

Juglone

Other names:	1,4-Dihydro-1,4-dioxo-5-hydroxynaphthalene 1,4-Naphthalenedione, 5-hydroxy- 1,4-Naphthoquinone, 5-hydroxy- 1,4-Naphthoquinone, 8-hydroxy- 5-Hydroxy-1,4-naftochinon 5-Hydroxy-1,4-naphthalenedione 5-Hydroxy-p-naphthoquinone 5-Hydroxynaphthoquinone 5-hydroxy-1,4-naphthoquinone 5-hydroxynaphthalene-1,4-dione 8-Hydroxy-1,4-naphthalenedione 8-Hydroxy-1,4-naphthoquinone Akhnot C.I. 75500 C.I. Natural Brown 7 Iuglon Juglane Juglon Jugnlon NCI 2323 NSC 153189 NSC 622948 Nucin Regianin Walnut Extract Yuglon
Inchi:	InChI=1S/C10H6O3/c11-7-4-5-9(13)10-6(7)2-1-3-8(10)12/h1-5,12H
InchiKey:	KQPYUDDGWXQXHS-UHFFFAOYSA-N
Formula:	C10H6O3
SMILES:	O=C1C=CC(=O)c2c(O)cccc21
Mol. weight [g/mol]:	174.15
CAS:	481-39-0

Physical Properties

Property code	Value	Unit	Source
gf	-177.38	kJ/mol	Joback Method

hf	-332.62	kJ/mol	Joback Method
hfus	16.30	kJ/mol	Joback Method
hvap	62.99	kJ/mol	Joback Method
ie	8.60	eV	NIST Webbook
ie	8.70 ± 0.02	eV	NIST Webbook
ie	8.85	eV	NIST Webbook
log10ws	-2.09		Crippen Method
logp	1.327		Crippen Method
mcvol	121.850	ml/mol	McGowan Method
pc	4966.33	kPa	Joback Method
tb	690.96	K	Joback Method
tc	965.28	K	Joback Method
tf	508.98	K	Joback Method
vc	0.404	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	314.20	J/mol×K	690.96	Joback Method
cpg	325.85	J/mol×K	736.68	Joback Method
cpg	336.61	J/mol×K	782.40	Joback Method
cpg	346.57	J/mol×K	828.12	Joback Method
cpg	355.79	J/mol×K	873.84	Joback Method
cpg	364.32	J/mol×K	919.56	Joback Method
cpg	372.25	J/mol×K	965.28	Joback Method

Sources

Measurement and correlation of the solubility of juglone in supercritical carbon dioxide:

<https://www.doi.org/10.1016/j.fluid.2011.08.024>

McGowan Method:

https://en.wikipedia.org/wiki/Joback_method

NIST Webbook:

<http://link.springer.com/article/10.1007/BF02311772>

Crippen Method:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C481390&Units=SI>

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

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
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.cheméo.com/cid/34-709-7/Juglone.pdf>

Generated by Cheméo on 2024-04-30 12:14:46.015859173 +0000 UTC m=+16768534.936436500.

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