

2,3-Dimethyl-1,2-dibromopentane

Inchi:	InChI=1S/C7H14Br2/c1-4-6(2)7(3,9)5-8/h6H,4-5H2,1-3H3
InchiKey:	OTWXHGYTYFOXAN-UHFFFAOYSA-N
Formula:	C7H14Br2
SMILES:	CCC(C)C(C)(Br)CBr
Mol. weight [g/mol]:	257.99

Physical Properties

Property code	Value	Unit	Source
gf	37.10	kJ/mol	Joback Method
hf	-149.18	kJ/mol	Joback Method
hfus	13.52	kJ/mol	Joback Method
hvap	42.36	kJ/mol	Joback Method
log10ws	-3.49		Crippen Method
logp	3.581		Crippen Method
mvol	144.490	ml/mol	McGowan Method
pc	3360.64	kPa	Joback Method
rinpol	1054.00		NIST Webbook
rinpol	1054.00		NIST Webbook
tb	488.21	K	Joback Method
tc	705.09	K	Joback Method
tf	275.67	K	Joback Method
vc	0.534	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	278.69	J/mol×K	488.21	Joback Method
cpg	334.80	J/mol×K	668.94	Joback Method
cpg	325.12	J/mol×K	632.80	Joback Method
cpg	314.73	J/mol×K	596.65	Joback Method
cpg	303.57	J/mol×K	560.50	Joback Method
cpg	291.58	J/mol×K	524.36	Joback Method
cpg	343.82	J/mol×K	705.09	Joback Method
dvisc	0.0003149	Paxs	488.21	Joback Method

dvisc	0.0004213	Paxs	452.79	Joback Method
dvisc	0.0005924	Paxs	417.36	Joback Method
dvisc	0.0008871	Paxs	381.94	Joback Method
dvisc	0.0014430	Paxs	346.52	Joback Method
dvisc	0.0026220	Paxs	311.09	Joback Method
dvisc	0.0055548	Paxs	275.67	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R559370&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
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/57-647-1/2-3-Dimethyl-1-2-dibromopentane.pdf>

Generated by Cheméo on 2024-04-23 15:28:44.646087717 +0000 UTC m=+16175373.566665028.

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