

4-Methoxy-2-(1-propenyl)-phenyl isobutyrate

Inchi:	InChI=1S/C14H18O3/c1-5-6-11-9-12(16-4)7-8-13(11)17-14(15)10(2)3/h5-10H,1-4H3/b6-
InchiKey:	HVYCNLFVRHNNHQR-WAYWQWQTSA-N
Formula:	C14H18O3
SMILES:	CC=Cc1cc(OC)ccc1OC(=O)C(C)C
Mol. weight [g/mol]:	234.29

Physical Properties

Property code	Value	Unit	Source
gf	-100.99	kJ/mol	Joback Method
hf	-383.78	kJ/mol	Joback Method
hfus	25.93	kJ/mol	Joback Method
hvap	61.49	kJ/mol	Joback Method
log10ws	-3.74		Crippen Method
logp	3.290		Crippen Method
mvol	193.370	ml/mol	McGowan Method
pc	2119.73	kPa	Joback Method
ripol	2462.00		NIST Webbook
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tb	658.79	K	Joback Method
tc	870.40	K	Joback Method
tf	373.31	K	Joback Method
vc	0.728	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	499.19	J/molxK	658.79	Joback Method
cpg	514.63	J/molxK	694.06	Joback Method
cpg	529.19	J/molxK	729.33	Joback Method
cpg	542.87	J/molxK	764.60	Joback Method
cpg	555.70	J/molxK	799.87	Joback Method
cpg	567.68	J/molxK	835.13	Joback Method
cpg	578.83	J/molxK	870.40	Joback Method
dvisc	0.0010266	Paxs	373.31	Joback Method

dvisc	0.0005495	Paxs	420.89	Joback Method
dvisc	0.0003339	Paxs	468.47	Joback Method
dvisc	0.0002225	Paxs	516.05	Joback Method
dvisc	0.0001587	Paxs	563.63	Joback Method
dvisc	0.0001193	Paxs	611.21	Joback Method
dvisc	0.0000935	Paxs	658.79	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=R418294&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
pc:	Critical Pressure
rip_{ol}:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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