

Phenol, 4-[1-ethyl-2-methyl-1-(1-methylpropyl)propyl]

Inchi:	InChI=1S/C16H26O/c1-6-13(5)16(7-2,12(3)4)14-8-10-15(17)11-9-14/h8-13,17H,6-7H2,1-
InchiKey:	DDJFQXYNOJOKOE-UHFFFAOYSA-N
Formula:	C16H26O
SMILES:	CCC(C)C(CC)(c1ccc(O)cc1)C(C)C
Mol. weight [g/mol]:	234.38

Physical Properties

Property code	Value	Unit	Source
gf	39.59	kJ/mol	Joback Method
hf	-333.66	kJ/mol	Joback Method
hfus	22.56	kJ/mol	Joback Method
hvap	64.43	kJ/mol	Joback Method
log10ws	-4.36		Crippen Method
logp	4.742		Crippen Method
mvol	218.410	ml/mol	McGowan Method
pc	1994.77	kPa	Joback Method
rinpol	1798.00		NIST Webbook
rinpol	1798.00		NIST Webbook
tb	668.67	K	Joback Method
tc	886.06	K	Joback Method
tf	380.64	K	Joback Method
vc	0.766	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	614.70	J/molxK	668.67	Joback Method
cpg	633.29	J/molxK	704.90	Joback Method
cpg	650.70	J/molxK	741.13	Joback Method
cpg	667.06	J/molxK	777.36	Joback Method
cpg	682.47	J/molxK	813.60	Joback Method
cpg	697.05	J/molxK	849.83	Joback Method
cpg	710.92	J/molxK	886.06	Joback Method
dvisc	0.0023307	Paxs	380.64	Joback Method

dvisc	0.0005882	Paxs	428.64	Joback Method
dvisc	0.0001959	Paxs	476.65	Joback Method
dvisc	0.0000798	Paxs	524.65	Joback Method
dvisc	0.0000378	Paxs	572.66	Joback Method
dvisc	0.0000201	Paxs	620.67	Joback Method
dvisc	0.0000117	Paxs	668.67	Joback Method

Sources

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=R592748&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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/66-222-2/Phenol-4-1-ethyl-2-methyl-1-1-methylpropyl-propyl.pdf>

Generated by Cheméo on 2024-04-24 14:38:37.104426602 +0000 UTC m=+16258766.025003914.

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