Pentane, 3-(bromomethyl)-

Other names:	1-Bromo-2-ethylbutane
	2-Ethylbutyl bromide
	3-(Bromomethyl)pentane
	Butane, 1-bromo-2-ethyl-
Inchi:	InChI=1S/C6H13Br/c1-3-6(4-2)5-7/h6H,3-5H2,1-2H3
InchiKey:	KKGUMGWNFARLSL-UHFFFAOYSA-N
Formula:	C6H13Br
SMILES:	CCC(CC)CBr
Mol. weight [g/mol]:	165.07
CAS:	3814-34-4

Physical Properties

Property code	Value	Unit	Source
gf	11.52	kJ/mol	Joback Method
hf	-146.12	kJ/mol	Joback Method
hfus	13.06	kJ/mol	Joback Method
hvap	35.00	kJ/mol	Joback Method
log10ws	-2.52		Crippen Method
logp	2.817		Crippen Method
mcvol	112.900	ml/mol	McGowan Method
pc	3392.03	kPa	Joback Method
tb	416.70	K	NIST Webbook
tc	590.50	K	Joback Method
tf	202.18	K	Joback Method
vc	0.427	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	197.64	J/mol×K	402.40	Joback Method
cpg	247.93	J/mol×K	559.15	Joback Method
cpg	238.81	J/mol×K	527.80	Joback Method
cpg	229.23	J/mol×K	496.45	Joback Method
cpg	219.19	J/mol×K	465.10	Joback Method

cpg	208.66	J/mol×K	433.75	Joback Method
cpg	256.62	J/mol×K	590.50	Joback Method
dvisc	0.0003340	Pa×s	402.40	Joback Method
dvisc	0.0004371	Pa×s	369.03	Joback Method
dvisc	0.0006036	Pa×s	335.66	Joback Method
dvisc	0.0008951	Pa×s	302.29	Joback Method
dvisc	0.0014638	Pa×s	268.92	Joback Method
dvisc	0.0027514	Pa×s	235.55	Joback Method
dvisc	0.0063698	Pa×s	202.18	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	ln(Pvp) = A + B/(T + C)
Coeff. A	1.53359e+01
Coeff. B	-3.85592e+03
Coeff. C	-5.69230e+01
Temperature range (K), min.	313.16
Temperature range (K), max.	441.58

Sources

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/ci990307l
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=C3814344&Units=SI

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
pvap:	Vapor pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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