

# Property Library for Decamethylcyclopenta-siloxane (D5) C<sub>10</sub>H<sub>30</sub>O<sub>5</sub>Si<sub>5</sub>

## FluidLAB with LibD5 for MATLAB®

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## Property Library for the Calculation of Decamethylcyclopentasiloxane (D5) FluidLAB for MATLAB<sup>®</sup> LibD5

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### 0. Package Contents

### Zip-file "CD\_FluidLAB\_LibD5.zip" including the following files:

FluidLAB_LibD5_Setup.exe	<ul> <li>Installation program for the FluidLAB Add-On for use in MATLAB<sup>®</sup></li> </ul>
LibD5.dll	- DLL with functions of the LibD5 library

**Documentation** 

FluidLAB\_LibD5\_Docu\_Eng.pdf - User's Guide

### **1. Property Functions**

### **1.1 Calculation Programs**

"D5" means Decamethylcyclopentasiloxane ( $C_{10}H_{30}O_5Si_5$ )

Functional Dependence	Function Name	Call from Fortran program	Call in DLL LibD5 as parameter	Property or Function	Unit of the result
$c_p = f(p, t, x)$	cp_ptx_D5	CPPTXD5(P,T,X)	C_CPPTXD5(CP,P,T,X)	Specific isobaric heat capacity	kJ/(kg K)
$c_{v} = f(p, t, x)$	cv_ptx_D5	CVPTXD5(P,T,X)	C_CVPTXD5(CV,P,T,X)	Specific isochoric heat capacity	kJ/(kg K)
$\left(\frac{\partial p}{\partial T}\right)_{v} = \mathbf{f}(\boldsymbol{p}, \boldsymbol{t}, \boldsymbol{x})$	dpdtv_ptx_D5	DPDTVD5(P,T,X)	C_DPDTVD5(DPDT,P,T,X)	Derivative of pressure with respect to temperature (at constant specific volume)	kPa/K
$\left(\frac{\partial p}{\partial v}\right)_T = f(p, t, x)$	dpdvt_ptx_D5	DPDTVD5(P,T,X)	C_DPDVTD5(DPDV,P,T,X)	Derivative of pressure with respect to specific volume (at constant temperature)	kPa/(m <sup>3</sup> /kg)
h = f(p, t, x)	h_ptx_D5	HPTXD5(P,T,X)	C_HPTXD5(H,P,T,X)	Specific enthalpy	kJ/kg
$\kappa = f(p, t, x)$	kappa_ptx_D5	KAPPAPTXD5(P,T,X)	C_KAPPAPTXD5(KAPPA,P,T,X)	Isentropic exponent	-
$p_{\rm s} = f(t)$	ps_t_D5	PSTD5(T)	C_PSTD5(PS,T)	Vapor pressure from temperature	bar
$\rho = f(p, t, x)$	rho_ptx_D5	RHOPTXD5(P,T,X)	C_RHOPTXD5(RHO,P,T,X)	Density	kg/m <sup>3</sup>
s = f(p, t, x)	s_ptx_D5	SPTXD5(P,T,X)	C_SPTXD5(S,P,T,X)	Specific entropy	kJ/(kg K)
t = f(p, h)	t_ph_D5	TPHD5(P,H)	C_TPHD5(T,P,H)	Backward function: Temperature from pressure and enthalpy	°C
t = f(p, s)	t_ps_D5	TPSD5(P,S)	C_TPSD5(T,P,S)	Backward function: Temperature from pressure and entropy	°C
$t_{\rm S} = f(p)$	ts_p_D5	TSPD5(P)	C_TSPD5(TS,P)	Boiling temperature from pressure	°C
u = f(p, t, x)	u_ptx_D5	UPTXD5(P,T,X)	C_UPTXD5(U,P,T,X)	Specific internal energy	kJ/kg
v = f(p, t, x)	v_ptx_D5	VPTXD5(P,T,X)	C_VPTXD5(V,P,T,X)	Specific volume	m³/kg
w = f(p, t, x)	w_ptx_D5	WPTXD5(P,T,X)	C_WPTXD5(W,P,T,X)	Isentropic speed of sound	m/s
x = f(p,h)	x_ph_D5	XPHD5(P,H)	C_XPHD5(X,P,H)	Backward function: Vapor fraction from pressure and enthalpy	kg/kg

Functional Dependence	Function Name	Call from Fortran program	Call in DLL LibD5 as parameter	Property or Function	Unit of the result
x = f(p,s)	x_ps_D5	XPSD5(P,S)	_ 、 、 、 、	Backward function: Vapor fraction from pressure and entropy	kg/kg
Z = f(p, t, x)	Z_ptx_D5	ZPTXD5(P,T,X)	C_ZPTXD5(W,P,T,X)	Compression factor	-

Units:

*t* in °C *p* in bar

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

#### Range of validity

Temperature range:	from $t = 26.85^{\circ}$ C to 399.85 °C
Pressure range:	from $p = 0.00001$ bar to 300 bar

#### **Reference state**

h = 0 kJ/kg and s = 0 kJ/(kg K) at  $t_B = 210.9 \text{ °C}$  on the boiling curve (x = 0;  $p_s = p_N = 1.01325 \text{ bar}$ )

#### 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:

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. Here the backward functions will also result in x = -1.

If the state point to be calculated is located in the two phase region (wet steam), either the value 0 or 1 has to be entered for x (x = 0 for boiling liquid, x = 1 for saturated steam). Here the backward functions will result in x = 0 or x = 1.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1. When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve. If it is not the case the calculation for the quantity of the chosen function to be calculated results in -1000.

Wet steam region: Temperature range from  $t = 26.85 \text{ }^{\circ}\text{C}$  to  $t_c = 346.0846 \text{ }^{\circ}\text{C}$ 

Pressure range from  $p_{s}$  (26.85 °C) = 0.001477917 bar to  $p_{c}$  = 11.6146 bar

#### Hint.

If the calculation results in – 1000, the entered values represent a state point beyond the range of validity of LibD5. For further information on each function and its range of validity see Chapter 3. The same information may also be accessed via the online help pages.

### 2 Application of FluidLAB in MATLAB<sup>®</sup>

The FluidLAB Add-In has been developed to calculate thermodynamic properties in MATLAB<sup>®</sup> more conveniently. Within MATLAB<sup>®</sup> it enables the direct call of functions relating to Decamethylcyclopentasiloxane from the LibD5 property library.

#### 2.1 Installing FluidLAB including LibD5

This section describes the installation of FluidLAB including the LibD5 property library.

Before you begin, it is best to close any Windows<sup>®</sup> applications, since Windows<sup>®</sup> may need to be rebooted during the installation process.

After you have downloaded and extracted the zip-file "CD\_FluidLAB\_LibD5.zip", you will see the folder

CD\_FluidLAB\_LibD5

in your Windows Explorer<sup>®</sup>, Norton Commander<sup>®</sup> or other similar program you are using.

Open this folder by double-clicking on it.

In this folder you will see the following three files:

FluidLAB\_LibD5\_Docu\_Eng.pdf FluidLAB\_LibD5\_Setup.exe LibD5.dll.

In order to run the installation of FluidLAB including the LibD5 property library, double-click on the file

FluidLAB\_LibD5\_Setup.exe.

Installation may start with a window noting that all Windows<sup>®</sup> programs should be closed. When this is the case, the installation can be continued. Click the "Next >" button.

In the following dialog box, "Destination Location", the default path offered automatically for the installation of FluidLAB is

C:\Program Files\FluidLAB\LibD5

By clicking the "B<u>r</u>owse..." button, you can change the installation directory before installation (see figure below).

漫 FluidLAB LibD5 📃	
Destination Location	1
Setup will install FluidLAB LibD5 in the following folder.	
To install into a different folder, click Browse, and select another folder.	
You can choose not to install FluidLAB LibD5 by clicking Cancel to exit Setup.	
Destination Folder	
C:\Program Files\FluidLAB\LibD5	
Wise Installation Wizard®	
< <u>B</u> ack <u>Next&gt;</u> Cancel	

#### Figure 2.1: "Destination Location"

If you wish to change directories, click the "Browse..." button and select your desired directory. The instructions in this documentation refer to the stated default directory. Leave this window by clicking the "Next >" button.

The dialog window "Start Installation" pops up. Click the "Next >" button to continue installation. The FluidLAB files are now being copied into the created directory on your hard drive. Click the "Finish >" button in the following window to complete installation.

The installation program has copied the following files for LibD5 into the directory "C:\Program Files\FluidLAB\LibD5":

advapi32.dll Dformd.dll	LibD5.dll msvcp60.dll
Dforrt.dll	msvcrt.dll
INSTALL.LOG	Unwise.exe
LC.dll	Unwise.ini
- MATLAB <sup>®</sup> -Interface-Program for calculable f	unctions
cp_ptx_D5	t_ph_D5
cv_ptx_D5	t_ps_D5
dpdtv_ptx_D5	ts_p_D5
dpdvt_ptx_D5	u_ptx_D5
h_ptx_D5	v_ptx_D5
Kappa_ptx_D5	w_ptx_D5
ps_t_D5	x_ph_D5
rho_ptx_D5	x_ps_D5
s_ptx_D5	Z_ptx_D5

Now, you have to overwrite the file "LibD5.dll" in your FluidLAB directory with the file of the same name provided on your CD with FluidLAB.

To do this, open the CD in "My Computer" and click on the file "LibD5.dll" in order to highlight it.

Then click on the "Edit" menu in your Explorer and select "Copy".

Now, open your FluidLAB directory (the standard being C:\Program Files\FluidLAB\LibD5) and insert the file "LibD5.dll" by clicking the "Edit" menu in your Explorer and then select "Paste".

Answer the question whether you want to replace the file by clicking the "Yes" button. Now, you have overwritten the file "LibD5.dll" successfully and the property functions are available in MATLAB<sup>®</sup>.

#### Licensing the LibD5 Property Library

The licensing procedure must be carried out when the prompt message appears. In this case, you will see the "License Information" window for LibD5 (see figure below).

License Information	×
LibD5	
Please type in your license key!	?
ОК	Cancel

Figure 2.2: "License Information" window

Here you are asked to type in the license key which you have obtained from the Zittau/Goerlitz University of Applied Sciences. If you do not have this, or have any questions, you will find contact information on the "Content" page of this User's Guide or by clicking the yellow question mark in the "License Information" window. Then the following window will appear:

👬 Help		×
Product:	In order to obtain a license for this product please contact us. LibD5	
Contact:	Zittau/Goerlitz University of Applied Sciences Faculty of Mechanical Engineering Department of Technical Thermodynamics Prof. Hans-Joachim Kretzschmar, Dr. Ines Stoecker Theodor-Koerner-Allee 16 02763 Zittau, Germany	
Phone: Fax: Email: www:	+49-3583-61-1846 +49-3583-61-1846 hj.kretzschmar@hs-zigr.de www.thermodynamics-zittau.de	

Figure 2.3: "Help" window

If you do not enter a valid license it is still possible to use MATLAB<sup>®</sup> by clicking "Cancel". In this case, the LibD5 property library will display the result "–11111111" for every calculation.

The "License Information" window will appear every time you use FluidLAB LibD5 until you enter a license code to complete registration. If you decide not to use FluidLAB LibD5, you can uninstall the program following the instructions given in section 2.4 of this User's Guide.

#### 2.2 Example: Calculation of h = f(p, t, x) in an M-File

Now we will calculate, step by step, the air-specific enthalpy h as a function of pressure p, temperature t and vapor fraction x using FluidLAB.

Please carry out the following instructions:

- Start Windows Explorer<sup>®</sup>, Total Commander<sup>®</sup>, My Computer or another file manager program. The following description refers to Windows Explorer<sup>®</sup>.
- Your Windows Explorer<sup>®</sup> should be set to "Details" for easier viewing. Click the "Views" button and select "Details."
- Switch into the program directory of FluidLAB, in which you will find the folder "\LibD5"; it is generally saved under: "C:\Program Files\FluidLAB"
- Create the folder "\LibD5\_Example" by clicking on "File" in the Explorer<sup>®</sup> menu, then "New" in the menu which appears and afterwards selecting "Folder". Name the new folder "\LibD5\_Example."
- You will now see the following window:

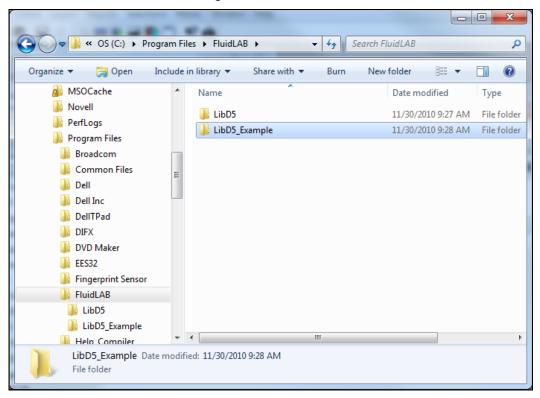


Figure 2.4: Folders "LibD5" and "LibD5\_Example"

- Switch into the directory "\LibD5" within "\FluidLAB", the standard being "C:\Program Files\FluidLAB\LibD5." - You will see the following window:

📀 🗢 🕻 « Program Files )	FluidLAB LibD5	<ul> <li>✓ 4→ Search LibD5</li> </ul>	<u>×</u> 0 -
Organize 🔻 🔳 Open with	Burn New folder		- 🗌 🔞
MSOCache Novell PerfLogs Program Files Common Files Common Files Dell Dell Inc DellTPad DIFX DVD Maker EES32 Fingerprint Sensor FluidLAB LibD5 LibD5_Example Help, Compiler	<ul> <li>advapi32.dll</li> <li>cp_ptx_D5.mexw32</li> <li>cv_ptx_D5.mexw32</li> <li>Dformd.dll</li> <li>Dforrt.dll</li> <li>Dfort.dll</li> <li>Dfotv_ptx_D5.mexw32</li> <li>dpdvt_ptx_D5.mexw32</li> <li>h_ptx_D5.mexw32</li> <li>INSTALL.LOG</li> <li>kappa_ptx_D5.mexw32</li> <li>LC.dll</li> <li>LibD5.dll</li> <li>msvcp60.dll</li> <li>msvcrt.dll</li> <li>ps_t_D5.mexw32</li> <li>rho_ptx_D5.mexw32</li> </ul>	<ul> <li>s_ptx_D5.mexw32</li> <li>t_ph_D5.mexw32</li> <li>t_ps_D5.mexw32</li> <li>ts_p_D5.mexw32</li> <li>u_ptx_D5.mexw32</li> <li>UNWISE.EXE</li> <li>UNWISE.INI</li> <li>v_ptx_D5.mexw32</li> <li>w_ptx_D5.mexw32</li> <li>x_ph_D5.mexw32</li> <li>x_ps_D5.mexw32</li> </ul>	
LibD5.dll Application extension	Date modified: 11/12/2010 10:46 AM Size: 88.0 KB	Date created: 11/30/2010 9:27 AM	I

Figure 2.5: Contents of the folder "LibD5"

You will now have to copy the following files into the directory "C:\Program Files\FluidLAB\LibD5\_Example" in order to calculate the function h = f(p, t, x).

- The following eight files are needed:
  - "advapi32.dll"
  - "Dformd.dll"
  - "Dforrt.dll"
  - "h\_ptx\_D5.mexw32"
  - "LC.dll"
  - "LibD5.dll"
  - "msvcp60.dll"
  - "msvcrt.dll."
- Click the file "h\_ptx\_D5.mexw32", then click "Edit" in the upper menu bar and select "Copy".
- Switch into the directory "C:\Program Files\FluidLAB\LibD5\_Example", click "Edit" and then "Paste".
- Repeat these steps in order to copy the other files listed above. You may also select all the above-named files and then copy them as a group (press the Control button to enable multiple

markings).

- You will see the following window:

🕒 🗢 📔 « Program Files 🕨	FluidLAB  LibD5_Example	Search LibD5_Example
Organize 🔻 Include in library	<ul> <li>Share with          <ul> <li>Burn New fold</li> </ul> </li> </ul>	ler 🔠 🔻 🗍 🔞
<ul> <li>Novell</li> <li>PerfLogs</li> <li>Program Files</li> <li>Broadcom</li> <li>Common Files</li> <li>Dell</li> <li>Dell Inc</li> <li>DellTPad</li> <li>DIFX</li> <li>DVD Maker</li> <li>EES32</li> <li>Fingerprint Sensor</li> <li>FluidLAB</li> <li>LibD5</li> <li>LibD5_Example</li> <li>Help_Compiler</li> <li>ICOL ite</li> </ul>	<ul> <li>Name</li> <li>         advapi32.dll     </li> <li>         Dformd.dll     </li> <li>         Dforrt.dll     </li> <li>         h_ptx_D5.mexw32     </li> <li>         LC.dll     </li> <li>         LibD5.dll     </li> <li>         msvcp60.dll     </li> <li>         msvcrt.dll     </li> </ul>	Date modified       Type         2/9/2009 11:51 AM       Application ex         6/20/2001 3:11 AM       Application ex         6/20/2001 3:10 AM       Application ex         11/23/2010 2:57 PM       MEXW32 File         3/30/2010 2:27 PM       Application ex         11/12/2010 10:46       Application ex         4/14/2008 2:00 PM       Application ex         4/14/2008 2:00 PM       Application ex

Figure 2.6: Contents of the folder "LibD5\_Example"

- Start MATLAB<sup>®</sup> (if you have not started it before).
- Click the button marked in the next figure in order to open the folder "\LibD5\_Example" in the "Current Folder" window.

MATLAB 7.11.0 (R2010b)	
<u>F</u> ile <u>E</u> dit De <u>b</u> ug <u>D</u> esktop <u>W</u> indow <u>H</u> elp	$\langle + \rangle$
🗄 🎦 🚰 🖌 🐂 🛍 🤊 🕅 🏣 🗊 🗐 🖉 Current Folder: C:\Program Files\MATLAB	
Shortcuts 🗷 How to Add 🗷 What's New	
Current Folder IF 🗖 🛪 🗙 Command Window	
→ MATLAB → → → ← /2 → /2 >>	
🗅 Name 🔺	
🛛 🕀 🌛 R2010b	

#### Figure 2.7: Selection of the working directory

- Find and select the directory "C:\Program Files\FluidLAB\LibD5\_Example" in the pop-up menu (see the following image).

Browse For Folder	
Select a new folder	
DVD Maker	
▷ 📕 EES32	
Fingerprint Sensor	
A Di FluidLAB	
LibD5	
📕 LibD5_Example 🚽	
Eolder: LibD5_Example	
Make New Folder OK Cancel	зł

Figure 2.8: Choosing the "LibD5\_Example" folder

- Confirm your selection by clicking the "OK" button.
- First of all you need to create an M–File in MATLAB<sup>®</sup>. Within MATLAB<sup>®</sup> click "Desktop", then select "Editor". Now click on the "New Script" button in the Editor Window.
- If the "Editor" window appears as a separate window, you can embed it into MATLAB<sup>®</sup> by clicking the insertion arrow (see next figure) in order to obtain a better view.



#### Figure 2.9: Embedding the "Editor" window

- In the following figure you will see the "Editor – Untitled" window.

MATLAB 7.11.0 (R2010b)		_			x
<u>File Edit Text Go C</u> ell T <u>o</u> ols De <u>b</u> ug	<u>D</u> esktop <u>W</u> indow <u>H</u> elp				
: 🗋 🖆 👗 🛍 ウ 😋 🎒 🖹 🧕	Current Folder: C:\Program Files\FluidLAB\Lib	D5_Examp	le 🔻 🖻	3	
Shortcuts 🗷 How to Add 🗷 What's New					
Current Folder 🏨 🖬 🕫 E	Editor - Untitled 💛 🖓	N ≤ 1	Workspace	<b>→</b> 1 🗖	× ×
📔 🛶 LibD5_Example 🛛 🔻 🔎 🖻 🏶 🕴 🎦	) 🚰 📓   🐰 🐂 🛍 🖅 • 💽 • 🚽 🕷 🔲	▼ ₹ X	1 🖬 🗃 🖏 🥾	Select data to plot	-
Name ▲	] ⊊   - 1.0 +   ÷ 1.1 ×   ‰ ‰		Name 🔺	Value	Min
advapi32.dll Dformd.dll			🛨 h	175.7065	175.7
Oforrt.dll h_ptx_D5.mexw32					
S LC.dll			٠	1	F
🚳 msvcp60.dll			Command History	<b>→</b> 1 🗖	× ×
🚳 msvcrt.dll			11/30/201	10 10:11 AM%	
		_			
		X N D			
fx >	>>				
h_ptx_D5.mexw32 (MEXW32 File)					
▲ <u>S</u> tart		script	L	Ln 1 Col 1 0	VR .::

#### Figure 2.10: Embedded "Editor" window

- Now type the following lines in the "Editor - Untitled" window:

Text to be written:	Explanation:			
% h_ptx_D5.m	file name as comment			
88	paragraph separation			
p=10; % pressure in bar	declaration of the			
t=300; % temperature in °C	variables pressure,			
x=-1; % vapor fraction in kg/kg	temperature, art and composition of mixture			
8 8	paragraph separation			
h=h_ptx_D5(p,t,x)	function call			
88	paragraph separation			

#### - Remarks:

- The program interprets the first line, starting with "%," to be a data description in "Current Directory."
- Paragraph separations which are mandatory are marked with "%%". This also serves to separate the declaration of variables and calculation instructions.
- The words which are printed in green, start with "%" and come after the variables are comments. They are not in fact absolutely necessary, but they are very helpful for your overview and to make the process more easily understood.
- Omit the semicolons after the numerical values if you wish to see the result for *h* and the input parameters.

The values of the function parameters in their corresponding units stand for:

- First operand: Value for p = 10(Range of validity: p = 0.00001 bar to 300 bar)
- Second operand: Value for t = 300 °C (Range of validity: t = 26.85 °C to 399.85 °C)
- Third operand: Value for x = -1 kg/kg

Since the wet steam region is calculated automatically by the subprograms, the following fixed details on the vapor fraction *x* are to be considered:

If the state point to be calculated is located in the single-phase region (liquid or superheated steam), e. g., pressure p and temperature t are given, the value –1 must be entered into the x cell as a pro-forma value.

In case, the state point to be calculated is located in the wet steam region, values between 0 and 1 have to be entered for x (the value 0 for boiling liquid, the value 1 for saturated steam).

Here, it is adequate to enter either the value given for t and p = -1, or the given value for p and t = -1, plus the value for x between 0 and 1.

However, if p and t and x are given when calculating wet steam, the program initially checks whether p and t meet the saturation-pressure curve. If this is not the case the enthalpy calculated later will result in -1000.

(D5 Saturation-pressure curve: t = 26.85 °C to  $t_{\rm C} = 346.0846$  °C

 $p_{\rm S}(26.85 \text{ °C}) = 0.001477917$  bar to  $p_{\rm C} = 11.6146$  bar)

- Save the "M-File" by clicking the "File" button and then click "Save As...".
- The menu "Save file as:" appears; In this menu, the folder name "LibD5\_Example" must be displayed in the "Save in:" field.
- Next to "File name" you have to type "Example\_h\_ptx\_D5.m" and afterwards click the "Save" button.

#### Note.

The name of the example file has to be different in comparison to the name of the used function. For example, the file could not be named "h\_ptx\_D5.m" in this case. Otherwise an error message will appear during the calculation.

- You will now see the following window:

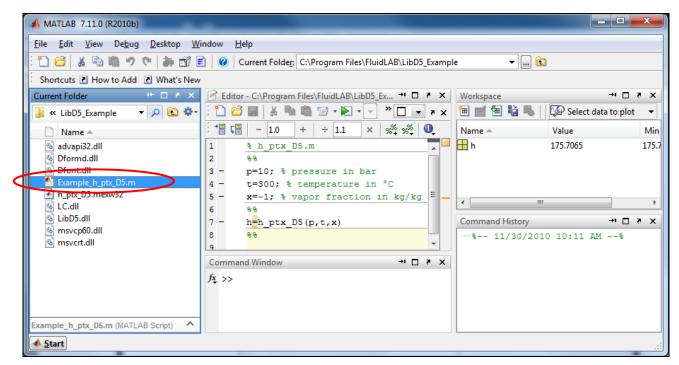


Figure 2.11: "Example\_h\_ptx\_D5.m" M-file

- Within the "Current Folder" window, the file "Example\_h\_ptx\_D5.m" appears.
- Right-click on this file and select "Run" in the menu which appears (see next image).

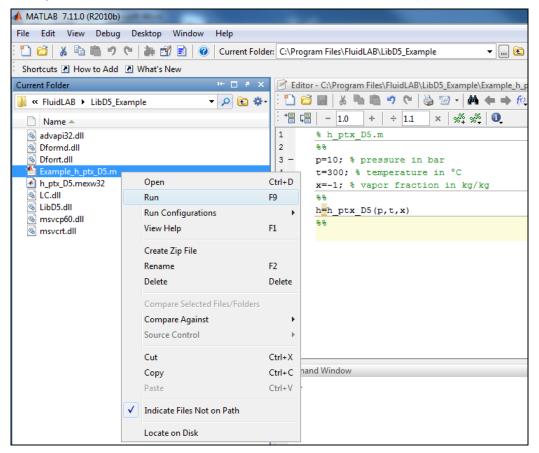


Figure 2.12: Running the "Example\_h\_ptx\_D5.m" M-file

- You will see the following window:

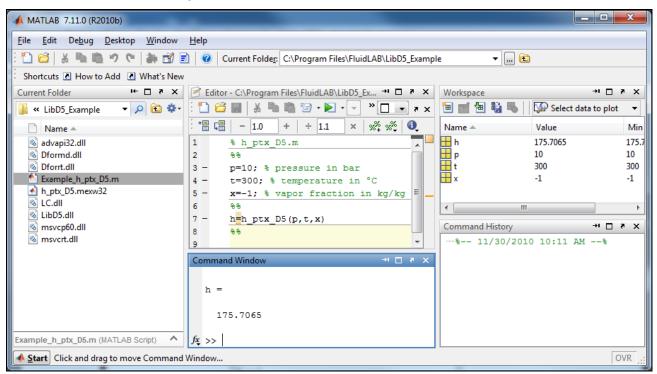


Figure 2.13: MATLAB® with calculated result

The result for h appears in the "Command Window".

 $\Rightarrow$  The result in our sample calculation here is: "h = 175.7065". The corresponding unit is kJ/kg (see table of the property functions in Chapter 1).

To be able to calculate other values, you have to copy the associated mexw32 files as well because MATLAB<sup>®</sup> can only access functions that are located in the "Current Directory" window. The example calculated can be found in the directory C:\Program Files\FluidLAB\LibD5\_Example," and you may use it as a basis for further calculations using FluidLAB.

#### 2.3 Example: Calculation of h = f(p, t, x) in the Command Window

- Start MATLAB<sup>®</sup> (if you have not started it already).
- Click the button marked in the following image in order to open the folder "\LibD5\_Example" in the window "Current Folder."

MATLAB 7.11.0 (R2010b)	É
<u>F</u> ile <u>E</u> dit De <u>b</u> ug <u>D</u> esktop <u>W</u> indow <u>H</u> elp	
🗄 🎦 🚰   👗 ங 🛍 🤊 🍽   🌦 🗊 🖹   🥝   Current Folde <u>r</u> : C:\Program Files\MATLAB	- 🛄 🖻 🖉
Shortcuts I How to Add I What's New	
Current Folder It 🛛 7 🗙 Command Window	
WATLAB ▶ ▼  P € \$\$ \$\$ \$\$ \$\$	
🗅 Name 🔺	
⊞ 📙 R2010b	

Figure 2.14: Selection of the working directory

- Find and select the directory "C:\Program Files\FluidLAB\LibD5\_Example" in the pop-up menu (see the following image).

Browse For Folder	×
Select a new folder	
DVD Maker	<b>^</b>
▷ 퉲 EES32	
Fingerprint Sensor	_
FluidLAB	
📗 LibD5	
LibD5_Example	-
•	F .
Eolder: LibD5_Example	
Make New Folder OK Ca	incel

Figure 2.15: Choosing the "LibD5\_Example" folder

- Confirm your selection by clicking the "OK" button.
- You will see the following window:

MATLAB 7.11.0 (R2010b)		
<u>File E</u> dit <u>V</u> iew De <u>b</u> ug <u>D</u> esktop <u>W</u> indow <u>H</u> elp		
🗄 🎦 🚰 🛛 😹 ங 🛍 🤊 🕅 🎥 🗊 🖹 🛛 🥑 Current Folder: C:\Program Files\FluidLAB\LibD5_Exam	nple 🔻 📖 🖻	
Shortcuts 🗷 How to Add 🗷 What's New		
Current Folder 🕛 र X Command Window 📲 🗖 र X	Workspace	× 5 ⊡ ++
📲 🤳 « LibD5_Example 🛛 🔻 🔎 🛍 🐲 🍂 >>	🛅 📷 🔁 🍓 🥾 🛛 💯 Sele	ct data to plot 🛛 👻
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Stormd.dll		
h_ptx_D5.mexw32		
🚳 LC.dll 🚳 LibD5.dll	<	
libbottini importanti in the second se	Command History	× 5 ⊡ ++
🚳 msvcrt.dll		AM%
Details		
Start Ready		
2 di neduj		

Figure 2.16: MATLAB<sup>®</sup> with necessary files

Corresponding to the table of the property functions in Chapter 1 you have to call up the function "**h\_ptx\_D5**" as follows for calculating h = f(p, t, x).

#### Write "h=h\_ptx\_D5(10,300,-1)" within the "Command Window"

The values of the function parameters in their corresponding units stand for:

- First operand: Value for p = 10 bar (Range of validity: p = 0.00001 bar to 300 bar)
- Second operand: Value for t = 300 °C(Range of validity: t = 26.85 °C to 399.85 °C)
- Third operand: Value for x = -1

Since the wet steam region is calculated automatically by the subprograms, the following fixed details on the vapor fraction *x* are to be considered:

If the state point to be calculated is located in the single-phase region (liquid or superheated steam), e. g., pressure p and temperature t are given, the value -1 must be entered into the x cell as a pro-forma value.

In case, the state point to be calculated is located in the wet steam region, values between 0 and 1 have to be entered for x (the value 0 for boiling liquid, the value 1 for saturated steam).

Here, it is adequate to enter either the value given for t and p = -1, or the given value for p and t = -1, plus the value for x between 0 and 1.

However, if p and t and x are given when calculating wet steam, the program initially checks whether p and t meet the saturation-pressure curve. If this is not the case the enthalpy calculated later will result in -1000.

(D5 Saturation-pressure curve: t = 26.85 °C to  $t_{\rm C} = 346.0846$  °C

 $p_{\rm S}(26.85 \text{ °C}) = 0.001477917$  bar to  $p_{\rm C} = 11.6146$  bar)

- Confirm your entry by pressing the "ENTER" button.
- You will see the following window:

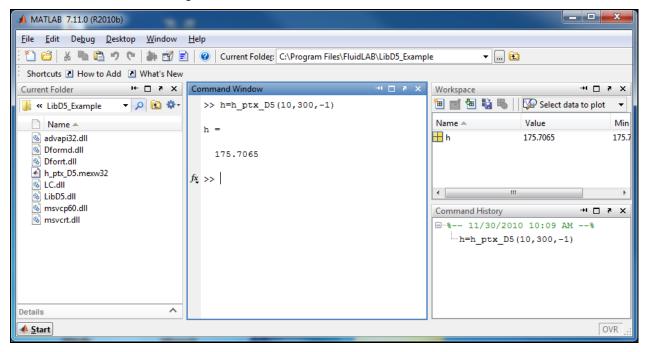


Figure 2.17: MATLAB® with calculated result

 $\Rightarrow$  In the "Command Window" you will see the result "h = 175.7065". The corresponding unit is kJ/kg (see table of the property functions in Chapter 1).

To be able to calculate other values, you will have to copy the respective mexw32 files into the working directory as well, because MATLAB<sup>®</sup> can only access functions that are located in the "Current Directory" window.

### 2.4 Using FluidLAB with SIMULINK

To use the functions of FluidLAB with the simulation program SIMULINK you have to start SIMULINK in MATLAB<sup>®</sup> by clicking on Simulink in the upper menu bar shown in Figure 2.19.

HOME	PLOTS	APPS				$\frown$				
New New Script Live Scrip	New Open	G Find Files	Import Save Data Workspace	<ul> <li>Bew Variable</li> <li>Dopen Variable ▼</li> <li>Open Variable ▼</li> <li>Clear Workspace ▼</li> </ul>	Analyze Code	<b>Pa</b> Simulink	(i) Preferences	Add-Ons	? Help	<ul> <li>➢ Community</li> <li>➢ Request Support</li> <li>☑ Learn MATLAB</li> </ul>
	FILE			VARIABLE	CODE	SIMULINK	ENVIRONMENT			RESOURCES
🗢 🌩 🔚 🔀	🚢 🕨 C: 🕨									
Current Folder		$\overline{\mathbf{v}}$	Command Wind	low						
🗋 Name 🔺			fx >>							

Figure 2.18: Starting Simulink

Then choose a blank model or a simulation in which you would like to use FluidLAB. Now you need to add a MATLAB function block that you can find in the library browser shown in Figure 2.20.

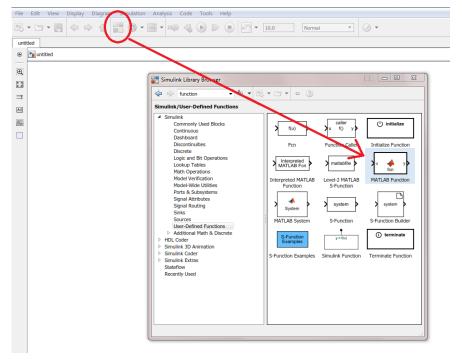


Figure 2.19: Simulink library browser and choosing a MATLAB Function

By dragging and dropping you can drag a Simulink block in your model. The function needs inputs and output that you can find in the Simulink library browser under sources and sinks. For this example constants were taken for the inputs and a display block were taken for outputting.

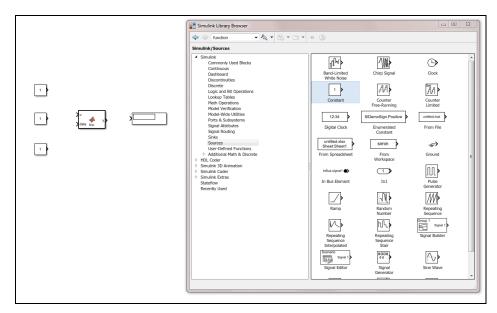
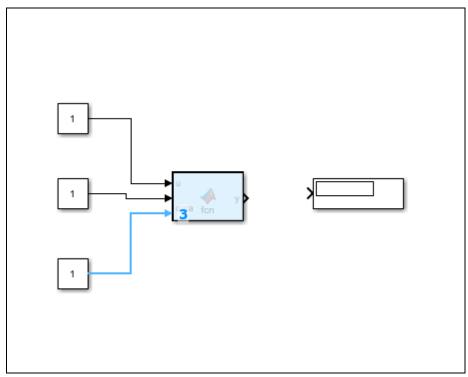
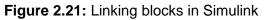


Figure 2.20: Inputs and outputs of the example

Now you have to link inputs and outputs to the MATLAB function block. By pressing and holding the left mouse button on the arrow of a block, you can draw a line and drag it to the MATLAB function block. With this method you can link all blocks together.





You can define the value of a constant block by double-click on them. If you want to calculate the example use the values you can find in section 2.3 and 2.4. With a double-click on the MATLAB function block you can define the function in MATLAB<sup>®</sup>. The following source code is for the example calculation and the table below describes the source code closer. You can adapt these few lines to call all other function of FluidLAB.

function h = fcn(p, t, x)

coder.extrinsic('addpath'); coder.extrinsic('h\_ptx\_D5'); addpath('C:\Program Files\FluidLAB\LibD5'); h = h ptx D5(p,t,x);

Matlab source code	Explanation
function $h = fcn(p, t, x)$	function header, you can define the function name and the inputs like p, t and x of the example
<pre>coder.extrinsic('addpath');</pre>	necessary to add a path
<pre>coder.extrinsic('h_ptx_D5');</pre>	Choose the function name of the FluidLAB function
<pre>addpath('C:\Program Files\FluidLAB\LibD5');</pre>	Add the installation path of FluidLAB
$h = h_ptx_D5(p,t,x);$	Linking the FluidLAB function to the MATLAB function block

You can copy and paste the sourcecode in MATLAB<sup>®</sup> or write it into the MATLAB<sup>®</sup> editor. The simulation will start by clicking the run button in Matlab or Simulink and you can see the example in the display block of the simulation which is shown in figure 2.23.

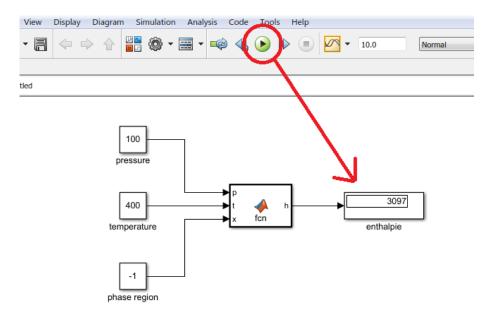


Figure 2.22: Starting the simulation and result of the calculation

Your result is may an other than shown in figure 2.22. If you want to calculate the example please use the values from section 2.3 and 2.4.

#### 2.5 Removing FluidLAB including LibD5

To remove the property library LibD5 from your hard disk drive in Windows<sup>®</sup>, click "Start" in the Windows<sup>®</sup> task bar, select "Settings" and click "Control Panel".

Now double-click on "Add or Remove Programs". In the list box of the "Add or Remove Programs" window that appears select "FluidLAB LibD5" by clicking on it and click the "Change/Remove" button.

In the following dialog box click "Automatic" and then click the "Next >" button.

Confirm the following menu "Perform Uninstall" by clicking the "Finish" button.

Finally, close the "Add or Remove Programs" and "Control Panel" windows. Now, FluidLAB has been removed.

If there is no library other than LibD5 installed, the directory "FluidLAB" will be removed as well.

## 3. Program Documentation

### Specific Isobaric Heat Capacity $c_p = f(p, t, x)$

Function Name:	cp_ptx_D5
Subroutine with function value: for call from Fortran	REAL*8 FUNCTION CPPTXD5(P,T,X) REAL*8 P,T,X
Subroutine with parameter: for call from DLL	INTEGER*4 FUNCTION C_CPPTXD5(CP,P,T,X) REAL*8 CP,P,T,X

#### Input Values:

**P** - Pressure *p* in bar

T - Temperature t in °C

X - Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

**CPPTXD5**, **CP** or **cp\_ptx\_D5** - specific isobaric heat capacity  $c_p$  in kJ/(kg K)

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}$ C to 399.85 °C
Pressure range:	from	p = 0.00001 bar to 300 bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located in the two phase region (wet steam), either the value 0 or 1 has to be entered for x (x = 0 for boiling liquid, x = 1 for saturated steam). The calculation for x values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1.

When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve. If it is not the case the calculation for the quantity of the chosen function to be calculated results in -1000.

Wet steam region: Temperature ranges from t = 26.85 °C to  $t_c = 346.0846$  °C

Pressure ranges from  $p_{\rm S}(26.85^{\circ}{\rm C}) = 0.001477917$  bar to  $p_{\rm C} = 11.6146$  bar

#### **Results for wrong input values**

Result CPPTXD5 = -1000, CP = -1000 or cp_ptx_D5 = -1000 for input values:					
Single phase region: ( <i>x</i> = -1)	p > 300 bar or $p < 0.00001$ bar or t > 399.85 °C or $t < 26.85$ °C				
Boiling or dew curve:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C				

References: [1]

### Specific Isochoric Heat Capacity $c_v = f(p, t, x)$

Function Name:	cv_ptx_D5
Subroutine with function value: for call from Fortran	REAL*8 FUNCTION CVPTXD5(P,T,X) REAL*8 P,T,X
Subroutine with parameter: for call from DLL	INTEGER*4 FUNCTION C_CVPTXD5(CV,P,T,X) REAL*8 CV,P,T,X

#### Input Values:

- P Pressure p in bar
- T Temperature t in °C
- X Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

CVPTXD5, CV or  $cv_ptx_D5$  - specific isochoric heat capacity  $c_v$  in kJ/(kg K)

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}C$	to	399.85 °C
Pressure range:	from	p = 0.00001	bar	to 300 bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located in the two phase region (wet steam), either the value 0 or 1 has to be entered for x (x = 0 for boiling liquid, x = 1 for saturated steam). The calculation for x values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1.

When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve. If it is not the case the calculation for the quantity of the chosen function to be calculated results in -1000.

Wet steam region:	Temperature ranges from $t = 26.85$ °C to $t_{c} = 346.0846$ °C
	Pressure ranges from $p_{s}(26.85^{\circ}C) = 0.001477917$ bar to $p_{c} = 11.6146$ bar

#### **Results for wrong input values**

Result CVPTXD5 = -1000, CV = -1000 or cv_ptx_D5 = -1000 for input values:			
Single phase region: ( <i>x</i> = -1)	<i>p</i> > 300 bar or <i>p</i> < 0.00001 bar or <i>t</i> > 399.85 °C or <i>t</i> < 26.85°C		
Boiling or dew curve:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C		

### Derivative of Pressure with Respect to Temperature (at

**Constant Specific Volume)** 

 $\frac{\partial \boldsymbol{p}}{\partial \boldsymbol{T}} = f(\boldsymbol{p}, \boldsymbol{t}, \boldsymbol{x})$ 

REAL\*8 DPDTV,P,T,X

Function Name:

#### dpdtv\_ptx\_D5

Subroutine with function value: for call from Fortran

REAL\*8 FUNCTION DPDTVPTXD5(P,T,X) REAL\*8 P,T,X

INTEGER\*4 FUNCTION C DPDTVPTXD5(DPDTV,P,T,X)

Derivative of pressure with respect to temperature (at constant specific volume) dpdtv in kPa/K

Subroutine with parameter: for call from DLL

#### **Input Values:**

- P Pressure p in bar
- T Temperature t in °C
- X Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}$ C to 399.85 °C
Pressure range:	from	p = 0.00001 bar to 300 bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located on the boiling curve, x = 0 must be entered. When calculating saturated steam (dew curve) x = 1 is entered as given value. The calculation for *x* values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1. When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve.

Boiling and dew curve:	Temperature ranges from $t = 26.85 \text{ °C}$ to $t_{c} = 346.0846 \text{ °C}$
	Pressure ranges from $p_{s}(26.85^{\circ}C) = 0.001477917$ bar to $p_{c} = 11.6146$ bar

#### Results for wrong input values

Result DPDTVPTXD5 = -1	000, DPDTV = -1000 or dpdtv_ptx_D5 = -1000 for input values:
Single phase region: ( <i>x</i> = -1)	p > 300 bar or $p < 0.00001$ bar or t > 399.85 °C or $t < 26.85$ °C
Boiling or dew curve:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C

References: [1]

### Derivative of Pressure with Respect to Specific Volume (at

### **Constant Temperature)**

 $\frac{\partial \boldsymbol{p}}{\partial \boldsymbol{v}} \bigg|_{\boldsymbol{\tau}} = \mathbf{f}(\boldsymbol{p}, \boldsymbol{t}, \boldsymbol{x})$ 

Function Name:

#### dpdvt\_ptx\_D5

Subroutine with function value: for call from Fortran

REAL\*8 FUNCTION DPDVTPTXD5(P,T,X) REAL\*8 P,T,X INTEGER\*4 FUNCTION C DPDVTPTXD5(DPDVT,P,T,X)

Subroutine with parameter: for call from DLL

#### **Input Values:**

- **P** Pressure *p* in bar
- **T** Temperature *t* in °C
- **X** Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

DPDVTPTXD4, DPDVT or dpdvt_ptx_D4 -	Derivative of pressure with respect to temperature		
	(at constant specific volume) dpdvt in kPa/(m <sup>3</sup> /kg)		

REAL\*8 DPDVT,P,T,X

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}$ C to 399.85 °C
Pressure range:	from	p = 0.00001 bar to 300 bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located on the boiling curve, x = 0 must be entered. When calculating saturated steam (dew curve) x = 1 is entered as given value. The calculation for *x* values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1. When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve.

Wet steam region:	Temperature ranges from $t = 26.85 \text{ °C}$ to $t_c = 346.0846 \text{ °C}$
	Pressure ranges from $p_{\rm s}(26.85^{\circ}{\rm C}) = 0.001477917$ bar to $p_{\rm c} = 11.6146$ bar

#### **Results for wrong input values**

Result <b>DPDVTPTXD5 = -1000</b> , <b>DPDVT = -1000</b> o	r dpdvt_ptx_D5 = -1000	for input values:
---	------------------------	-------------------

Single phase region:	<i>p</i> > 300 bar or <i>p</i> < 0.00001 bar or
( <i>x</i> = -1)	<i>t</i> > 399.85 °C or <i>t</i> < 26.85°C
Boiling or dew curve:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C

### Specific Enthalpy h = f(p, t, x)

Function Name:

Subroutine with function value: for call from Fortran

Subroutine with parameter: for call from DLL

#### **Input Values:**

**P** - Pressure *p* in bar

**T** - Temperature *t* in °C

**X** - Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

HPTXD5, H or h\_ptx\_D5 - specific enthalpy h in kJ/kg

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}C$	to	399	9.85	°C
Pressure range:	from	p = 0.00001	bar	to	300	bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

h\_ptx\_D5

REAL\*8 P,T,X

REAL\*8 H,P,T,X

REAL\*8 FUNCTION HPTXD5(P,T,X)

INTEGER\*4 FUNCTION C\_HPTXD5(H,P,T,X)

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.

If the state point to be calculated is located in the two phase region (wet steam), either the value 0 or 1 has to be entered for x (x = 0 for boiling liquid, x = 1 for saturated steam). The calculation for x values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1.

When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve. If it is not the case the calculation for the quantity of the chosen function to be calculated results in -1000.

Wet steam region:	Temperature ranges from $t = 26.85$ °C to $t_c = 346.0846$ °C
	Pressure ranges from $p_{s}(26.85^{\circ}C) = 0.001477917$ bar to $p_{c} = 11.6146$ bar

#### Results for wrong input values

Result **HPTXD5 = -1000**, **H = -1000** or **h\_ptx\_D5 = -1000** for input values:

Single phase region:	<i>p</i> > 300 bar or <i>p</i> < 0.00001 bar or
( <i>x</i> = -1)	<i>t</i> > 399.85 °C or <i>t</i> < 26.85°C
Boiling or dew curve:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C

### Isentropic Exponent $\kappa = f(p, t, x)$

Function Name:

kappa\_ptx\_D5

REAL\*8 KAPPA, P, T, X

Subroutine with function value: for call from Fortran

REAL\*8 FUNCTION KAPPAPTXD5(P,T,X) REAL\*8 P,T,X

INTEGER\*4 FUNCTION C\_KAPPAPTXD5(KAPPA,P,T,X)

Subroutine with parameter: for call from DLL

#### **Input Values:**

- P Pressure p in bar
- **T** Temperature *t* in °C
- X Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

**KAPPAPTXD5**, **KAPPA** or **kappa\_ptx\_D5** - Isentropic exponent  $\kappa = \frac{w^2}{p \cdot v}$ 

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}C$	to	399	).85	°C
Pressure range:	from	p = 0.00001	bar	to	300	bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located on the boiling curve, x = 0 must be entered. When calculating saturated steam (dew curve) x = 1 is entered as given value. The calculation for *x* values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1. When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve.

Boiling and dew curve: Temperature ranges from t = 26.85 °C to  $t_c = 346.0846$  °C Pressure ranges from  $p_s(26.85$ °C) = 0.001477917 bar to  $p_c = 11.6146$  bar

#### **Results for wrong input values**

Result KAPPAPTXD5, KAPPA = -1000 or kappa\_ptx\_D5 = -1000 for input values:

Single phase region:	<i>p</i> > 300 bar or <i>p</i> < 0.00001 bar or
( <i>x</i> = -1)	<i>t</i> > 399.85 °C or <i>t</i> < 26.85°C
Boiling or dew curve:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C

#### References: [1]

### Vapor Pressure $p_s = f(t)$

Function Name:

ps\_t\_D5

REAL\*8 PS,T

Subroutine with function value: for call from Fortran

REAL\*8 FUNCTION PSTD5(T) REAL\*8 T INTEGER\*4 FUNCTION C\_PSTD5(PS,T)

Subroutine with parameter: for call from DLL

#### **Input Values:**

T - Temperature t in °C

#### Result

**PSTD5**, **PS** or **ps\_t\_D5** - Vapor pressure  $p_s$  in bar

#### Range of validity

Temperature range: from t = 26.85 °C to  $t_c = 346.0846$  °C

#### **Results for wrong input values**

Result **PSTD5 = -1000**, **PS = -1000** or **ps\_t\_D5 = -1000** for input values:

 $t = 26.85 \text{ °C to } t_{\text{c}} = 346.0846 \text{ °C}$ 

References: [1]

### Density $\rho = f(p, t, x)$

Function Name:

Subroutine with function value: for call from Fortran

Subroutine with parameter: for call from DLL

#### **Input Values:**

**P** - Pressure *p* in bar

**T** - Temperature *t* in °C

X - Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

**RHO\_PTX\_D5**, **RHO** or **rho\_ptx\_D5** - Density  $\rho$  in kg/m<sup>3</sup>

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}C$	to	399.	.85	°C
Pressure range:	from	<i>p</i> = 0.00001	bar	to	300	bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located in the two phase region (wet steam), either the value 0 or 1 has to be entered for x (x = 0 for boiling liquid, x = 1 for saturated steam). The calculation for x values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1.

When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve. If it is not the case the calculation for the quantity of the chosen function to be calculated results in -1000.

Wet steam region:Temperature ranges from t = 26.85 °C to  $t_c = 346.0846$  °CPressure ranges from  $p_s(26.85$ °C) = 0.001477917 bar to  $p_c = 11.6146$  bar

#### **Results for wrong input values**

Result RHOPTXD5 = -1000, RHO = -1000 or rho_ptx_D5 = -1000 for input values:			
Single phase region: ( <i>x</i> = -1)	p > 300 bar or $p < 0.00001$ bar or t > 399.85 °C or $t < 26.85$ °C		
Wet steam region:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C		

rho\_ptx\_D5

REAL\*8 FUNCTION RHOPTXD5(P,T,X) REAL\*8 P,T,X

**INTEGER\*4 FUNCTION C\_RHOPTXD5(RHO,P,T,X)** REAL\*8 RHO,P,T,X

### Specific Entropy s = f(p, t, x)

**Function Name:** 

Subroutine with function value: for call from Fortran

Subroutine with parameter: for call from DLL

REAL\*8 FUNCTION SPTXD5(P,T,X) REAL\*8 P,T,X INTEGER\*4 FUNCTION C\_SPTXD5(S,P,T,X) REAL\*8 S,P,T,X

s\_ptx\_D5

#### **Input Values:**

P - Pressure p in bar

T - Temperature t in °C

**X** - Vapor fraction *x* (kg of saturated steam)/(kg wet steam)

#### Result

SPTXD5, S or s\_ptx\_D5 - Specific entropy s in kJ/kg K

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}$ C to 399.85 °C
Pressure range:	from	p = 0.00001 bar to 300 bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located in the two phase region (wet steam), either the value 0 or 1 has to be entered for x (x = 0 for boiling liquid, x = 1 for saturated steam). The calculation for x values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1.

When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve. If it is not the case the calculation for the quantity of the chosen function to be calculated results in -1000.

Wet steam region:Temperature ranges from t = 26.85 °C to  $t_c = 346.0846$  °CPressure ranges from  $p_s(26.85$ °C) = 0.001477917 bar to  $p_c = 11.6146$  bar

#### Results for wrong input values

Result <b>SPTXD5 = -1000</b> , <b>S = -1000</b> or <b>s_ptx_D5 = -1000</b> for input values:			
Single phase region: ( <i>x</i> = -1)	<i>p</i> > 300 bar or <i>p</i> < 0.00001 bar or <i>t</i> > 399.85 °C or <i>t</i> < 26.85°C		
Boiling or dew curve:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C		

**References:** [1]

# Backward Function: Temperature t = f(p,h)

Function Name:

Subroutine with function value: for call from Fortran

Subroutine with parameter: for call from DLL

t\_ph\_D5 REAL\*8 FUNCTION TPHD5(P,H) REAL\*8 P,H INTEGER\*4 FUNCTION C\_TPHD5(T,P,H) REAL\*8 T,P,H

#### **Input Values:**

**P** - Pressure *p* in bar

H - Specific enthalpy h in kJ/kg

#### Result

TPHD5, T or t\_ph\_D5 - Temperature t in °C

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}C$	to	399.85 °C
Pressure range:	from	p = 0.00001	bar	to 300 bar

#### Details on the calculation of wet steam

The wet steam region is calculated automatically. That means the given values of p and h are taken as a basis and the subprogram will determine whether the state point to be calculated is located within the single-phase region (liquid or steam) or the wet steam region. Afterwards the calculation of the appropriate state region will be carried out.

Wet steam region : Pr essure ranges from  $p_s(26.85^{\circ}C) = 0.001477917$  bar to  $p_c = 11.6146$  bar

#### **Results for wrong input values**

Result **T\_PH\_D5**, **T = -1000** or **t\_ph\_D5 = -1000** for input values:

Single phase region:	<i>p</i> > 300 bar or <i>p</i> < <i>p</i> <sub>S</sub> (26.85°C) = 0.001477917 bar or
( <i>x</i> = -1)	at result <i>t</i> > 399.85 °C, <i>t</i> < 26.85°C
Boiling or dew curve:	at $p > p_c = 11.6146$ bar or $p < p_s(26.85^{\circ}C) = 0.001477917$ bar and at result $t > t_c = 346.0846^{\circ}C$ or $t < 26.85^{\circ}C$

# Backward Function: Temperature t = f(p, s)

Function Name:

Subroutine with function value: for call from Fortran

REAL\*8 FUNCTION TPSD5(P,S) REAL\*8 P,S INTEGER\*4 FUNCTION C\_TPSD5(T,P,S) REAL\*8 T,P,S

t\_ps\_D5

#### **Input Values:**

for call from DLL

P - Pressure p in bar

Subroutine with parameter:

S - Specific entropy s in kJ/(kg K)

#### Result

TPSD5, T or t\_ps\_D5 - Temperature t in °C

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}C$	to	399	).85 °C	С
Pressure range:	from	p = 0.00001	bar	to	300 b	bar

#### Details on the calculation of wet steam

The wet steam region is calculated automatically. That means the given values of p and s are taken as a basis and the subprogram will determine whether the state point to be calculated is located within the single-phase region (liquid or steam) or the wet steam region. Afterwards the calculation of the appropriate state region will be carried out.

Wet steam region : Pr essure ranges from  $p_s(26.85^{\circ}C) = 0.001477917$  bar to  $p_c = 11.6146$  bar

#### **Results for wrong input values**

#### Result **T\_PS\_D5**, **T = -1000** or **t\_ps\_D5 = -1000** for input values:

Single phase region: ( <i>x</i> = -1)	$p > 300$ bar or $p < p_{s}(26.85^{\circ}C) = 0.001477917$ bar or at result $t > 399.85^{\circ}C$ , $t < 26.85^{\circ}C$
Boiling or dew curve:	at $p > p_c = 11.6146$ bar or $p < p_s(26.85^{\circ}C) = 0.001477917$ bar and at result $t > t_c = 346.0846^{\circ}C$ or $t < 26.85^{\circ}C$

# Boiling Temperature $t_s = f(p)$

Function Name:	ts_p_D5
Subroutine with function value: for call from Fortran	REAL*8 FUNCTION TSPD5(P) REAL*8 P
Subroutine with parameter: for call from DLL	INTEGER*4 FUNCTION C_TSPD5(TS,P) REAL*8 TS,P
Input Values:	
<b>P</b> - Pressure <i>p</i> in bar	

#### Result

**TSPD5**, **TS** or **ts\_p\_D5** - Boiling temperature  $t_s$  in °C

#### Range of validity

Pressure ranges: from  $p_{\rm S}(26.85^{\circ}{\rm C}) = 0.001477917$  bar to  $p_{\rm C} = 11.6146$  bar

#### **Results for wrong input values**

Result **TSPD5 = -1000**, **TS = -1000** or **ts\_p\_D5 = -1000** for input values:

 $p < p_{\rm S}(26.85^{\circ}{\rm C}) = 0.001477917$  bar or  $p > p_{\rm C} = 11.6146$  bar

# Specific Internal Energy *u* = f(*p*,*t*,*x*)

#### Function Name:

Subroutine with function value: for call from Fortran

Subroutine with parameter: for call from DLL

# INTEGER\*4 FUNCTION C\_UPTXD5(U,P,T,X) REAL\*8 U,P,T,X

**REAL\*8 FUNCTION UPTXD5(P,T,X)** 

u\_ptx\_D5

REAL\*8 P,T,X

#### **Input Values:**

- P Pressure p in bar
- T Temperature t in °C
- X Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

UPTXD5, U or u\_ptx\_D5 - Specific internal energy u in kJ/kg

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}C$	to	399	).85	°C
Pressure range:	from	<i>p</i> = 0.00001	bar	to	300	bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located in the two phase region (wet steam), either the value 0 or 1 has to be entered for x (x = 0 for boiling liquid, x = 1 for saturated steam). The calculation for x values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1.

When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve. If it is not the case the calculation for the quantity of the chosen function to be calculated results in -1000.

Wet steam region:Temperature ranges from t = 26.85 °C to  $t_c = 346.0846$  °CPressure ranges from  $p_s(26.85$ °C) = 0.001477917 bar to  $p_c = 11.6146$  bar

#### Results for wrong input values

Result <b>UPTXD5 = -1000</b> , <b>U</b>	J = -1000 or u_ptx_D5 = -1000 for input values:
Single phase region: ( <i>x</i> = -1)	<i>p</i> > 300 bar or <i>p</i> < 0.00001 bar or <i>t</i> > 399.85 °C or <i>t</i> < 26.85°C
Wet steam region:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C

# Specific Volume v = f(p,t,x)

**Function Name:** 

Subroutine with function value: for call from Fortran

REAL\*8 FUNCTION VPTXD5(P,T,X) REAL\*8 P,T,X

INTEGER\*4 FUNCTION C\_VPTXD5(V,P,T,X)

v\_ptx\_D5

REAL\*8 V,P,T,X

Subroutine with parameter: for call from DLL

#### **Input Values:**

**P** - Pressure *p* in bar

**T** - Temperature *t* in °C

X - Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

**VPTXD5**, **V** or **v\_ptx\_D5** - Specific volume v in m<sup>3</sup>/kg

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}C$	to	399	9.85 °	°C
Pressure range:	from	p = 0.00001	bar	to	300	bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located in the two phase region (wet steam), either the value 0 or 1 has to be entered for x (x = 0 for boiling liquid, x = 1 for saturated steam). The calculation for x values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1.

When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve. If it is not the case the calculation for the quantity of the chosen function to be calculated results in -1000.

Wet steam region: Temperature ranges from t = 26.85 °C to  $t_c = 346.0846$  °C Pressure ranges from  $p_s(26.85$ °C) = 0.001477917 bar to  $p_c = 11.6146$  bar

#### **Results for wrong input values**

Single phase region:	<i>p</i> > 300 bar or <i>p</i> < 0.00001 bar or
( <i>x</i> = -1)	<i>t</i> > 399.85 °C or <i>t</i> < 26.85°C
Wet steam region:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C

# Isentropic Speed of Sound w = f(p, t, x)

**Function Name:** 

Subroutine with function value: for call from Fortran

REAL\*8 FUNCTION WPTXD5(P,T,X) REAL\*8 P,T,X INTEGER\*4 FUNCTION C\_WPTXD5(W,P,T,X)

w\_ptx\_D5

REAL\*8 W,P,T,X

Subroutine with parameter: for call from DLL

#### **Input Values:**

- **P** Pressure *p* in bar
- **T** Temperature *t* in °C
- X Vapor fraction x (kg of saturated steam)/(kg wet steam)

#### Result

WPTXD5, W or w\_ptx\_D5 - Speed of sound w in m/s

#### Range of validity

Temperature range:	from	$t = 26.85^{\circ}C$	to	399.85 °C
Pressure range:	from	p = 0.00001	bar	to 300 bar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located on the boiling curve, x = 0 must be entered. When calculating saturated steam (dew curve) x = 1 is entered as given value. The calculation for *x* values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1. When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve.

Boiling and dew curve: Temperature ranges from t = 26.85 °C to  $t_c = 346.0846$  °C Pressure ranges from  $p_s(26.85$ °C) = 0.001477917 bar to  $p_c = 11.6146$  bar

#### Results for wrong input values

Result WPTXD5 = -1000, W = -1000 or w\_ptx\_D5 = -1000 for input values:

Single phase region:	p > 300 bar or $p < 0.00001$ bar or
(x = -1)	t > 399.85 °C or $t < 26.85$ °C
Boiling or dew curve:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C

# Backward Function: Vapor Fraction x = f(p,h)

Function Name:x\_ph\_D5Subroutine with function value:<br/>for call from FortranREAL\*8 FUNCTION XPHD5(P,H)<br/>REAL\*8 P,HSubroutine with parameter:<br/>for call from DLLINTEGER\*4 FUNCTION C\_XPHD5(X,P,H)<br/>REAL\*8 X,P,H

#### **Input Values:**

**P** - Pressure *p* in bar

H - Specific enthalpy h in kJ/kg

#### Result

XPHD5, X or x\_ph\_D5 - Vapor fraction x in (kg saturated steam/kg wet steam)

#### Range of validity

Temperature range:	from $t = 26.85^{\circ}$ C to $t_{c} = 346.0846^{\circ}$ C
Pressure range:	from $p_{\rm s}(26.85^{\circ}{\rm C}) = 0.001477917$ bar to $p_{\rm c} = 11.6146$ bar

#### Details on the calculation of wet steam

The wet steam region is calculated automatically. That means the given values of p and h are taken as a basis and the subprogram will determine whether the state point to be calculated is located within the single-phase region (liquid or superheated steam) or the wet steam region. In case of wet steam, x will be calculated, otherwise the result is set to x = -1.

Wet steam region : Pr essure ranges from  $p_s(26.85^{\circ}C) = 0.001477917$  bar to  $p_c = 11.6146$  bar

#### Results for wrong input values

Result X\_PH\_D5 = -1, X = -1 or x\_ph\_D5 = -1 for input values:

If the state point is located in the single phase region:  $p > p_c = 11.6146$  bar or p < 0.001477917 bar

# Backward Function: Vapor Fraction x = f(p, s)

Function Name:

Subroutine with function value: for call from Fortran

Subroutine with parameter: for call from DLL

x\_ps\_D5 REAL\*8 FUNCTION XPSD5(P,S) REAL\*8 P,S INTEGER\*4 FUNCTION C\_XPSD5(X,P,S) REAL\*8 X,P,S

#### **Input Values:**

- P Pressure p in bar
- S Specific entropy s in kJ/(kg K)

#### Result

XPSD5, X or x\_ps\_D5 - Vapor fraction x in (kg saturated steam/kg wet steam)

#### Range of validity

Temperature range:	from $t = 26.85^{\circ}$ C to $t_{c} = 346.0846^{\circ}$ C
Pressure range:	from $p_{\rm s}(26.85^{\circ}{\rm C}) = 0.001477917$ bar to $p_{\rm c} = 11.6146$ bar

#### Details on the calculation of wet steam

The wet steam region is calculated automatically. That means the given values of p and h are taken as a basis and the subprogram will determine whether the state point to be calculated is located within the single-phase region (liquid or superheated steam) or the wet steam region. In case of wet steam, x will be calculated, otherwise the result is set to x = -1.

Wet steam region : Pr essure ranges from  $p_s(26.85^{\circ}C) = 0.001477917$  bar to  $p_c = 11.6146$  bar

#### Results for wrong input values

Result  $X_PS_D5$ , X = -1 or  $x_ps_D5 = -1$  for input values:

If the state point is located in the single phase region:  $p > p_c = 11.6146$  bar or  $p < p_s(26.85^{\circ}C) = 0.001477917$  bar

# Compression Factor Z = f(p, t, x)

Function Name:

Subroutine with function value: for call from Fortran

Subroutine with parameter: for call from DLL

REAL\*8 FUNCTION ZPTXD5(P,T,X) REAL\*8 P,T,X INTEGER\*4 FUNCTION C\_ZPTXD5(Z,P,T,X) REAL\*8 Z,P,T,X

Z\_ptx\_D5

#### **Input Values:**

- P Pressure p in bar
- T Temperature t in °C
- **X** Vapor fraction *x* (kg of saturated steam)/(kg wet steam)

#### Result

ZPTXD5, Z or Z\_ptx\_D5 - Speed of sound w in m/s

#### Range of validity

Temperature range:	from $t = 26.85^{\circ}$ C to 399.85 °C	;
Pressure range:	from $p = 0.00001$ bar to 300 b	ar

#### Details on the vapor fraction x and on the calculation of wet steam

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:

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.

If the state point to be calculated is located on the boiling curve, x = 0 must be entered. When calculating saturated steam (dew curve) x = 1 is entered as given value. The calculation for *x* values between 0 and 1 is not possible.

If the state point to be calculated is located in the two phase region, it is adequate to enter either the given value for t and p = -1000, or the given value for p and t = -1000, plus the value for x between 0 and 1. When calculating wet steam and p and t and x are entered as given values, the program will consider p and t to be appropriate to represent the saturation-pressure curve.

Boiling and dew curve: Temperature ranges from t = 26.85 °C to  $t_c = 346.0846$  °C Pressure ranges from  $p_s(26.85$ °C) = 0.001477917 bar to  $p_c = 11.6146$  bar

#### Results for wrong input values

Result **ZPTXD5 = -1000**, **Z = -1000** or **Z\_ptx\_D5 = -1000** for input values:

Single phase region:	p > 300 bar or $p < 0.00001$ bar or
( <i>x</i> = -1)	t > 399.85 °C or $t < 26.85$ °C
Boiling or dew curve:	at $p = -1000$ and $t > t_c = 346.0846$ °C or $t < 26.85$ °C at $t = -1000$ and $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar or at $p > p_c = 11.6146$ bar or $p < p_s(26.85$ °C) = 0.001477917 bar and $t > t_c = 346.0846$ °C or $t < 26.85$ °C





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# **Property Libraries** for Calculating Heat Cycles, Boilers, Turbines and Refrigerators

#### Water and Steam

#### Library LibIF97

- Industrial Formulation
   Library LibSBIL\_95
   IAPWS-IF97 (Revision 2007) Extremely fast property calculations
- Supplementary Standards
   IAPWS-IF97-S01
  - IAPWS-IF97-S03rev
  - IAPWS-IF97-S04
- IAPWS-IF97-S05
- IAPWS Revised Advisory Note No. 3 on Thermodynamic Derivatives (2008)

# Library LibSBTL\_IF97

# Library LibSBTL\_95

Extremely fast property calculations according to the IAPWS Guideline 2015 Spline-based Table Look-up Method (SBTL) applied to the Industrial Formulation IAPWS-IF97 and to the Scientific Formulation IAPWS-95

for Computational Fluid Dynamics and simulating non-stationary processes

#### **Humid Combustion Gas Mixtures**

#### Library LibHuGas

Model: Ideal mixture of the real fluids:

 $CO_2$  - Span, Wagner H<sub>2</sub>O - IAPWS-95 O<sub>2</sub> - Schmidt, Wagner N<sub>2</sub> - Span et al. Ar - Tegeler et al. and of the ideal gases: SO<sub>2</sub>, CO, Ne

(Scientific Formulation of Bücker et al.) Consideration of:

- Dissociation from VDI 4670
- Poynting effect

#### **Humid Air**

#### Library LibHuAir

Model: Ideal mixture of the real fluids:

- Dry air from Lemmon et al.
  Steam, water and ice from
- IAPWS-IF97 and IAPWS-06

Consideration of:

- Condensation and freezing of steam
- Dissociation from VDI 4670Poynting effect from
- ASHRAE RP-1485

#### Carbon Dioxide Including Dry Ice

#### Library LibCO2

Formulation of Span and Wagner (1996)

#### Seawater

#### Library LibSeaWa

IAPWS Industrial Formulation 2013

#### lce

#### Library LibICE

Ice from IAPWS-06, Melting and sublimation pressures from IAPWS-08, Water from IAPWS-IF97, Steam from IAPWS-95 and -IF97

# Ideal Gas Mixtures

#### Library LibldGasMix

Model: Ideal mixture of the ideal gases:

Ar	NO	He	Propylene
Ne	H <sub>2</sub> O	F <sub>2</sub>	Propane
N <sub>2</sub>	SO <sub>2</sub>	NH <sub>3</sub>	Iso-Butane
<b>O</b> <sub>2</sub>	H <sub>2</sub>	Methane	n-Butane
CO	H₂S	Ethane	Benzene
CO <sub>2</sub>	ОН	Ethylene	Methanol
Air			

Consideration of: • Dissociation from the VDI Guideline 4670

#### Library LibIDGAS

Model: Ideal gas mixture from VDI Guideline 4670

Consideration of: • Dissociation from the VDI Guideline 4670

## Humid Air

#### Library ASHRAE LibHuAirProp

Model: Virial equation from ASHRAE Report RP-1485 for real mixture of the real fluids: - Dry air

- Steam
- Consideration of:
- Enhancement of the partial saturation pressure of water vapor at elevated total pressures

www.ashrae.org/bookstore

### Dry Air Including Liquid Air Library LibRealAir

Formulation of Lemmon et al. (2000)

#### Refrigerants

#### Ammonia

# Library LibNH3

Formulation of Tillner-Roth et al. (1993)

R134a

#### Library LibR134a

Formulation of Tillner-Roth and Baehr (1994)

#### Iso-Butane

#### Library LibButane\_Iso

Formulation of Bücker and Wagner (2006)

### n-Butane

#### Library LibButane\_n

Formulation of Bücker and Wagner (2006)

#### Mixtures for Absorption Processes

#### Ammonia/Water Mixtures

#### Library LibAmWa

IAPWS Guideline 2001 of Tillner-Roth and Friend (1998) Helmholtz energy equation for the mixing term (also useable for calculating the Kalina Cycle)

#### Water/Lithium Bromide Mixtures

#### Library LibWaLi

Formulation of Kim and Infante Ferreira (2004) Gibbs energy equation for the mixing term

#### Liquid Coolants

#### Liquid Secondary Refrigerants

#### Library LibSecRef

Liquid solutions of water with		
$C_2H_6O_2$	Ethylene glycol	
$C_3H_8O_2$	Propylene glycol	
C₂H₅OH	Ethanol	
CH₃OH	Methanol	
C <sub>3</sub> H <sub>8</sub> O <sub>3</sub>	Glycerol	
K <sub>2</sub> CO <sub>3</sub>	Potassium carbonate	
CaCl <sub>2</sub>	Calcium chloride	
MgCl <sub>2</sub>	Magnesium chloride	
NaCl	Sodium chloride	
$C_2H_3KO_2$	Potassium acetate	
CHKO <sub>2</sub>	Potassium formate	
LiCl	Lithium chloride	
NH <sub>3</sub>	Ammonia	

Formulation of the International Institute of Refrigeration (IIR 2010)

#### Ethanol

Library LibC2H5OH

Formulation of Schroeder (2012)

#### Methanol Library LibCH3OH

Formulation of de Reuck and Craven (1993)

### Propane Library LibPropane

Formulation of Lemmon et al. (2009)

#### Siloxanes as ORC Working Fluids

Octamethylcyclotetrasiloxane  $C_8H_{24}O_4Si_4$  Library LibD4 Decamethylcyclopentasiloxane  $C_{10}H_{30}O_5Si_5$  Library LibD5 Tetradecamethylhexasiloxane  $C_{14}H_{42}O_5Si_6$  Library LibMD4M Hexamethyldisiloxane  $C_6H_{18}OSi_2$  Library LibMM Formulation of Colonna et al. (2006)

Dodecamethylcyclohexasiloxane  $C_{12}H_{36}O_6Si_6$  Library LibD6 Decamethyltetrasiloxane  $C_{10}H_{30}O_3Si_4$  Library LibMD2M Dodecamethylpentasiloxane  $C_{12}H_{36}O_4Si_5$  Library LibMD3M Octamethyltrisiloxane  $C_8H_{24}O_2Si_3$  Library LibMDM Formulation of Colonna et al. (2008)

#### Nitrogen and Oxygen Libraries LibN2 and LibO2

Formulations of Span et al. (2000) and Schmidt and Wagner (1985)

#### Hydrogen Library LibH2

Formulation of Leachman et al. (2009)

#### Helium

Library LibHe

Formulation of Arp et al. (1998)

#### Hydrocarbons

Decane  $C_{10}H_{22}$  Library LibC10H22 Isopentane  $C_5H_{12}$  Library LibC5H12\_ISO Neopentane  $C_5H_{12}$  Library LibC5H12\_NEO Isohexane  $C_6H_{14}$  Library LibC6H14 Toluene  $C_7H_8$  Library LibC7H8 Formulation of Lemmon and Span (2006)

#### **Further Fluids**

Carbon monoxide CO Library LibCO Carbonyl sulfide COS Library LibCOS Hydrogen sulfide  $H_2S$  Library LibH2S Nitrous oxide  $N_2O$  Library LibN2O Sulfur dioxide SO<sub>2</sub> Library LibSO2 Acetone  $C_3H_6O$  Library LibC3H6O Formulation of Lemmon and Span (2006)

## For more information please contact:

KCE-ThermoFluidProperties UG (limited liability) & Co. KG Professor Hans-Joachim Kretzschmar

Wallotstr. 3 01307 Dresden, Germany

Internet: www.thermofluidprop.com E-mail: info@thermofluidprop.com Phone: +49-351-27597860 Mobile: +49-172-7914607 Fax: +49-3222-4262250

### The following thermodynamic and transport properties can be calculated<sup>a</sup>:

#### **Thermodynamic Properties**

- Vapor pressure ps
- Saturation temperature T<sub>s</sub>
- Density ρ
- Specific volume v
- Enthalpy h
- Internal energy u
- Entropy s
- Exergy e
- Isobaric heat capacity c<sub>p</sub>
- Isochoric heat capacity  $c_v$
- Isentropic exponent  $\kappa$
- Speed of sound w
- Surface tension σ

#### **Transport Properties**

- Dynamic viscosity  $\eta$
- Kinematic viscosity v
- Thermal conductivity  $\lambda$
- Prandtl number Pr
- *p*, *T*(*v*,*h*) *p*, *T*(*v*,*u*)

• T, v, s(p,h)

• *T*, *v*, *h*(*p*,*s*)

• p, T, v (h,s)

**Backward Functions** 

#### Thermodynamic Derivatives

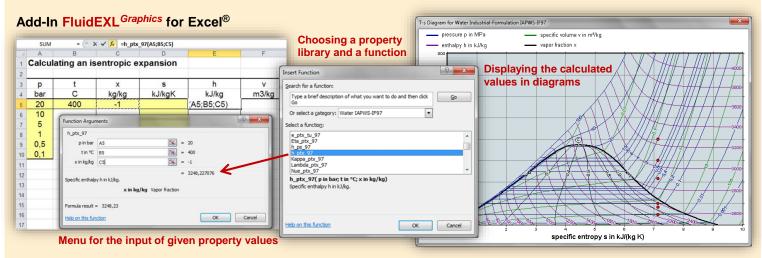
 Partial derivatives can be calculated.

<sup>a</sup> Not all of these property functions are available in all property libraries.



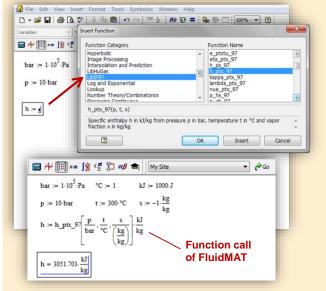


# Property Software for Calculating Heat Cycles, Boilers, Turbines and Refrigerators



#### Add-In FluidMAT for Mathcad®

The property libraries can be used in Mathcad<sup>®</sup>.



#### Add-In FluidLAB for MATLAB®

Using the Add-In FluidLAB the property functions can be called in MATLAB<sup>®</sup>.

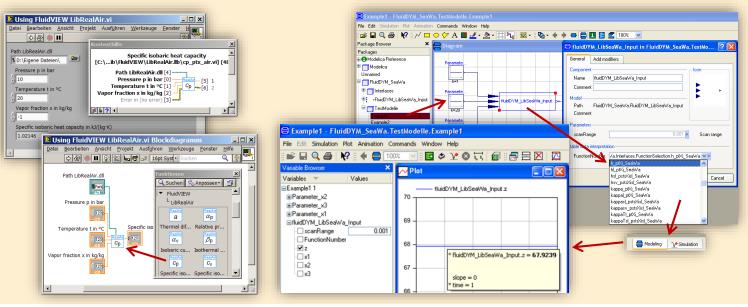
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Al Files	Image: Command Window       Image: Command Window         Image: Command Window       Image: Command Window

#### Add-On FluidVIEW for LabVIEW™

The property functions can be calculated in LabVIEW™.

#### Add-In FluidDYM for DYMOLA® (Modelica) and SimulationX®

The property functions can be called in DYMOLA® and SimulationX®.



#### Add-In FluidEES for Engineering Equation Solver<sup>®</sup>

#### ? × Function Inform ○ EES library routines Math functions Fluid properties External routines ○ Boiling and Condensation 💌 Solid/liquid properties CIEBNOIS CONTRACTOR CO CURVEFIT1D n\Fuer\_EES\HuAirProp\_SI\Beisp Tables Plots Windows Help Exa Equations Window ulating the Enthalpy - h\_ptWHuAirPn p=11 Main t=20 Unit Settings: [kJ]/[C]/[kPa]/[kg]/[degrees] W=( h = 45.4866 [kJ/kg] p = 101.3 [kPa] t = 20 [C] W = 0.01 [kg/kg] CAL No unit problems were detected. Calculation time = .1 sec.

#### App International Steam Tables for iPhone, iPad, iPod touch, Android Smartphones and Tablets

International Steam Tables

IAPWS-IF97

p,x t,x p,h p,s

Spe Den

# Online Property Calculator at www.thermofluidprop.com

Zittau's	Fluid Property	Calculator		
Fluid:	Water and Steam IAPWS-IF	97 - LiblF97 💌	12XXXV	
Function:	Specific enthalpy h(p.t.x)	• //		
Unit System:	SI 💌			
Enter given	values: Range of validity			
Pressure p		100	bar	-
Temperature	et	400	-C	·
Vapor fractio	on x le vapor fraction x	-1	kg/kg	
	Contraction in the second s	e / Recalculate		K
Result:	1981 141	THARTS	111411C	XX
Specific ent	halpy h	= 3097.38	kJ/kg	
Engineering E here.	ormation on property libraries quation Solver®, DYMOLA® iculating steam properties on description	(Modelica). Simulation	nX®, and LabView®	click
Faculty of Me Department of	University of Applied Sciences chanical Engineering of Technical Thermodynamics pachim Kretzschmar cker	Tel. +49-3583-61-184 Fax: +49-3583-61-184 E-mail: info@thermoo www.thermodynamics	46 dynamics-zittau.de	A A

### Property Software for Pocket Calculators



#### For more information please contact:

KCE-ThermoFluidProperties UG (limited liability) & Co. KG Professor Hans-Joachim Kretzschmar

Wallotstr. 3 01307 Dresden, Germany Internet: www.thermofluidprop.com E-mail: info@thermofluidprop.com Phone: +49-351-27597860 Mobile: +49-172-7914607 Fax: +49-3222-4262250

# The following thermodynamic and transport properties<sup>a</sup> can be calculated in Excel<sup>®</sup>, MATLAB<sup>®</sup>, Mathcad<sup>®</sup>, Engineering Equation Solver<sup>®</sup> (EES), DYMOLA<sup>®</sup> (Modelica), SimulationX<sup>®</sup> and LabVIEW<sup>™</sup>:

#### **Thermodynamic Properties**

- Vapor pressure p<sub>s</sub>
- Saturation temperature T<sub>s</sub>
- Density  $\rho$
- Specific volume v
- Enthalpy h
- Internal energy u
- Entropy s
- Exergy e
- Isobaric heat capacity  $c_p$
- Isochoric heat capacity  $c_v$
- Isentropic exponent κ
- Speed of sound w
- Surface tension σ

#### Transport Properties

- Dynamic viscosity  $\eta$
- Kinematic viscosity v
- Thermal conductivity  $\lambda$
- Prandtl number Pr

#### **Backward Functions**

- *T*, *v*, *s*(*p*,*h*)
- T, v, h (p,s)
- p, T, v(h,s)
- p, T (v,h)
- p, T (v,u)

#### Thermodynamic Derivatives

 Partial derivatives can be calculated.

<sup>a</sup> Not all of these property functions are available in all property libraries.

# 5. References

 Colonna, P.; Nannan, N. R.; Guardone, A.;Lemmon, E. W. Multiparameter equations of state for selected siloxanes Fluid Phase Equilibria, 244, (2006) S. 193-211

# 6. Satisfied Customers

### Date: 05/2018

The following companies and institutions use the property libraries

- FluidEXL<sup>Graphics</sup> for Excel<sup>®</sup>
- FluidLAB for MATLAB®
- FluidMAT for Mathcad®
- FluidEES for Engineering Equation Solver<sup>®</sup> EES
- FluidDYM for Dymola  $^{\ensuremath{\mathbb{R}}}$  (Modelica) and Simulation  $X^{\ensuremath{\mathbb{R}}}$
- FluidVIEW for LabVIEW<sup>™</sup>.

### 2018

Universität Madrid, Madrid, Spanien	05/2018
HS Zittau/ Görlitz, Fakultät Wirtschaft, Zittau	05/2018
HS Niederrhein, Krefeld	05/2018
GRS, Köln	03/2018
RONAL AG, Härklingen, Schweiz	02/2018
Ingenieurbüro Leipert, Riegelsberg	02/2018
AIXPROCESS, Aachen	02/2018
KRONES, Neutraubling	02/2018
Doosan Lentjes, Ratingen	01/2018

Compact Kältetechnik, Dresden	12/2017
Endress + Hauser Messtechnik GmbH +Co. KG, Hannover	12/2017
TH Mittelhessen, Gießen	11/2017
Haarslev Industries, Søndersø, Denmark	11/2017
Hochschule Zittau/Görlitz, Fachgebiet Energiesystemtechnik	11/2017
ATESTEO, Alsdorf	10/2017
Wijbenga, PC Geldermalsen, Netherlands	10/2017
Fels-Werke GmbH, Elbingerode	10/2017
KIT Karlsruhe, Institute für Neutronenphysik und Reaktortechnik	09/2017
Air-Consult, Jena	09/2017
Papierfabrik Koehler, Oberkirch	09/2017
ZWILAG, Würenlingen, Switzerland	09/2017
TLK-Thermo Universität Braunschweig, Braunschweig	08/2017
Fichtner IT Consulting AG, Stuttgart	07/2017
Hochschule Ansbach, Ansbach	06/2017
RONAL, Härkingen, Switzerland	06/2017
BORSIG Service, Berlin	06/2017

BOGE Kompressoren, Bielefeld	06/2017
STEAG Energy Services, Zwingenberg	06/2017
CES clean energy solutions, Wien, Austria	04/2017
Princeton University, Princeton, USA	04/2017
B2P Bio-to-Power, Wadersloh	04/2017
TU Dresden, Institute for Energy Engineering, Dresden	04/2017
SAINT-GOBAIN, Vaujours, France	03/2017
TU Bergakademie Freiberg, Chair of Thermodynamics, Freiberg	03/2017
SCHMIDT + PARTNER, Therwil, Switzerland	03/2017
KAESER Kompressoren, Gera	03/2017
F&R, Praha, Czech Republic	03/2017
ULT Umwelt-Lufttechnik, Löbau	02/2017
JS Energie & Beratung, Erding	02/2017
Kelvion Brazed PHE, Nobitz-Wilchwitz	02/2017
MTU Aero Engines, München	02/2017
Hochschule Zittau/Görlitz, IPM	01/2017
CombTec ProCE, Zittau	01/2017
SHELL Deutschland Oil, Wesseling	01/2017
MARTEC Education Center, Frederikshaven, Denmark	01/2017
SynErgy Thermal Management, Krefeld	01/2017

BOGE Druckluftsysteme, Bielefeld	12/2016
BFT Planung, Aachen	11/2016
Midiplan, Bietigheim-Bissingen	11/2016
BBE Barnich IB	11/2016
Wenisch IB,	11/2016
INL, Idaho Falls	11/2016
TU Kältetechnik, Dresden	11/2016
Kopf SynGas, Sulz	11/2016
INTVEN, Bellevne (USA)	11/2016
DREWAG Dresden, Dresden	10/2016
AGO AG Energie+Anlagen, Kulmbach	10/2016
Universität Stuttgart, ITW, Stuttgart	09/2016
Pöyry Deutschland GmbH, Dresden	09/2016
Siemens AG, Erlangen	09/2016
BASF über Fichtner IT Consulting AG	09/2016
B+B Engineering GmbH, Magdeburg	09/2016
Wilhelm Büchner Hochschule, Pfungstadt	08/2016

	Webasto Thermo & Comfort SE, Gliching	08/2016
	TU Dresden, Dresden	08/2016
	Endress+Hauser Messtechnik GmbH+Co. KG, Hannover	08/2016
	D + B Kältetechnik, Althausen	07/2016
	Fichtner IT Consulting AG, Stuttgart	07/2016
	AB Electrolux, Krakow, Poland	07/2016
	ENEXIO Germany GmbH, Herne	07/2016
	VPC GmbH, Vetschau/Spreewald	07/2016
	INWAT, Lodz, Poland	07/2016
	E.ON SE, Düsseldorf	07/2016
	Planungsbüro Waidhas GmbH, Chemnitz	07/2016
	EEB Enerko, Aldershoven	07/2016
	IHEBA Naturenergie GmbH & Co. KG, Pfaffenhofen	07/2016
	SSP Kälteplaner AG, Wolfertschwenden	07/2016
	EEB ENERKO Energiewirtschaftliche Beratung GmbH, Berlin	07/2016
	BOGE Kompressoren Otto BOGE GmbH & Co KG, Bielefeld	06/2016
	Universidad Carlos III de Madrid, Madrid, Spain	04/2016
	INWAT, Lodzi, Poland	04/2016
	Planungsbüro WAIDHAS GmbH, Chemnitz	04/2016
	STEAG Energy Services GmbH, Laszlo Küppers, Zwingenberg	03/2016
	WULFF & UMAG Energy Solutions GmbH, Husum	03/2016
	FH Bielefeld, Bielefeld	03/2016
	EWT Eckert Wassertechnik GmbH, Celle	03/2016
	ILK Institut für Luft- und Kältetechnik GmbH, Dresden 02/2016, 06/2	2016 (2x)
	IEV KEMA - DNV GV – Energie, Dresden	02/2016
	Allborg University, Department of Energie, Aalborg, Denmark	02/2016
	G.A.M. Heat GmbH, Gräfenhainichen	02/2016
	Institut für Luft- und Kältetechnik, Dresden 02/2016, 05/2016	06/2016
	Bosch, Stuttgart	02/2016
	INL Idaho National Laboratory, Idaho, USA 11/2016	01/2016
	FriedI ID, Wien, Austria	01/2016
	Technical University of Dresden, Dresden	01/2016
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	EES Enerko, Aachen	12/2015
	Ruldolf IB, Strau, Austria	12/2015
	Allborg University, Department of Energie, Aalborg, Denmark	12/2015

Bosch, Lohmar	10/2015
Team Turbo Machines, Rouen, France	09/2015
BTC – Business Technology Consulting AG, Oldenburg	07/2015
KIT Karlsruhe Institute of Technology, Eggenstein-Leopoldshafen	07/2015
ILK, Dresden	07/2015
Schniewindt GmbH & Co. KG, Neuenwalde	08/2015
2014	
PROJEKTPLAN, Dohna	04/2014
Technical University of Vienna, Austria	04/2014
MTU Aero Engines AG, Munich	04/2014
GKS, Schweinfurt	03/2014
Technical University of Nuremberg	03/2014
EP-E, Niederstetten	03/2014
Rückert NatUrgas GmbH, Lauf	03/2014
YESS-World, South Korea	03/2014
ZAB, Dessau	02/2014
KIT-TVT, Karlsruhe	02/2014
Stadtwerke Neuburg	02/2014
COMPAREX, Leipzig for RWE Essen	02/2014
Technical University of Prague, Czech Republic	02/2014
HS Augsburg	02/2014
Envi-con, Nuremberg	01/2014
DLR, Stuttgart	01/2014
Doosan Lentjes, Ratingen	01/2014
Technical University of Berlin	01/2014
Technical University of Munich	01/2014
Technical University of Braunschweig	01/2014
M&M Turbinentechnik, Bielefeld	01/2014
2013	
TRANTER-GmbH, Artern	12/2013
SATAKE, Shanghai, China	12/2013
VOITH, Kunshan, China	12/2013
ULT, Löbau	12/2013
MAN, Copenhagen, Dänemark	11/2013
DREWAG, Dresden	11/2013
Haarslev Industries, Herlev, Dänemark	11/2013
STEAG, Herne	11/2013, 12/2013
Ingersoll-Rand, Oberhausen	11/2013
Wilhelm-Büchner HS, Darmstadt	10/2013

IAN/ Characita	10/0012
IAV, Chemnitz	10/2013
Technical University of Regensburg	10/2013
PD-Energy, Bitterfeld	09/2013
Thermofin, Heinsdorfergrund	09/2013
SHI, New Jersey, USA	09/2013
M&M Turbinentechnik, Bielefeld	08/2013
BEG-BHV, Bremerhaven	08/2013
TIG-Group, Husum	08/2013
COMPAREX, Leipzig for RWE Essen	08/2013, 11/2013 12/2013
	08/2013
University of Budapest, Hungary	
Siemens, Frankenthal	08/2013, 10/2013
	11/2013
VGB, Essen	07/2013, 11/2013
Brunner Energieberatung, Zurich, Switzerland	07/2013
Technical University of Deggendorf	07/2013
University of Maryland, USA	07/2013, 08/2013
University of Princeton, USA	07/2013
NIST, Boulder, USA	06/2013
IGUS GmbH, Dresden	06/2013
BHR Bilfinger, Essen	06/2013
SÜDSALZ, Bad Friedrichshall	06/2013, 12/2013
Technician School of Berlin	05/2013
KIER, Gajeong-ro, Südkorea	05/2013
Schwing/Stetter GmbH, Memmingen	05/2013
Vattenfall, Berlin	05/2013
AUTARK, Kleinmachnow	05/2013
STEAG, Zwingenberg	05/2013
Hochtief, Düsseldorf	05/2013
University of Stuttgart	04/2013
Technical University -Bundeswehr, Munich	04/2013
Rerum Cognitio Forschungszentrum, Frankfurt	04/2013
Kältetechnik Dresen + Bremen, Alfhausen	04/2013
University Auckland, New Zealand	04/2013
MASDAR Institut, Abu Dhabi, United Arab Emirates	03/2013
Simpelkamp, Dresden	02/2013
VEO, Eisenhüttenstadt	02/2013
ENTEC, Auerbach	02/2013
Caterpillar, Kiel	02/2013
Technical University of Wismar	02/2013
Technical University of Dusseldorf	02/2013
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ILK, Dresden Fichtner IT, Stuttgart Schnepf Ingeniuerbüro, Nagold Schütz Engineering, Wadgassen Endress & Hauser, Reinach, Switzerland Oschatz GmbH, Essen frischli Milchwerke, Rehburg-Loccum	01/2013, 08/2013 01/2013, 11/2013 01/2013 01/2013 01/2013 01/2013 01/2013 01/2013
2012	
Voith, Bayreuth	12/2012
Technical University of Munich	12/2012
Dillinger Huette	12/2012
University of Stuttgart	11/2012
Siemens, Muehlheim	11/2012
Sennheiser, Hannover	11/2012
Oschatz GmbH, Essen	10/2012
Fichtner IT, Stuttgart	10/2012, 11/2012
Helbling Technik AG, Zurich, Switzerland	10/2012
University of Duisburg	10/2012
Rerum Cognitio Forschungszentrum, Frankfurt	09/2012
Pöyry Deutschland GmbH, Dresden	08/2012
Extracciones, Guatemala	08/2012
RWE, Essen	08/2012
Weghaus Consulting Engineers, Wuerzburg	08/2012
GKS, Schweinfurt	07/2012
COMPAREX, Leipzig for RWE Essen	07/2012
GEA, Nobitz	07/2012
Meyer Werft, Papenburg	07/2012
STEAG, Herne	07/2012
GRS, Cologne	06/2012
Fichtner IT Consult, Chennai, India	06/2012
Siemens, Freiburg	06/2012
Nikon Research of America, Belmont, USA	06/2012
Niederrhein University of Applied Sciences, Krefeld	06/2012
STEAG, Zwingenberg	06/2012
Mainova, Frankfurt on Main via Fichtner IT Consult	05/2012
Endress & Hauser	05/2012
PEU, Espenheim	05/2012
Luzern University of Applied Sciences, Switzerland	05/2012

BASF, Ludwigshafen (general license) via Fichtner IT Consult	05/2012
SPX Balcke-Dürr, Ratingen	05/2012, 07/2012
Gruber-Schmidt, Wien, Austria	04/2012
Vattenfall, Berlin	04/2012
ALSTOM, Baden	04/2012
SKW, Piesteritz	04/2012
TERA Ingegneria, Trento, Italy	04/2012
Siemens, Erlangen	04/2012, 05/2012
LAWI Power, Dresden	04/2012
Stadtwerke Leipzig	04/2012
SEITZ, Wetzikon, Switzerland	03/2012, 07/2012
M & M, Bielefeld	03/2012
Sennheiser, Wedemark	03/2012
SPG, Montreuil Cedex, France	02/2012
German Destilation, Sprendlingen	02/2012
Lopez, Munguia, Spain	02/2012
Endress & Hauser, Hannover	02/2012
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WIPAK, Walsrode	02/2012
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