

12-nor-Preziza-7(15)-en-2-one

Inchi:	InChI=1S/C14H20O/c1-9-10-6-7-14(8-10)11(13(9,2)3)4-5-12(14)15/h10-11H,1,4-8H2,2-3
InchiKey:	ROIPFIPZYGGKEB-PEMBBPQUSA-N
Formula:	C14H20O
SMILES:	C=C1C2CCC3(C2)C(=O)CCC3C1(C)C
Mol. weight [g/mol]:	204.31

Physical Properties

Property code	Value	Unit	Source
gf	136.85	kJ/mol	Joback Method
hf	-169.53	kJ/mol	Joback Method
hfus	9.05	kJ/mol	Joback Method
hvap	48.64	kJ/mol	Joback Method
log10ws	-3.53		Crippen Method
logp	3.348		Crippen Method
mcvol	172.810	ml/mol	McGowan Method
pc	2472.73	kPa	Joback Method
rinpol	1593.00		NIST Webbook
rinpol	1563.00		NIST Webbook
rinpol	1566.00		NIST Webbook
rinpol	1593.00		NIST Webbook
tb	611.27	K	Joback Method
tc	853.73	K	Joback Method
tf	419.78	K	Joback Method
vc	0.660	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	487.80	J/molxK	611.27	Joback Method
cpg	508.81	J/molxK	651.68	Joback Method
cpg	528.58	J/molxK	692.09	Joback Method
cpg	547.44	J/molxK	732.50	Joback Method
cpg	565.72	J/molxK	772.91	Joback Method
cpg	583.74	J/molxK	813.32	Joback Method

Sources

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=R198584&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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