2-(4-Methoxyphenyl)ethanol

Other names:	4-Methoxyphenethyl alcohol
	2-(p-Methoxyphenyl)ethanol
	p-Methoxyphenethyl alcohol
	4-Methoxy-«beta»-phenethyl alcohol
	Benzeneethanol, 4-methoxy-
	p-Methoxyphenylethyl alcohol
	4-Methoxyphenylethanol
Inchi:	InChI=1S/C9H12O2/c1-11-9-4-2-8(3-5-9)6-7-10/h2-5,10H,6-7H2,1H3
InchiKey:	AUWDOZOUJWEPBA-UHFFFAOYSA-N
Formula:	C9H12O2
SMILES:	COc1ccc(CCO)cc1
Mol. weight [g/mol]:	152.19
CAS:	702-23-8

Physical Properties

Property code	Value	Unit	Source
gf	-114.14	kJ/mol	Joback Method
hf	-288.48	kJ/mol	Joback Method
hfus	17.99	kJ/mol	Joback Method
hvap	57.66	kJ/mol	Joback Method
log10ws	-1.66		Crippen Method
logp	1.230		Crippen Method
mcvol	125.650	ml/mol	McGowan Method
pc	3497.14	kPa	Joback Method
rinpol	1374.00		NIST Webbook
rinpol	1374.00		NIST Webbook
ripol	2341.00		NIST Webbook
ripol	2355.00		NIST Webbook
ripol	2341.00		NIST Webbook
tb	608.20	K	NIST Webbook
tc	746.03	K	Joback Method
tf	313.18	K	Joback Method
vc	0.469	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	289.51	J/mol×K	551.58	Joback Method
cpg	300.67	J/mol×K	583.99	Joback Method
cpg	311.28	J/mol×K	616.40	Joback Method
cpg	321.36	J/mol×K	648.80	Joback Method
cpg	330.92	J/mol×K	681.21	Joback Method
cpg	339.96	J/mol×K	713.62	Joback Method
cpg	348.51	J/mol×K	746.03	Joback Method
dvisc	0.0055162	Pa×s	313.18	Joback Method
dvisc	0.0018885	Pa×s	352.91	Joback Method
dvisc	0.0008032	Pa×s	392.65	Joback Method
dvisc	0.0003997	Pa×s	432.38	Joback Method
dvisc	0.0002237	Pa×s	472.11	Joback Method
dvisc	0.0001370	Pa×s	511.85	Joback Method
dvisc	0.0000901	Pa×s	551.58	Joback Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C702238&Units=SI
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

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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume