

SYNTHIA™ Retrosynthesis Software

DISCOVER Greener pathways

MilliporeSigma is the U.S. and Canada Life Science business of Merck KGaA, Darmstadt, Germany.

the path to a Greener Future

Designing an efficient process for synthesizing a target compound with minimal environmental impact presents numerous challenges. Our powerful retrosynthesis software, SYNTHIA[™], can greatly facilitate prediction of more sustainable synthetic routes, starting from commercially available raw materials.

Learn how our advanced algorithms help you optimize pathway design based on price, step count, and atom economy, leading to more efficient, economical, and ecological routes.

Sustainability Information on Building Blocks

Building blocks that support Green Chemistry are clearly labeled in pathways.



Greener Alternative Product High standards, environmental goals, and ecosystem partnerships.





Spend-based emission factor Reported in kg CO₂/EUR and according to the Environmentally Extended Input-Output (EEIO) model.



EPA Safer Chemical Ingredients List Includes chemicals evaluated through the Safer Choice Program



For more information, please visit: WWW.SYNTHIAONLINE.COM



Building block details with one click:

- Link to PubChem
- Chemical Hazard Symbols
- Greener Alternative Product Symbol
- EPA Safer Chemical Ingredients Labels
- Spend-based emission factor in kgCO₂ eq/EUR

Greener Alternative Products



This symbol indicates that the compound is part of our Greener Alternative Product portfolio and belongs to one of the 4 categories shown below.

4 Categories of Greener Alternative Products



Re-Engineered Products are developed by our scientists to significantly reduce environmental impact



Enabling Products help make greener alternatives possible through new technologies



12 Principles Aligned Products follow at least one of the 12 Principles of Green Chemistry



Design for Sustainability (DfS) Developed Products demonstrate significant sustainability improvements

Enzymatic Reactions

Enzymatic reactions are found in our expert-coded "Enzymatic Transformations" database. Reaction information includes reported conditions and enzyme details along with reference links.

Benefit

Another benefit is reaction conditions for enzymatic transformations are often greener themselves - not using harsh chemical reagents or solvents required in other processes.



Shared Path Library

Shared Path Library in SYNTHIA[™] allows you to plan a synthetic pathway for an entire library of target molecules simultaneously via multi-target analysis.

Simply upload a library of target molecules you want to synthesize, and SYNTHIA[™] will identify and prioritize synthetic routes that have the highest number of common intermediates and starting materials.

- Multi-target analysis for efficient, autonomous retrosynthetic analysis of entire target library
- Prioritization of routes with shared synthetic steps

This functionality enables quick prediction of the optimal strategy for synthesizing an entire library of target molecules using shared synthetic steps.



Customize for Sustainable Pathways

Requirements for synthesis projects can differ greatly. That's why SYNTHIA[™] offers numerous customization options for adjusting results according to your project's needs. This is especially important in the context of sustainable chemistry.

SYNTHIA[™] offers limitless options for customization:

- Avoid using gaseous reagents/reactants
- Eliminate metal catalysis
- Limit the number of synthesis steps
- Minimize the cost of starting materials
- Specify protecting group requirements
- Define bond disconnections
- Exclude toxic or hazardous molecules
- Select catalyst/reagents type
- Promote or exclude specific chemistry reaction classes, such as enzyme-catalyzed reactions
- Choose starting materials: commercial, known, or internal inventory

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0.04	908.70
-+-	
0.04	908.7
Pathway Similarity (i)	
-	(10
10	10
Number of Reactions in Path	
1 1	
1	2
Protecting Groups	
0	5
0	

YOUR BENEFITS

- Reduce risk of overlooking a more sustainable pathway
- Limit use of hazardous reactants/reagents
- Save time finding the best synthetic route, including for target molecule libraries
- Find pathways that meet your synthetic needs
- Quickly filter molecule sets based on synthesizability, or connect with other tools using SYNTHIA[™] API

12 principles of Green Chemistry



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MS_FL5674EN Ver. 1.0 01/2025