

2-Bromobenzyl alcohol, isopropyl ether

Inchi:	InChI=1S/C10H13BrO/c1-8(2)12-7-9-5-3-4-6-10(9)11/h3-6,8H,7H2,1-2H3
InchiKey:	GRVVRTKOGFUJML-UHFFFAOYSA-N
Formula:	C10H13BrO
SMILES:	CC(C)OCc1ccccc1Br
Mol. weight [g/mol]:	229.11

Physical Properties

Property code	Value	Unit	Source
gf	42.98	kJ/mol	Joback Method
hf	-135.84	kJ/mol	Joback Method
hfus	18.26	kJ/mol	Joback Method
hvap	49.25	kJ/mol	Joback Method
log10ws	-3.97		Crippen Method
logp	3.374		Crippen Method
mcvol	151.370	ml/mol	McGowan Method
pc	3079.57	kPa	Joback Method
rinpola	1371.00		NIST Webbook
tb	548.00	K	Joback Method
tc	772.32	K	Joback Method
tf	308.43	K	Joback Method
vc	0.561	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	322.42	J/molxK	548.00	Joback Method
cpg	384.12	J/molxK	734.93	Joback Method
cpg	373.32	J/molxK	697.55	Joback Method
cpg	361.77	J/molxK	660.16	Joback Method
cpg	349.46	J/molxK	622.77	Joback Method
cpg	336.35	J/molxK	585.39	Joback Method
cpg	394.20	J/molxK	772.32	Joback Method
dvisc	0.0001912	Paxs	548.00	Joback Method
dvisc	0.0002433	Paxs	508.07	Joback Method

dvisc	0.0003225	Paxs	468.14	Joback Method
dvisc	0.0004505	Paxs	428.22	Joback Method
dvisc	0.0006742	Paxs	388.29	Joback Method
dvisc	0.0011067	Paxs	348.36	Joback Method
dvisc	0.0020654	Paxs	308.43	Joback Method

Sources

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=U378181&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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/28-893-0/2-Bromobenzyl-alcohol-isopropyl-ether.pdf>

Generated by Cheméo on 2024-04-19 02:02:48.778423716 +0000 UTC m=+15781417.699001047.

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