

2-Pentene, 3,4,4-trimethyl-

Other names:	3,4,4-Trimethyl-2-pentene 3,4,4-Trimethylpentene-2 3,4,4-trimethylpent-2-ene NSC 73944
Inchi:	InChI=1S/C8H16/c1-6-7(2)8(3,4)5/h6H,1-5H3
InchiKey:	FZQMZRXXKWHQJAG-UHFFFAOYSA-N
Formula:	C8H16
SMILES:	CC=C(C)C(C)(C)C
Mol. weight [g/mol]:	112.21
CAS:	598-96-9

Physical Properties

Property code	Value	Unit	Source
gf	90.99	kJ/mol	Joback Method
hf	-109.77	kJ/mol	Joback Method
hfus	7.95	kJ/mol	Joback Method
hvap	32.14	kJ/mol	Joback Method
log10ws	-2.78		Crippen Method
logp	2.999		Crippen Method
mcvol	119.280	ml/mol	McGowan Method
pc	2741.15	kPa	Joback Method
rinpol	751.00		NIST Webbook
rinpol	752.00		NIST Webbook
rinpol	745.50		NIST Webbook
rinpol	746.50		NIST Webbook
rinpol	752.00		NIST Webbook
rinpol	746.50		NIST Webbook
rinpol	752.00		NIST Webbook
rinpol	752.00		NIST Webbook
tb	385.10 ± 1.00	K	NIST Webbook
tb	385.15 ± 0.20	K	NIST Webbook
tb	385.24 ± 0.20	K	NIST Webbook
tb	385.10 ± 1.00	K	NIST Webbook
tb	385.00 ± 4.00	K	NIST Webbook
tb	381.40 ± 7.00	K	NIST Webbook
tb	385.40 ± 0.40	K	NIST Webbook
tb	385.20 ± 0.40	K	NIST Webbook

tb	385.60 ± 1.00	K	NIST Webbook
tb	383.00 ± 4.00	K	NIST Webbook
tb	384.50 ± 1.00	K	NIST Webbook
tb	385.00 ± 7.00	K	NIST Webbook
tb	385.10 ± 1.00	K	NIST Webbook
tc	570.42	K	Joback Method
tf	163.30	K	Joback Method
vc	0.454	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	219.11	J/mol×K	383.25	Joback Method
cpg	233.89	J/mol×K	414.44	Joback Method
cpg	247.86	J/mol×K	445.64	Joback Method
cpg	261.06	J/mol×K	476.83	Joback Method
cpg	273.52	J/mol×K	508.03	Joback Method
cpg	285.30	J/mol×K	539.22	Joback Method
cpg	296.41	J/mol×K	570.42	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.41614e+01
Coeff. B	-3.29361e+03
Coeff. C	-4.82220e+01
Temperature range (K), min.	285.62
Temperature range (K), max.	420.39

Sources

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C598969&Units=SI>

The Yaws Handbook of Vapor Pressure:

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.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
hvac:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mccol:	McGowan's characteristic volume
pc:	Critical Pressure
pvac:	Vapor 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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