=CC1COC(=O)/C1=C/c1ccc(C)cc1
3. 3-brom-but-1-en

### Products:

1. C=CC(C)C(c1ccc(C)cc1)C1(C(C)=O)C(=O)OCC1C=C

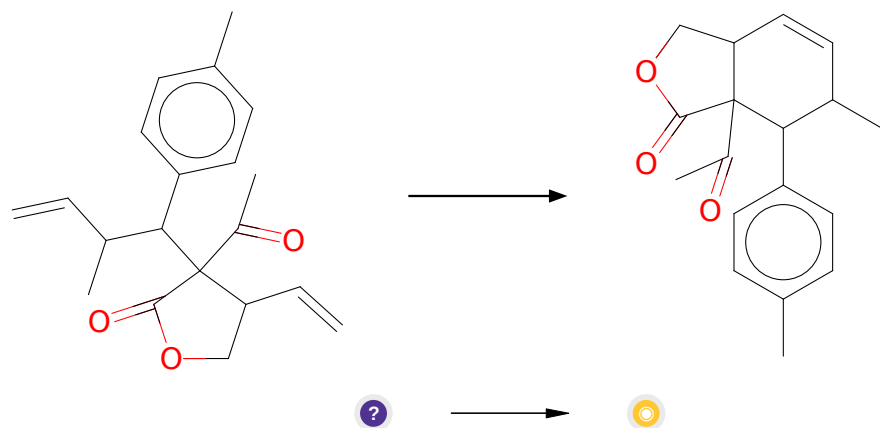
**Typical conditions:** 1.RCuLi.2.AcCl.HMPA

**Protections:** none

**Reference:** [10.3987/COM-99-S143](#) AND [10.1021/ja00148a023](#) AND [10.1016/S0040-4039\(01\)80891-1](#)

**Retrosynthesis ID:** 20523

## 2.4.3 Ring-Closing Metathesis



**Substrates:**

1. C=CC(C)C(c1ccc(C)cc1)C1(C(C)=O)C(=O)OCC1C=C

**Products:**

1. CC(=O)C12C(=O)OCC1C=CC(C)C2c1ccc(C)cc1

**Typical conditions:** catalyst e.g. Hoveyda-Grubbs . solvent e.g. CH<sub>2</sub>Cl<sub>2</sub>

**Protections:** none

**Reference:** DOI: [10.1002/anie.200800693](https://doi.org/10.1002/anie.200800693) and [10.1021/acs.orglett.8b04003](https://doi.org/10.1021/acs.orglett.8b04003) and [10.1021/jo0264729](https://doi.org/10.1021/jo0264729) and [10.1021/ja072334v](https://doi.org/10.1021/ja072334v) and [10.1002/ejoc.201001102](https://doi.org/10.1002/ejoc.201001102)

**Retrosynthesis ID:** 31014187

## 2.5 Path 5

Score: 76.25

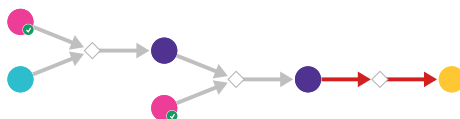
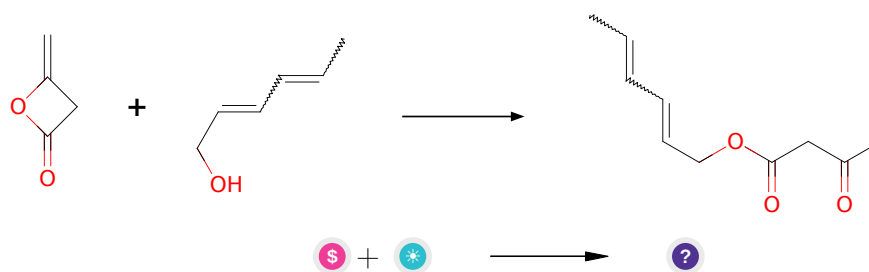


Figure 5: Outline of path 5

### 2.5.1 Reaction of alcohols with diketene



**Substrates:**

1. diketene - *available at Sigma-Aldrich*
2. sorbic alcohol

**Products:**

1. CC=CC=CCOC(=O)CC(C)=O

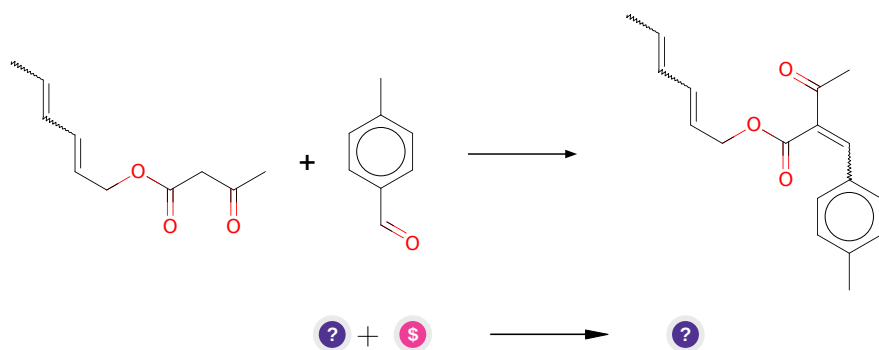
**Typical conditions:** DCM.heat

**Protections:** none

**Reference:** WO2012/31028 A2 (p.39) AND [10.1021/ol051945u](#) AND [10.1021/ol0069756](#) AND [10.1002/adsc.200800532](#)

**Retrosynthesis ID:** 14881

### 2.5.2 Knoevenagel Condensation



**Substrates:**

1. CC=CC=CCOC(=O)CC(C)=O
2. p-Tolualdehyde - [available at Sigma-Aldrich](#)

**Products:**

1. CC=CC=CCOC(=O)C(=Cc1ccc(C)cc1)C(C)=O

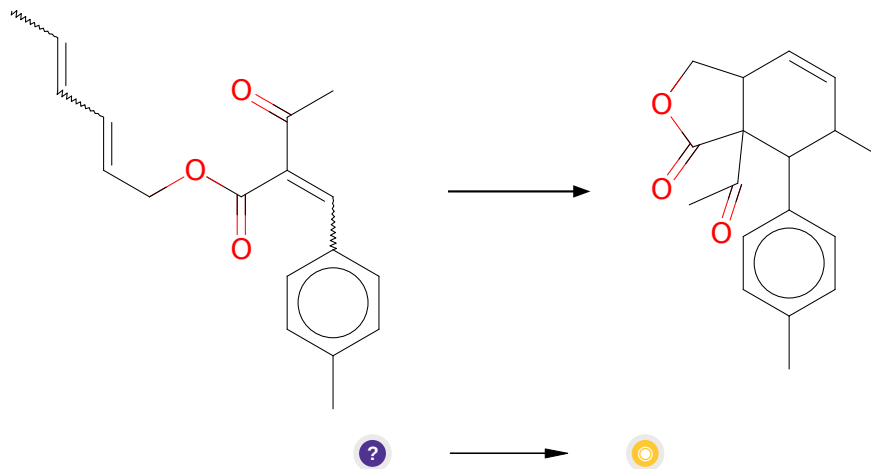
**Typical conditions:** base e.g.piperidine. solvent

**Protections:** none

**Reference:** [10.1002/0471264180.or015.02](#) and [10.13005/ojc/350154](#)

**Retrosynthesis ID:** 252

### 2.5.3 Diels-Alder



**Substrates:**

1. CC=CC=CCOC(=O)C(=Cc1ccc(C)cc1)C(C)=O

**Products:**

1. CC(=O)C12C(=O)OCC1C=CC(C)C2c1ccc(C)cc1

**Typical conditions:** Lewis acid or chiral Lewis acid. Solvent.

**Protections:** none

**Reference:** DOI: [10.1002/1521-3773\(20020517\)41:10<1668::AID-ANIE1668>3.0.CO;2-Z](https://doi.org/10.1002/1521-3773(20020517)41:10<1668::AID-ANIE1668>3.0.CO;2-Z) AND [10.1021/ja062508t](https://doi.org/10.1021/ja062508t)

**Retrosynthesis ID:** 18116