

Phenol, 4-propoxy-

Other names:	Phenol, p-propoxy- p-Propoxyphenol 4-Propoxyphenol Hydroquinone monopropyl ether 4-n-Propoxyphenol
Inchi:	InChI=1S/C9H12O2/c1-2-7-11-9-5-3-8(10)4-6-9/h3-6,10H,2,7H2,1H3
InchiKey:	KIIIPQXXLVCCQP-UHFFFAOYSA-N
Formula:	C9H12O2
SMILES:	CCCOc1ccc(O)cc1
Mol. weight [g/mol]:	152.19
CAS:	18979-50-5

Physical Properties

Property code	Value	Unit	Source
gf	-122.31	kJ/mol	Joback Method
hf	-302.09	kJ/mol	Joback Method
hfus	20.08	kJ/mol	Joback Method
hvap	53.33	kJ/mol	Joback Method
log10ws	-1.98		Crippen Method
logp	2.181		Crippen Method
mcvol	125.650	ml/mol	McGowan Method
pc	3815.10	kPa	Joback Method
tb	535.04	K	Joback Method
tc	754.68	K	Joback Method
tf	351.56	K	Joback Method
vc	0.415	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	290.27	J/mol×K	535.04	Joback Method
cpg	302.86	J/mol×K	571.65	Joback Method
cpg	314.65	J/mol×K	608.25	Joback Method
cpg	325.69	J/mol×K	644.86	Joback Method

cpg	336.05	J/mol×K	681.47	Joback Method
cpg	345.78	J/mol×K	718.08	Joback Method
cpg	354.94	J/mol×K	754.68	Joback Method
dvisc	0.0022554	Pa×s	351.56	Joback Method
dvisc	0.0009440	Pa×s	382.14	Joback Method
dvisc	0.0004496	Pa×s	412.72	Joback Method
dvisc	0.0002372	Pa×s	443.30	Joback Method
dvisc	0.0001359	Pa×s	473.88	Joback Method
dvisc	0.0000833	Pa×s	504.46	Joback Method
dvisc	0.0000540	Pa×s	535.04	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=C18979505&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
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/64-803-9/Phenol-4-propoxy.pdf>

Generated by Cheméo on 2025-12-05 12:35:57.935313969 +0000 UTC m=+4686355.465354639.

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