

Phenindione

Other names:	1H-Indene-1,3(2H)-dione, 2-phenyl-
	Athrombon
	Bindan
	Danedion
	Danilon
	Danilone
	Diadilan
	Dindevan
	Dineval
	Diophindane
	Emandion
	Emandione
	Fenhydren
	Fenilin
	Fenindion
	Hedulín
	1,3-Indandione, 2-phenyl-
	Indema
	Indion
	Indon
	Phenhydren
	Phenillin
	2-Phenyl-1,3-diketohydrindene
	Phenylén
	Phenylin
	2-Phenylindandione
	2-Phenyl-1,3-indandione
	Phenylindanedione
	2-Phenyl-1,3-indanedione
	Phenylindione
	PID
	Pindione
	Rectadione
	Theradione
	Tromazal
	Trombol
	Eridione
	Phenylene
	Phenyllin
	2-Phenylindan-1,3-dione

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

CAS:

2-Phenyl-1H-indene-1,3(2H)-dione

Cronodione

Hemolidione

Thrombasal

1,3(2H)-Indenedione, 2-phenyl-

InChI=1S/C15H10O2/c16-14-11-8-4-5-9-12(11)15(17)13(14)10-6-2-1-3-7-10/h1-9,13H

NFBAXHOPROOJAW-UHFFFAOYSA-N

C15H10O2

O=C1c2ccccc2C(=O)C1c1ccccc1

222.24

83-12-5

Physical Properties

Property code	Value	Unit	Source
gf	106.18	kJ/mol	Joback Method
hf	-93.94	kJ/mol	Joback Method
hfus	19.45	kJ/mol	Joback Method
hvap	62.61	kJ/mol	Joback Method
log10ws	-3.85		Crippen Method
logp	2.849		Crippen Method
mcvol	166.970	ml/mol	McGowan Method
pc	3049.04	kPa	Joback Method
rinpol	2055.00		NIST Webbook
tb	743.32	K	Joback Method
tc	1020.04	K	Joback Method
tf	478.55	K	Joback Method
vc	0.630	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	461.07	J/mol×K	743.32	Joback Method
cpg	477.06	J/mol×K	789.44	Joback Method
cpg	491.51	J/mol×K	835.56	Joback Method
cpg	504.46	J/mol×K	881.68	Joback Method
cpg	515.95	J/mol×K	927.80	Joback Method
cpg	526.02	J/mol×K	973.92	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=C83125&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
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/84-995-5/Phenindione.pdf>

Generated by Cheméo on 2025-12-06 01:48:39.882920428 +0000 UTC m=+4733917.412961093.

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