

Pulegone, 2,4-dinitrophenyl hydrazone

Inchi:	InChI=1S/C16H20N4O4/c1-10(2)12-5-4-11(3)15(8-12)18-17-14-7-6-13(19(21)22)9-16(14)
InchiKey:	WLFDIFNUOQDTAZ-OBGWFSINSA-N
Formula:	C16H20N4O4
SMILES:	CC(C)=C1CCC(C)C(=NNc2ccc([N+](=O)[O-])cc2[N+](=O)[O-])C1
Mol. weight [g/mol]:	332.35

Physical Properties

Property code	Value	Unit	Source
hf	33.56	kJ/mol	Joback Method
hvap	99.87	kJ/mol	Joback Method
log10ws	-6.21		Crippen Method
logp	4.427		Crippen Method
mcvol	247.880	ml/mol	McGowan Method
pc	1882.17	kPa	Joback Method
tb	1061.20	K	Joback Method
tc	1334.91	K	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=B6002097&Units=SI

Legend

hf:	Enthalpy of formation 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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

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