

Furan, 2,3-dihydro-5-(2-methylpropyl)-

Inchi:	InChI=1S/C8H14O/c1-7(2)6-8-4-3-5-9-8/h4,7H,3,5-6H2,1-2H3
InchiKey:	FBQBYCIPUSEWMX-UHFFFAOYSA-N
Formula:	C8H14O
SMILES:	CC(C)CC1=CCCO1
Mol. weight [g/mol]:	126.20
CAS:	56755-31-8

Physical Properties

Property code	Value	Unit	Source
gf	-7.49	kJ/mol	Joback Method
hf	-218.60	kJ/mol	Joback Method
hfus	14.63	kJ/mol	Joback Method
hvap	39.04	kJ/mol	Joback Method
log10ws	-2.26		Crippen Method
logp	2.337		Crippen Method
mcvol	114.290	ml/mol	McGowan Method
pc	3254.14	kPa	Joback Method
tb	433.04	K	Joback Method
tc	634.64	K	Joback Method
tf	219.91	K	Joback Method
vc	0.426	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	232.86	J/molxK	433.04	Joback Method
cpg	297.89	J/molxK	601.04	Joback Method
cpg	286.28	J/molxK	567.44	Joback Method
cpg	273.99	J/molxK	533.84	Joback Method
cpg	261.02	J/molxK	500.24	Joback Method
cpg	247.31	J/molxK	466.64	Joback Method
cpg	308.86	J/molxK	634.64	Joback Method
dvisc	0.0003163	Paxs	433.04	Joback Method
dvisc	0.0004171	Paxs	397.52	Joback Method

dvisc	0.0005806	Paxs	362.00	Joback Method
dvisc	0.0008685	Paxs	326.47	Joback Method
dvisc	0.0014335	Paxs	290.95	Joback Method
dvisc	0.0027198	Paxs	255.43	Joback Method
dvisc	0.0063466	Paxs	219.91	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=C56755318&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
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/34-045-4/Furan-2-3-dihydro-5-2-methylpropyl.pdf>

Generated by Cheméo on 2024-04-25 18:55:00.888475451 +0000 UTC m=+16360549.809052779.

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