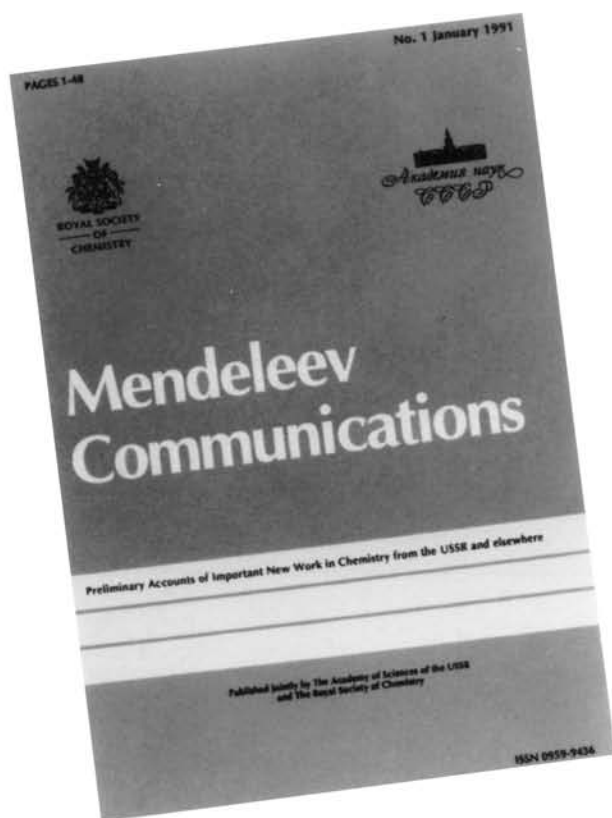


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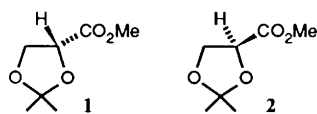


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Chiral Building Blocks

Aldrich lists a variety of building blocks for asymmetric synthesis. Listed on this page are just a few of the many useful products in this category. Please consult the 1990-1991 Aldrich Catalog/Handbook for additional chiral molecules.

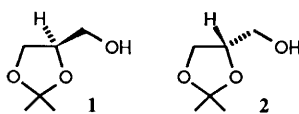


Optically active precursors for *O*-protected α -hydroxy ketones.¹ Chiral synthons for the preparation of terpenes.²

(1) Dumont, R.; Pfander, H. *Helv. Chim. Acta* **1983**, *66*, 814.
(2) Larcheveque, M.; Petit, Y. *Synthesis* **1986**, 60.

34,548-2 Methyl (*R*)-(+)-2,2-dimethyl-1,3-dioxolane-4-carboxylate, 98% (1)
1g \$14.00; 5g \$44.20

25,460-6 Methyl (*S*)-(-)-2,2-dimethyl-1,3-dioxolane-4-carboxylate, 98% (2)
1g \$11.60; 5g \$39.10

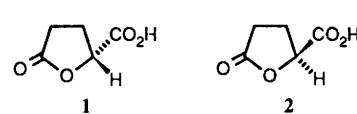


Important starting materials for the synthesis of chiral phospholipids.

Rosario-Jansen, T.; Jiang, R.-T.; Tsai, M.-D. *Biochemistry* **1988**, *27*, 4619.

24,180-6 (*R*)-(-)-2,2-Dimethyl-1,3-dioxolane-4-methanol, 98% (1) 250mg \$12.00
1g \$32.50

23,774-4 (*S*)-(+)-2,2-Dimethyl-1,3-dioxolane-4-methanol, 98% (2) 1g \$12.60

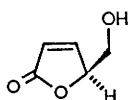


Useful precursors for many chiral synthons.¹ The (*S*)-isomer was used as a chiral template in the synthesis of (-)-steganone.²

(1) Ravid, U.; Silverstein, R.M.; Smith, L.R. *Tetrahedron Lett.* **1980**, *21*, 2709. (2) Larcheveque, M.; Petit, Y. *Synthesis* **1986**, 60.

31,047-6 (*R*)-(-)-5-Oxo-2-tetrahydrofuran-2-carboxylic acid, 98% (1) 1g \$12.50
5g \$40.40

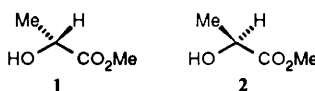
30,146-9 (*S*)-(+)-5-Oxo-2-tetrahydrofuran-2-carboxylic acid, 98% (2) 1g \$14.00
5g \$45.50



Key intermediate for the efficient total synthesis of the potent antiviral agent, 3'-azido-3'-deoxythymidine (AZT).¹ Also used as the starting material in Hanessian's replicating lactone strategy for acyclic stereocontrol.²

(1) Chu, C.K.; Beach, J.W.; Ullas, G.V.; Kosugi, Y. *Tetrahedron Lett.* **1988**, *29*, 5349. (2) Hanessian, S. *Aldrichim. Acta* **1989**, *22*, 3.

34,686-1 (*S*)-(-)-5-(Hydroxymethyl)-2(5*H*)-furanone, 98% 250mg \$16.20
1g \$45.30

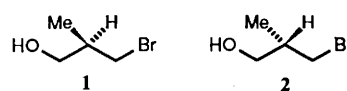


For the preparation of α -amino esters of high optical purity via displacement of their derived triflates.¹ The (*S*)-isomer was recently used in the first total synthesis of the redox coenzyme Factor 420 (F₄₂₀).²

(1) Feenstra, R.W.; Stokkingreef, E.H.M.; Nivard, R.J.F.; Ottenheijm, H.C.J. *Tetrahedron* **1988**, *44*, 5583. (2) Tanaka, K.; Kimachi, T.; Kawase, M.; Yoneda, F. *J. Chem. Soc., Chem. Commun.* **1988**, 524.

27,776-2 Methyl (*R*)-(+)-lactate, 98% (1) 1g \$10.50; 5g \$30.30

23,034-0 Methyl (*S*)-(-)-lactate, 98% (2) 5g \$10.70; 100g \$28.80

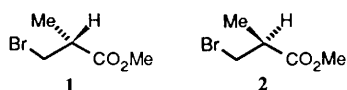


Reaction with triphenylphosphine produces interesting optically active α -hydroxyphosphonium salts from which the corresponding ylids can be generated.

Kozikowski, A.P.; Chen, Y.-Y.; Wang, B.C.; Xu, Z.-B. *Tetrahedron* **1984**, *40*, 2345.

32,506-6 (*R*)-(-)-3-Bromo-2-methyl-1-propanol, 97% (1) 1g \$11.90; 5g \$37.70

32,505-8 (*S*)-(+)-3-Bromo-2-methyl-1-propanol, 99% (2) 1g \$11.10; 5g \$33.10

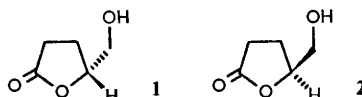


Precursors for optically active zinc homoenolates which react with electrophiles to yield chiral methyl esters.

Nakamura, E.; Sekiya, K.; Kuwajima, I. *Tetrahedron Lett.* **1987**, *28*, 337.

32,509-0 Methyl (*R*)-(+)-3-bromo-2-methylpropanoate, 97% (1) 1g \$10.75
5g \$35.70

32,492-2 Methyl (*S*)-(-)-3-bromo-2-methylpropanoate, 97% (2) 1g \$9.20
5g \$30.70

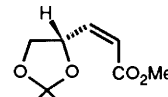


Chiral starting materials for enantioselective synthesis, for example, (+)-umbellactone and (*S*)-(+)-angelicalactone.^{1,2}

(1) Ortuno, R.M.; Alonso, D.; Cardellach, J.; Font, J. *Tetrahedron* **1987**, *43*, 2191. (2) Ortuno, R.M.; Bigorna, J.; Font, J. *ibid.* **1987**, *43*, 2199.

34,354-4 (*R*)-(-)-Dihydro-5-(hydroxymethyl)-2(3*H*)-furanone, 97% (1) 250mg \$16.20; 1g \$54.00

34,890-2 (*S*)-(+)-Dihydro-5-(hydroxymethyl)-2(3*H*)-furanone, 98% (2) 250mg \$16.20; 1g \$54.00



Valuable intermediate in stereocontrolled syntheses of *d*-pentitols.

Minami, N.; Ko, S.S.; Kishi, Y. *J. Am. Chem. Soc.* **1982**, *104*, 1109.

34,769-8 Methyl (*S*)-(+)-3-(2,2-dimethyl-1,3-dioxolan-4-yl)-*cis*-2-propenoate 1g \$20.50
5g \$68.00



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