



Data Sheet

Product Name: Repaglinide
Cat. No.: CS-0979
CAS No.: 135062-02-1
Molecular Formula: C27H36N2O4

Molecular Weight: 452.59
Target: Others
Pathway: Others

Solubility: DMSO : \geq 50 mg/mL (110.48 mM)

BIOLOGICAL ACTIVITY:

Repaglinide is an **insulin secretagogue** for the treatment of type-2 diabetes mellitus^[1]. **In Vitro**: Repaglinide reduces postprandial glucose levels by enhancing the early phase of insulin secretion and increasing the total amount of insulin secreted^[1]. **In Vivo**: Repaglinide (AG-EE 623ZW) is very rapidly absorbed (t_{max} less than 1 hour) with a $t_{1/2}$ of less than one hour. Furthermore, Repaglinide is inactivated in the liver and more than 90 % excreted via the bile. Repaglinide (1 mg/kg p.o.) is effective (P<0.001) as an insulin-releasing agent in a rat model (low-dose streptozotocin) of type 2 diabetes.

References:

[1]. Wang LC, et al. Characteristics of repaglinide and its mechanism of action on insulin secretion in patients with newly diagnosed type-2 diabetes mellitus. Medicine (Baltimore). 2018 Sep;97(38):e12476.

CAIndexNames:

 $Benzoic\ acid,\ 2-ethoxy-4-[2-[[(1S)-3-methyl-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino]-2-oxoethyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]-1-[2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl]amino[-2-(1-piperidinyl)phenyl]butyl[-2-(1-piperidinyl)phenyl]butyl[-2-(1-piperidinyl)phenyl]butyl[-2-(1-piperidinyl)phenyl]butyl[-2-(1-piperidinyl)phenyl]butyl[-2-(1-piperidinyl)phenyl[-2-(1-piperidinyl)phe$

SMILES:

CC(C)C[C@H](NC(CC1=CC(OCC)=C(C(O)=O)C=C1)=O)C2=CC=CC=C2N3CCCC3

Caution: Product has not been fully validated for medical applications. For research use only.

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