

3,3,7,11-TETRAMETHYLTRICYCLO[5.4.0.0(4,11)]UNDECAN-1-OL

Other names:	3,3,7,11-Tetramethyltricyclo[5.4.0.0(4,11)]undecan-1-ol
Inchi:	InChI=1S/C15H26O/c1-12(2)10-15(16)13(3)7-5-8-14(15,4)11(12)6-9-13/h11,16H,5-10H2
InchiKey:	QOXUIQMPPDIDGM-UHFFFAOYSA-N
Formula:	C15H26O
SMILES:	CC1(C)CC2(O)C3(C)CCCC2(C)C1CC3
Mol. weight [g/mol]:	222.37

Physical Properties

Property code	Value	Unit	Source
hfus	5.85	kJ/mol	Joback Method
logp	3.754		Crippen Method
mcvol	195.500	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	590.59	J/mol×K	655.16	Joback Method
cpg	610.07	J/mol×K	692.02	Joback Method
cpg	629.03	J/mol×K	728.88	Joback Method
cpg	647.94	J/mol×K	765.73	Joback Method
cpg	667.27	J/mol×K	802.59	Joback Method
cpg	687.47	J/mol×K	839.45	Joback Method
cpg	709.02	J/mol×K	876.31	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
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=C117591807&Units=SI>

Legend

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

Latest version available from:

<https://www.chemeo.com/cid/49-927-9/3-3-7-11-TETRAMETHYLTRICYCLO-5-4-0-0-4-11-UNDECAN-1-OL.pdf>

Generated by Cheméo on 2026-07-21 20:12:37.789283578 +0000 UTC m=+175853.631168986.

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