

6,7-dimethoxy-m-cymene

Inchi:	InChI=1S/C12H18O2/c1-9(2)10-5-6-12(14-4)11(7-10)8-13-3/h5-7,9H,8H2,1-4H3
InchiKey:	WMZSSQXSRIEJGC-UHFFFAOYSA-N
Formula:	C12H18O2
SMILES:	COCc1cc(C(C)C)ccc1OC
Mol. weight [g/mol]:	194.27

Physical Properties

Property code	Value	Unit	Source
gf	-69.13	kJ/mol	Joback Method
hf	-347.14	kJ/mol	Joback Method
hfus	18.95	kJ/mol	Joback Method
hvap	50.34	kJ/mol	Joback Method
log10ws	-3.16		Crippen Method
logp	2.965		Crippen Method
mcvol	167.920	ml/mol	McGowan Method
pc	2278.41	kPa	Joback Method
rinpol	1467.00		NIST Webbook
rinpol	1467.00		NIST Webbook
tb	555.00	K	Joback Method
tc	757.29	K	Joback Method
tf	305.92	K	Joback Method
vc	0.629	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	398.99	J/mol×K	555.00	Joback Method
cpg	414.88	J/mol×K	588.71	Joback Method
cpg	430.06	J/mol×K	622.43	Joback Method
cpg	444.52	J/mol×K	656.14	Joback Method
cpg	458.28	J/mol×K	689.86	Joback Method
cpg	471.33	J/mol×K	723.57	Joback Method
cpg	483.66	J/mol×K	757.29	Joback Method
dvisc	0.0014180	Paxs	305.92	Joback Method

dvisc	0.0007434	Paxs	347.43	Joback Method
dvisc	0.0004474	Paxs	388.95	Joback Method
dvisc	0.0002969	Paxs	430.46	Joback Method
dvisc	0.0002118	Paxs	471.97	Joback Method
dvisc	0.0001596	Paxs	513.49	Joback Method
dvisc	0.0001254	Paxs	555.00	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=R340993&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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

Latest version available from:

<https://www.chemeo.com/cid/64-203-5/6-7-dimethoxy-m-cymene.pdf>

Generated by Cheméo on 2024-04-27 07:02:51.711862273 +0000 UTC m=+16490620.632439588.

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