

Naphtho[2,3-k]fluoranthene

Other names:	Naphto[2,3-k]fluoranthene
Inchi:	InChI=1S/C24H14/c1-2-6-17-12-19-14-23-21-10-4-8-15-7-3-9-20(24(15)21)22(23)13-18(
InchiKey:	UFZAKNUHTLXSMY-UHFFFAOYSA-N
Formula:	C24H14
SMILES:	<chem>c1ccc2cc3cc4c(cc3cc2c1)-c1cccc2cccc-4c12</chem>
Mol. weight [g/mol]:	302.37
CAS:	207-18-1

Physical Properties

Property code	Value	Unit	Source
gf	752.58	kJ/mol	Joback Method
hf	561.85	kJ/mol	Joback Method
hfus	38.47	kJ/mol	Joback Method
hvap	81.51	kJ/mol	Joback Method
log10ws	-10.34		Crippen Method
logp	6.794		Crippen Method
mcvol	232.260	ml/mol	McGowan Method
pc	2241.88	kPa	Joback Method
rinpol	536.96		NIST Webbook
rinpol	543.09		NIST Webbook
rinpol	568.10		NIST Webbook
rinpol	543.06		NIST Webbook
rinpol	543.06		NIST Webbook
rinpol	543.06		NIST Webbook
rinpol	543.09		NIST Webbook
rinpol	543.09		NIST Webbook
tb	882.32	K	Joback Method
tc	1151.38	K	Joback Method
tf	606.52	K	Joback Method
vc	0.911	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
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cpg	661.66	J/molxK	882.32	Joback Method
cpg	739.36	J/molxK	1106.54	Joback Method
cpg	722.35	J/molxK	1061.69	Joback Method
cpg	706.42	J/molxK	1016.85	Joback Method
cpg	691.24	J/molxK	972.01	Joback Method
cpg	676.44	J/molxK	927.16	Joback Method
cpg	757.84	J/molxK	1151.38	Joback Method
dvisc	0.0039222	Paxs	882.32	Joback Method
dvisc	0.0040664	Paxs	836.35	Joback Method
dvisc	0.0042338	Paxs	790.39	Joback Method
dvisc	0.0044300	Paxs	744.42	Joback Method
dvisc	0.0046630	Paxs	698.45	Joback Method
dvisc	0.0049439	Paxs	652.49	Joback Method
dvisc	0.0052883	Paxs	606.52	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=C207181&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
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
mccvol:	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

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