

(-)-Silphiperfolan-6«alpha»-ol

Inchi:	InChI=1S/C15H26O/c1-10-5-8-15-11(2)14(4,16)9-13(15,3)7-6-12(10)15/h10-12,16H,5-9H
InchiKey:	SNSNYEAITDGGIF-IRQRARHPSA-N
Formula:	C15H26O
SMILES:	CC1CCC23C1CCC2(C)CC(C)(O)C3C
Mol. weight [g/mol]:	222.37

Physical Properties

Property code	Value	Unit	Source
hfus	13.22	kJ/mol	Joback Method
logp	3.610		Crippen Method
mcvol	195.500	ml/mol	McGowan Method
rinpol	1501.00		NIST Webbook
ripol	1958.00		NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	595.75	J/mol×K	650.25	Joback Method
cpg	615.74	J/mol×K	685.82	Joback Method
cpg	634.97	J/mol×K	721.39	Joback Method
cpg	653.76	J/mol×K	756.96	Joback Method
cpg	672.45	J/mol×K	792.52	Joback Method
cpg	691.37	J/mol×K	828.09	Joback Method
cpg	710.84	J/mol×K	863.66	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=R327549&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
rinpol:	Non-polar retention indices
ripol:	Polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/70-336-1/Silphiperfolan-6-alpha-ol.pdf>

Generated by Cheméo on 2026-07-21 15:27:06.055184688 +0000 UTC m=+158721.897070080.

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