

2-Cyclohexen-1-one, 3,4,4-trimethyl-

Inchi:	InChI=1S/C9H14O/c1-7-6-8(10)4-5-9(7,2)3/h6H,4-5H2,1-3H3
InchiKey:	LJILDAULEWAKOE-UHFFFAOYSA-N
Formula:	C9H14O
SMILES:	CC1=CC(=O)CCC1(C)C
Mol. weight [g/mol]:	138.21
CAS:	17299-41-1

Physical Properties

Property code	Value	Unit	Source
gf	-58.40	kJ/mol	Joback Method
hf	-250.92	kJ/mol	Joback Method
hfus	4.95	kJ/mol	Joback Method
hvap	40.11	kJ/mol	Joback Method
log10ws	-2.38		Crippen Method
logp	2.322		Crippen Method
mcvol	124.080	ml/mol	McGowan Method
pc	3170.40	kPa	Joback Method
rinpol	1198.00		NIST Webbook
ripol	1720.00		NIST Webbook
tb	490.20	K	NIST Webbook
tc	727.23	K	Joback Method
tf	303.97	K	Joback Method
vc	0.464	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	275.69	J/molxK	497.07	Joback Method
cpg	291.98	J/molxK	535.43	Joback Method
cpg	307.32	J/molxK	573.79	Joback Method
cpg	321.80	J/molxK	612.15	Joback Method
cpg	335.52	J/molxK	650.51	Joback Method
cpg	348.57	J/molxK	688.87	Joback Method
cpg	361.03	J/molxK	727.23	Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	371.20	K	1.70	NIST Webbook

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C17299411&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
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvp:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tbrp:	Boiling point at reduced pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/13-895-4/2-Cyclohexen-1-one-3-4-4-trimethyl.pdf>

Generated by Cheméo on 2024-04-24 18:58:31.672115039 +0000 UTC m=+16274360.592692353.
Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.