

Ibuprofen, ethyl ester

Inchi:	InChI=1S/C15H22O2/c1-5-17-15(16)12(4)14-8-6-13(7-9-14)10-11(2)3/h6-9,11-12H,5,10H
InchiKey:	HXTFUVWJFLDLJP-UHFFFAOYSA-N
Formula:	C15H22O2
SMILES:	CCOC(=O)C(C)c1ccc(CC(C)C)cc1
Mol. weight [g/mol]:	234.33

Physical Properties

Property code	Value	Unit	Source
gf	-60.60	kJ/mol	Joback Method
hf	-383.23	kJ/mol	Joback Method
hfus	24.00	kJ/mol	Joback Method
hvap	60.30	kJ/mol	Joback Method
log10ws	-3.75		Crippen Method
logp	3.552		Crippen Method
mcvol	205.890	ml/mol	McGowan Method
pc	1923.67	kPa	Joback Method
rinpol	1347.00		NIST Webbook
rinpol	1347.00		NIST Webbook
tb	649.67	K	Joback Method
tc	855.34	K	Joback Method
tf	339.91	K	Joback Method
vc	0.779	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	547.01	J/molxK	649.67	Joback Method
cpg	564.30	J/molxK	683.95	Joback Method
cpg	580.59	J/molxK	718.23	Joback Method
cpg	595.92	J/molxK	752.51	Joback Method
cpg	610.31	J/molxK	786.79	Joback Method
cpg	623.79	J/molxK	821.07	Joback Method
cpg	636.37	J/molxK	855.34	Joback Method
dvisc	0.0023739	Paxs	339.91	Joback Method

dvisc	0.0010314	Paxs	391.54	Joback Method
dvisc	0.0005442	Paxs	443.16	Joback Method
dvisc	0.0003281	Paxs	494.79	Joback Method
dvisc	0.0002177	Paxs	546.42	Joback Method
dvisc	0.0001550	Paxs	598.04	Joback Method
dvisc	0.0001165	Paxs	649.67	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U389552&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
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:	Log ₁₀ of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
m_{cvol}:	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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