

«alpha»-Alaskene

Inchi:

InChI=1S/C15H24/c1-11(2)14-6-5-13(4)15(14)9-7-12(3)8-10-15/h7,13H,5-6,8-10H2,1-4H

InchiKey:

HMKLOOMRRZKSNM-AFYWWNPRSA-N

Formula:

C15H24

SMILES:

CC1=CCC2(CC1)C(=C(C)C)CCC2C

Mol. weight [g/mol]:

204.35

Physical Properties

Property code	Value	Unit	Source
gf	200.27	kJ/mol	Joback Method
hf	-104.18	kJ/mol	Joback Method
hfus	16.02	kJ/mol	Joback Method
hvap	50.17	kJ/mol	Joback Method
log10ws	-5.11		Crippen Method
logp	4.869		Crippen Method
mcvol	191.890	ml/mol	McGowan Method
pc	2069.88	kPa	Joback Method
rinpol	1495.00		NIST Webbook
rinpol	1525.00		NIST Webbook
rinpol	1514.00		NIST Webbook
rinpol	1512.00		NIST Webbook
rinpol	1539.00		NIST Webbook
rinpol	1539.00		NIST Webbook
rinpol	1557.00		NIST Webbook
rinpol	1504.00		NIST Webbook
rinpol	1504.00		NIST Webbook
rinpol	1525.00		NIST Webbook
rinpol	1514.00		NIST Webbook
rinpol	1532.00		NIST Webbook
rinpol	1532.00		NIST Webbook
rinpol	1532.00		NIST Webbook
rinpol	1512.00		NIST Webbook
ripol	1747.00		NIST Webbook
ripol	1763.50		NIST Webbook
ripol	1747.00		NIST Webbook
ripol	1763.50		NIST Webbook
tb	584.06	K	Joback Method
tc	811.24	K	Joback Method

tf	314.19	K	Joback Method
vc	0.726	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	501.06	J/mol×K	584.06	Joback Method
cpg	523.47	J/mol×K	621.92	Joback Method
cpg	544.48	J/mol×K	659.79	Joback Method
cpg	564.25	J/mol×K	697.65	Joback Method
cpg	582.97	J/mol×K	735.51	Joback Method
cpg	600.80	J/mol×K	773.38	Joback Method
cpg	617.92	J/mol×K	811.24	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R207891&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mccol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
ripol:	Polar retention indices

tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/62-264-0/alpha-Alaskene.pdf>

Generated by Cheméo on 2025-12-23 06:44:48.599637941 +0000 UTC m=+6220486.129678595.

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