Benzene, 2-iodo-1,4-dimethyl-

Other names:	1-Iodo-2,5-dimethylbenzene
	2,5-Dimethyliodobenzene
	2-Iodo-1,4-dimethylbenzene
	2-Iodo-p-xylene
	p-Xylene, 2-iodo-
Inchi:	InChI=1S/C8H9I/c1-6-3-4-7(2)8(9)5-6/h3-5H,1-2H3
InchiKey:	WYZVNUSNUCABRF-UHFFFAOYSA-N
Formula:	C8H9I
SMILES:	Cc1ccc(C)c(I)c1
Mol. weight [g/mol]:	232.06
CAS:	1122-42-5

Physical Properties

Property code	Value	Unit	Source
gf	167.75	kJ/mol	Joback Method
hf	82.01	kJ/mol	Joback Method
hfus	14.14	kJ/mol	Joback Method
hvap	46.38	kJ/mol	Joback Method
log10ws	-3.57		Crippen Method
logp	2.908		Crippen Method
mcvol	125.640	ml/mol	McGowan Method
pc	3431.89	kPa	Joback Method
tb	512.22	K	Joback Method
tc	761.77	K	Joback Method
tf	289.44	K	Joback Method
vc	0.464	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	222.94	J/mol×K	512.22	Joback Method
cpg	234.35	J/mol×K	553.81	Joback Method
cpg	244.98	J/mol×K	595.40	Joback Method
cpg	254.87	J/mol×K	637.00	Joback Method

cpg	264.08	J/mol×K	678.59	Joback Method
cpg	272.63	J/mol×K	720.18	Joback Method
cpg	280.59	J/mol×K	761.77	Joback Method
dvisc	0.0021926	Pa×s	289.44	Joback Method
dvisc	0.0012902	Pa×s	326.57	Joback Method
dvisc	0.0008460	Pa×s	363.70	Joback Method
dvisc	0.0005998	Pa×s	400.83	Joback Method
dvisc	0.0004508	Pa×s	437.96	Joback Method
dvisc	0.0003543	Pa×s	475.09	Joback Method
dvisc	0.0002884	Pa×s	512.22	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	ln(Pvp) = A + B/(T + C)
Coeff. A	1.36324e+01
Coeff. B	-3.82899e+03
Coeff. C	-8.01920e+01
Temperature range (K), min.	367.12
Temperature range (K), max.	540.36

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=C1122425&Units=SI

The Yaws Handbook of Vapor Pressure: https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure

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
pvap:	Vapor pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume