

cis-Muurola-5-ene-4-«alpha»-ol

Inchi:	InChI=1S/C15H26O/c1-10(2)15(16)8-7-12(4)13-6-5-11(3)9-14(13)15/h9-13,16H,5-8H2,1H
InchiKey:	UAJHVFUBIKCODA-HNQLUHSGSA-N
Formula:	C15H26O
SMILES:	CC1C=C2C(CC1)C(C)CCC2(O)C(C)C
Mol. weight [g/mol]:	222.37

Physical Properties

Property code	Value	Unit	Source
hfus	19.72	kJ/mol	Joback Method
logp	3.776		Crippen Method
mcvol	202.060	ml/mol	McGowan Method
rinpol	1554.00		NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	597.80	J/mol×K	659.94	Joback Method
cpg	617.79	J/mol×K	694.13	Joback Method
cpg	636.77	J/mol×K	728.33	Joback Method
cpg	654.86	J/mol×K	762.52	Joback Method
cpg	672.18	J/mol×K	796.72	Joback Method
cpg	688.83	J/mol×K	830.91	Joback Method
cpg	704.93	J/mol×K	865.11	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R606996&Units=SI

Joback Method:

https://en.wikipedia.org/wiki/Joback_method

McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

Legend

cpg:	Ideal gas heat capacity
hfus:	Enthalpy of fusion at standard conditions
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rmpol:	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/71-398-2/cis-Muurola-5-ene-4-alpha-ol.pdf>

Generated by Cheméo on 2026-07-21 14:00:13.899210231 +0000 UTC m=+153509.741095626.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.