

Property Library for Decane

LibC10H22

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Property Functions

Functional Dependence	Function Name	Call from Fortran Program	Property or Function	Unit of the Result
a = f(p,t,x)	a_ptx_C10H22	A_PTX_C10H22(P,T,X)	Thermal diffusivity	m²/s
$c_p = f(p,t,x)$	cp_ptx_C10H22	CP_PTX_C10H22(P,T,X)	Specific isobaric heat capacity	kJ/(kg K)
$c_V = f(p,t,x)$	cv_ptx_C10H22	CV_PTX_C10H22(P,T,X)	Specific isochoric heat capacity	kJ/(kg K)
$\varepsilon = f(p,t,x)$	eps_ptx_C10H22	EPS_PTX_C10H22(P,T,X)	Dielectric constant	-
$\eta = f(p,t,x)$	eta_ptx_C10H22	ETA_PTX_C10H22(P,T,X)	Dynamic viscosity	Pa . s
h = f(p,t,x)	h_ptx_C10H22	H_PTX_C10H22(P,T,X)	Specific enthalpy	kJ/kg
$\kappa = f(p,t,x)$	ka_ptx_C10H22	KA_PTX_C10H22(P,T,X)	Isentropic exponent	-
$\lambda = f(p,t,x)$	lam_ptx_C10H22	LAM_PTX_C10H22(P,T,X)	Thermal conductivity	W/(m . K)
v = f(p,t,x)	ny_ptx_C10H22	NY_PTX_C10H22(P,T,X)	Kinematic viscosity	m²/s
Pr = f(p,t,x)	pr_ptx_C10H22	PR_PTX_C10H22(P,T,X)	Prandtl-number	-
$p_s = f(t)$	ps_t_C10H22	PS_T_C10H22(T)	Vapor pressure from temperature	bar
ρ = f(p , t , x)	rho_ptx_C10H22	RHO_PTX_C10H22(P,T,X)	Density	kg/m ³
s = f(p,t,x)	s_ptx_C10H22	S_PTX_C10H22(P,T,X)	Specific entropy	kJ/(kg K)
σ = f(t)	sigma_t_C10H22	SIGMA_T_C10H22(T)	Surface tension from temperature	N/m
t = f(p,h)	t_ph_C10H22	T_PH_C10H22(P,H)	Backward function: Temperature from pressure and enthalpy	°C
t = f(p,s)	t_ps_C10H22	T_PS_C10H22(P,S)	Backward function: Temperature from pressure and entropy	°C
$t_{\rm S} = f(p)$	ts_p_C10H22	TS_P_C10H22(P)	Saturation temperature from pressure	°C
u = f(p,t,x)	u_ptx_C10H22	U_PTX_C10H22(P,T,X)	Specific internal energy	kJ/kg
v = f(p,t,x)	v_ptx_C10H22	V_PTX_C10H22(P,T,X)	Specific volume	m³/kg
w = f(p,t,x)	w_ptx_C10H22	W_PTX_C10H22(P,T,X)	Isentropic speed of sound	m/s

Functional Dependence	Function Name	Call from Fortran Program	Property or Function	Unit of the Result
x = f(p,h)	x_ph_C10H22	X_PH_C10H22(P,H)	Backward function: Vapor fraction from pressure and enthalpy	kg/kg
x = f(p,s)	x_ps_C10H22	X_PS_C10H22(P,S)	Backward function: Vapor fraction from pressure and entropy	kg/kg

Units: t in °C

p in bar

x in (kg saturated steam)/(kg wet steam)

Range of validity

for transport properties (a, η , λ , ν , Pr):

from - 29.65 °C to 401.85 °C

Temperature range: from 1.404184 x 10⁻⁵ bar to 2000 bar Pressure range:

for other properties:

Temperature range: from - 29.65 °C to 401.85 °C

from 1.404184 x 10⁻⁵ bar to 8000 bar Pressure range:

Reference state

h = 0 kJ/kg and s = 0 kJ/(kg K) at p = 1,01325 bar on the saturated liquid line (x = 0)

Details on the vapor fraction x

The wet steam region is calculated automatically by the subprograms. For this purpose the following fixed details on the vapor fraction *x* are to be considered:

Single-phase region

If the state point to be calculated is located in the single-phase region (liquid or superheated steam) x = -1 must be entered as a pro-forma value.

Wet-steam region

If the state point to be calculated is located in the wet steam region, a value for x between 0 and 1 (x = 0 for saturated liquid, x = 1 for saturated steam) must be entered. In this case, the backward functions result in the appropriate value between 0 and 1 for x. When calculating wet steam either the given value for t and t = -1000 or the given value for t and t = -1000 and in both cases the value for t between 0 and 1 must be entered.

If p and t and x are entered as given values, the program considers p and t to be appropriate to represent the vapor pressure curve. If this is not the case the calculation for the property of the chosen function results in -1.

Wet steam region: Temperature ranges from $t_{min} = -29.65$ °C to $t_{c} = 344.55$ °C

Pressure ranges from $p_{min} = 1.404184 \text{ x } 10^{-5} \text{ bar to } p_{c} = 21.01369 \text{ bar}$

Note:

If the input values are located outside the range of validity, the calculated function will always result in –1000. Please find more exact details on every function and its corresponding range of validity in the enclosed program documentation in Chapter 3.