

Pentane, 1-iodo-3-methyl

Inchi:	InChI=1S/C6H13I/c1-3-6(2)4-5-7/h6H,3-5H2,1-2H3
InchiKey:	VYP SHMOLHNQXBM-UHFFFAOYSA-N
Formula:	C6H13I
SMILES:	CCC(C)CCl
Mol. weight [g/mol]:	212.07

Physical Properties

Property code	Value	Unit	Source
gf	55.32	kJ/mol	Joback Method
hf	-95.58	kJ/mol	Joback Method
hfus	12.18	kJ/mol	Joback Method
hvap	37.94	kJ/mol	Joback Method
log10ws	-3.04		Crippen Method
logp	2.858		Crippen Method
mcvol	121.220	ml/mol	McGowan Method
pc	3059.17	kPa	Joback Method
rinpol	1058.00		NIST Webbook
rinpol	995.00		NIST Webbook
rinpol	1058.00		NIST Webbook
rinpol	995.00		NIST Webbook
ripol	1214.00		NIST Webbook
ripol	1214.00		NIST Webbook
tb	429.38	K	Joback Method
tc	634.59	K	Joback Method
tf	200.44	K	Joback Method
vc	0.454	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	209.62	J/molxK	429.38	Joback Method
cpg	260.91	J/molxK	600.38	Joback Method
cpg	251.72	J/molxK	566.18	Joback Method
cpg	242.02	J/molxK	531.98	Joback Method

cpg	231.78	J/mol×K	497.78	Joback Method
cpg	220.99	J/mol×K	463.58	Joback Method
cpg	269.61	J/mol×K	634.59	Joback Method
dvisc	0.0003483	Paxs	429.38	Joback Method
dvisc	0.0004646	Paxs	391.22	Joback Method
dvisc	0.0006596	Paxs	353.07	Joback Method
dvisc	0.0010195	Paxs	314.91	Joback Method
dvisc	0.0017766	Paxs	276.75	Joback Method
dvisc	0.0036980	Paxs	238.60	Joback Method
dvisc	0.0101753	Paxs	200.44	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R25530&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

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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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