

Benzoic acid, 4-(2-methylbutyl)oxy-, methyl ester

Inchi:	InChI=1S/C13H18O3/c1-4-10(2)9-16-12-7-5-11(6-8-12)13(14)15-3/h5-8,10H,4,9H2,1-3H
InchiKey:	BAMCFGQBADMYOF-UHFFFAOYSA-N
Formula:	C13H18O3
SMILES:	CCC(C)COc1ccc(C(=O)OC)cc1
Mol. weight [g/mol]:	222.28

Physical Properties

Property code	Value	Unit	Source
gf	-180.00	kJ/mol	Joback Method
hf	-468.89	kJ/mol	Joback Method
hfus	23.53	kJ/mol	Joback Method
hvap	58.65	kJ/mol	Joback Method
log10ws	-3.27		Crippen Method
logp	2.898		Crippen Method
mcvol	183.580	ml/mol	McGowan Method
pc	2239.76	kPa	Joback Method
rinpol	1726.00		NIST Webbook
rinpol	1726.00		NIST Webbook
tb	626.77	K	Joback Method
tc	832.59	K	Joback Method
tf	354.60	K	Joback Method
vc	0.692	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	469.86	J/mol×K	626.77	Joback Method
cpg	485.46	J/mol×K	661.07	Joback Method
cpg	500.22	J/mol×K	695.38	Joback Method
cpg	514.14	J/mol×K	729.68	Joback Method
cpg	527.21	J/mol×K	763.98	Joback Method
cpg	539.46	J/mol×K	798.29	Joback Method
cpg	550.88	J/mol×K	832.59	Joback Method
dvisc	0.0014954	Paxs	354.60	Joback Method

dvisc	0.0007775	Paxs	399.96	Joback Method
dvisc	0.0004619	Paxs	445.32	Joback Method
dvisc	0.0003021	Paxs	490.69	Joback Method
dvisc	0.0002123	Paxs	536.05	Joback Method
dvisc	0.0001577	Paxs	581.41	Joback Method
dvisc	0.0001222	Paxs	626.77	Joback Method

Sources

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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U375328&Units=SI

Legend

cpg:	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:	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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