

2-Methyl-6-hepten-5-one

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

Physical Properties

Property code	Value	Unit	Source
gf	-27.04	kJ/mol	Joback Method
hf	-200.88	kJ/mol	Joback Method
hfus	13.27	kJ/mol	Joback Method
hvap	39.09	kJ/mol	Joback Method
log10ws	-2.06		Crippen Method
logp	2.178		Crippen Method
mcvol	120.850	ml/mol	McGowan Method
pc	2850.52	kPa	Joback Method
rinpola	896.00		NIST Webbook
tb	432.55	K	Joback Method
tc	615.62	K	Joback Method
tf	213.09	K	Joback Method
vc	0.465	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	239.57	J/molxK	432.55	Joback Method
cpg	295.62	J/molxK	585.10	Joback Method
cpg	285.43	J/molxK	554.59	Joback Method
cpg	274.75	J/molxK	524.08	Joback Method
cpg	263.55	J/molxK	493.57	Joback Method
cpg	251.83	J/molxK	463.06	Joback Method
cpg	305.33	J/molxK	615.62	Joback Method
dvisc	0.0002805	Paxs	432.55	Joback Method
dvisc	0.0003689	Paxs	395.97	Joback Method

dvisc	0.0005129	Paxs	359.40	Joback Method
dvisc	0.0007685	Paxs	322.82	Joback Method
dvisc	0.0012767	Paxs	286.24	Joback Method
dvisc	0.0024614	Paxs	249.67	Joback Method
dvisc	0.0059449	Paxs	213.09	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R326006&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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/66-995-5/2-Methyl-6-hepten-5-one.pdf>

Generated by Cheméo on 2024-04-17 01:37:56.189650876 +0000 UTC m=+15607125.110228189.

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