

Indipone

Inchi:	InChI=1S/C15H24O/c1-10(2)14(16)15-8-7-12(4)13(15)6-5-11(3)9-15/h9-10,12-13H,5-8H
InchiKey:	NUTDFMLLZCEHML-OPFPJEHXSА-N
Formula:	C15H24O
SMILES:	CC1=CC2(C(=O)C(C)C)CCC(C)C2CC1
Mol. weight [g/mol]:	220.35

Physical Properties

Property code	Value	Unit	Source
gf	36.39	kJ/mol	Joback Method
hf	-302.46	kJ/mol	Joback Method
hfus	18.26	kJ/mol	Joback Method
hvap	55.18	kJ/mol	Joback Method
log10ws	-4.06		Crippen Method
logp	3.984		Crippen Method
mcvol	197.760	ml/mol	McGowan Method
pc	2030.89	kPa	Joback Method
rinqol	1496.00		NIST Webbook
tb	622.03	K	Joback Method
tc	842.78	K	Joback Method
tf	352.00	K	Joback Method
vc	0.749	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	548.52	J/molxK	622.03	Joback Method
cpg	569.66	J/molxK	658.82	Joback Method
cpg	589.58	J/molxK	695.61	Joback Method
cpg	608.41	J/molxK	732.40	Joback Method
cpg	626.31	J/molxK	769.20	Joback Method
cpg	643.45	J/molxK	805.99	Joback Method
cpg	659.96	J/molxK	842.78	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=R641439&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
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
hvac:	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
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/79-151-7/Indipone.pdf>

Generated by Cheméo on 2024-04-29 09:36:18.348784439 +0000 UTC m=+16672627.269361750.

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