

Benzene,2-cyclopenten-1-yl-

Other names:	3-Phenyl-1-cyclopentene
Inchi:	InChI=1S/C11H12/c1-2-6-10(7-3-1)11-8-4-5-9-11/h1-4,6-8,11H,5,9H2
InchiKey:	ZOZIUVMYACMJGS-UHFFFAOYSA-N
Formula:	C11H12
SMILES:	<chem>C1=CC(c2ccccc2)CC1</chem>
Mol. weight [g/mol]:	144.21
CAS:	37689-22-8

Physical Properties

Property code	Value	Unit	Source
gf	220.66	kJ/mol	Joback Method
hf	84.42	kJ/mol	Joback Method
hfus	13.44	kJ/mol	Joback Method
hvap	42.91	kJ/mol	Joback Method
ie	9.20 ± 0.05	eV	NIST Webbook
log10ws	-3.24		Crippen Method
logp	3.120		Crippen Method
mcvol	126.930	ml/mol	McGowan Method
pc	3352.86	kPa	Joback Method
tb	492.20	K	Joback Method
tc	730.84	K	Joback Method
tf	251.81	K	Joback Method
vc	0.470	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	268.90	J/molxK	492.20	Joback Method
cpg	347.12	J/molxK	691.07	Joback Method
cpg	333.84	J/molxK	651.30	Joback Method
cpg	319.46	J/molxK	611.52	Joback Method
cpg	303.88	J/molxK	571.75	Joback Method
cpg	287.06	J/molxK	531.97	Joback Method
cpg	359.36	J/molxK	730.84	Joback Method

dvisc	0.0002887	Paxs	492.20	Joback Method
dvisc	0.0003579	Paxs	452.13	Joback Method
dvisc	0.0004628	Paxs	412.07	Joback Method
dvisc	0.0006324	Paxs	372.00	Joback Method
dvisc	0.0009318	Paxs	331.94	Joback Method
dvisc	0.0015273	Paxs	291.88	Joback Method
dvisc	0.0029293	Paxs	251.81	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C37689228&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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
ie:	Ionization energy
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/50-340-8/Benzene-2-cyclopenten-1-yl.pdf>

Generated by Cheméo on 2024-04-24 21:44:48.006846772 +0000 UTC m=+16284336.927424087.

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