

PROSIMPLUS APPLICATION EXAMPLE SIMULATION OF A CROSS-FLOW PLATE-FIN HEAT EXCHANGER (PFHE) WITH PROSEC CAPE-OPEN UNIT OPERATION

INTEREST OF THIS EXAMPLE

This example shows a cross-flow plate-fin heat exchanger used in a gas-gas application. This heat exchanger is modeled using ProSec, ProSim's CAPE-OPEN compliant unit operation dedicated to the simulation of plate-fin heat exchangers. Regarding the thermodynamic and physico-chemical data needed, they are automatically calculated using the thermodynamic calculation server of the process simulation software.

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CORRESPONDING PROSIMPLUS FILE		COPROSEC_EX_EN-Cross-flow-heat-exchanger.pmp3			

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Version: September 2018 Page: 2 / 11

TABLE OF CONTENTS

1.	PR	OCESS MODELING	3
	1.1.	Process description	3
	1.2.	Process flowsheet	4
	1.3.	Compounds	5
	1.4.	Thermodynamic model	5
	1.5.	Operating parameters	5
	1.5	.1. Process feed	5
	1.5	.2. Plate-fin heat exchanger E101	6
2.	RE	SULTS	9
	2.1.	Mass and energy balances	9
	2.2.	Global results	9
	2.3.	Plate-fin heat exchanger profiles	10
3.	RE	FERENCES	11

Version: September 2018 Page: 3 / 11

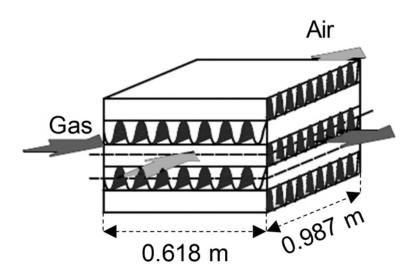
1. Process modeling

1.1. Process description

This example deals with a gas-gas cross-flow plate-fin heat exchanger (PFHE). Compact heat exchangers are characterized by a large heat transfer surface area per unit volume. Amongst different varieties of compact heat exchangers, crossflow plate-fin heat exchangers are widely used in aerospace, automobile, cryogenic and chemical process plants for their low weight and volume, high efficiency. It will be simulated with ProSec CAPE-OPEN unit operation. In ProSec, cross-flow heat exchangers are modeled as counter-flow heat exchangers and an efficiency coefficient [ROE10] is computed and applied to take into account the cross-flow behavior. The assumptions of the model are:

- ✓ Common wall temperature,
- ✓ Longitudinal conduction not taken into account,
- ✓ Two fluids,
- ✓ Continuous thermodynamic option not available.

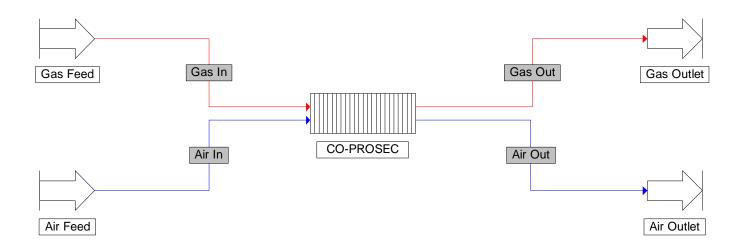
The example is extracted from [MIS09]. The sketch is shown below. The two fluids are modeled as air. The "Gas" is the main stream and the "Air" is the cross-flow stream.



The thermodynamic and physico-chemical data needed for the simulation are automatically calculated by the unit operation. In this example, as the ProSec CAPE-OPEN unit operation is used in ProSimPlus simulation environment, it uses Simulis Thermodynamics, the calculation server for thermophysical properties and phase equilibria calculations available in ProSimPlus.

Version: September 2018 Page: 4 / 11

1.2. Process flowsheet



Version: September 2018 Page: 5 / 11

1.3. Compounds

The compounds taken into account in the simulation, their chemical formula and CAS numbers are shown in the following table. Their pure component properties are extracted from the standard data base provided with ProSimPlus [ROW17].

Compound	Chemical formula	CAS number	
Oxygen	O ₂	7782-44-7	
Nitrogen	N_2	7727-37-9	

1.4. Thermodynamic model

The PSRK profile, which is based on a predictive combined approach model [HOL91], [GME91], [CHE02] is selected.

1.5. Operating parameters

1.5.1. Process feed

	Gas	Air	
Temperature (°C)	240 4		
Pressure (bar)	1	1	
Total flow rate (kg/s)	kg/s) 0.8962 0.8296		
Molar fraction			
Oxygen	0.21		
Nitrogen	0.79		

Version: September 2018 Page: 6 / 11

1.5.2. Plate-fin heat exchanger

✓ General parameters

Parameters	Value	
Type of exchanger	CO-ProSec	
Number of body	1	
Orientation	Horizontal	
Fin data base	2015 -> Now	
Material		
Туре	Other	
a1	160	
a2 = a3 = a4 = a5 = a6 =a7	000	
Used width (m)	0.987	
Thickness of the side bars (mm)	0.001	
Thickness of the end bars (mm)	0.001	
Thickness of the separation plates (mm)	0.500	

√ Streams parameters

	Stre	eam	
Parameter	Gas	Air	
Flow direction	From top to bottom	From bottom to top	
Cross flow	No	Yes	
Heat exchange correlation	HTFS85		

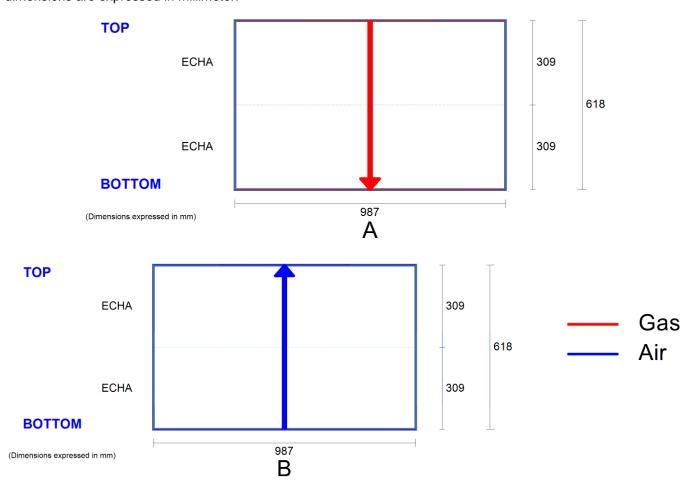
✓ Fin characteristics

Name	Fin #1
Origin	User
Calculation mode	From geometry
Reference	1001
Туре	Serrated
Height (mm)	9.799
Thickness (mm)	0.101
Fins number per meter	654.5
Serration length (mm)	7.551

Version: September 2018 Page: 7 / 11

✓ Reference passages

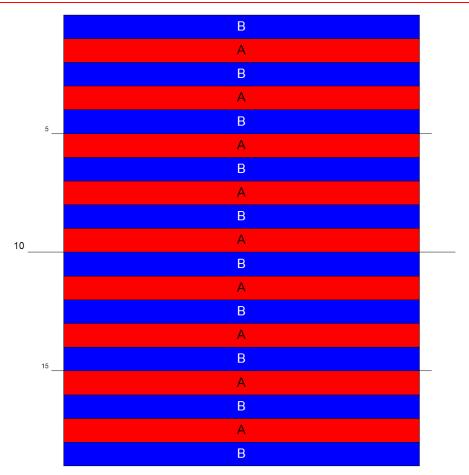
Fin #1 is used for each stream. The following figures show the two reference passages of the heat exchanger. The dimensions are expressed in millimeter.



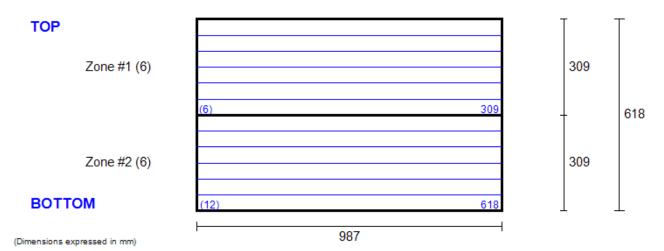
✓ Stacking

Parameters	Value
Sequence 1	
Number of repetitions of the sequence	9
Sequence	ВА
Sequence 2	
Number of repetitions of the sequence	1
Sequence	В

Version: September 2018 Page: 8 / 11



✓ Number of meshs for each elementary zone (dimensions are expressed in millimeter)



Version: September 2018 Page: 9 / 11

2. RESULTS

2.1. Mass and energy balances

Streams	Air In	Air Out	Gas In	Gas Out	
Total flow	kg/s	0.8296	0.8296	0.8962	0.8962
Mole fractions					
OXYGEN		0.21	0.21	0.21	0.21
NITROGEN	0.79	0.79	0.79	0.79	
Physical state		Vapor	Vapor	Vapor	Vapor
Temperature	°C	4	201.81	240	57.828
Pressure	kPa	100	95.282	100	97.997
Enthalpic flow kW		-17.831	149.49	196.94	29.624
Vapor molar fraction		1	1	1	1

2.2. Global results

The heat duty computed by ProSec is 167 kW. [MIS09] indicates a value of 160 kW. The deviation is explained by the differences in the assumptions:

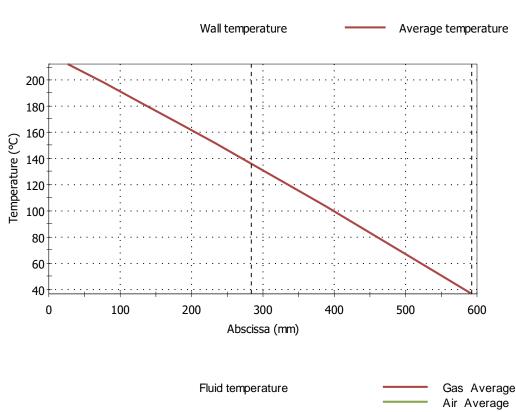
- ✓ Constant physico-chemical properties for [MIS09] vs temperature dependent for ProSec,
- ✓ The correlations used to compute the Colburn and Fanning factors are not the same.

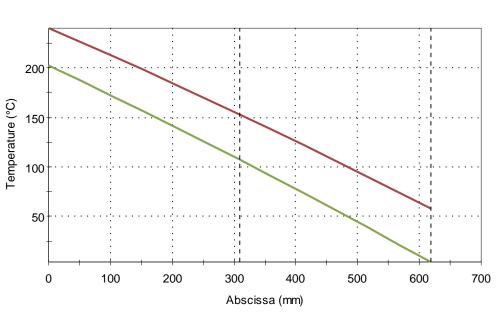
The pressure drops computed by ProSec (2 003 Pa for the gas stream and 4 718 Pa for the air stream) are closed to the ones calculated by [MIS09].

Version: September 2018 Page: 10 / 11

2.3. Plate-fin heat exchanger profiles

Several profiles (wall temperature, fluid temperature, pressure, heat transfer coefficient, vaporization ratio...) in the heat exchanger are available after the simulation from ProSec edition window ("Results" tab). The following figures show the average wall temperature profile and the fluid mean temperature profiles along the length of the heat exchanger.





Version: September 2018 Page: 11 / 11

3. REFERENCES

[CHE02]	CHEN J., FISCHER K., GMEHLING J., "Modification of PSRK Mixing Rules and Results for Vapor – Liquid Equilibria, Enthalpy of Mixing and Activity Coefficients at Infinite Dilution", Fluid Phase Equilib., 200, 411-429 (2002)
[GME95]	GMEHLING J., "From UNIFAC to Modified UNIFAC to PSRK with the Help of DDB", Fluid Phase Equilib., 107, 1-29 (1995)
[HOL91]	HOLDERBAUM T., GMEHLING J., "PSRK: A Group Contribution Equation of State based on UNIFAC", Fluid Phase Equilib., 70, 251-265 (1991)
[MIS09]	MISHRA M., DAS P.K., "Thermoeconomic Design-Optimisation of Crossflow Plate-Fin Heat Exchanger Using Genetic Algorithm", Int. J. Exergy, 6, 837-852 (2009)
[ROE10]	ROETZEL W., SPANG B., "VDI Heat Atlas, Chapter C1: Thermal Design of Heat Exchangers", Springer-Verlag, 2nd Edition (2010)
[ROW17]	ROWLEY R.L., WILDING W.V., OSCARSON J.L., GILES N.F., "DIPPR® Data Compilation of Pure Chemical Properties", Design Institute for Physical Properties, AIChE, New York, NY (2017)