# Benchmarking a Cryogenic Code to the FREIA Helium Liquefier

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The FREIA He Liquefier





#### Background and project motivation

- Helium liquefier from Linde group
- Liquid He crucial for research infrastructure
- Simulations of thermodynamics better understanding of liquefaction process
- Main objective: build a theoretical model of the FREIA liquefier based upon the principle of enthalpy conservation



Linde liquid He production and recovery system – lindeengineering.com





# Thermodynamics Background

- Intensive quantities: T,P
- Extensive quantities: U, H, S, V
- Enthalpy

H = U + PV

• Equation of state for ideal gas

PV = nRT

- Ideal gas vs real gas
- More complex relations for real gases use CoolProp library





## **Thermodynamics Background**

- Heat capacity  $C=rac{\partial Q}{\partial T}$  Ideal monatomic gas:  $c_V=rac{3}{2}R$  ,  $c_P=rac{5}{2}R$
- Entropy real and ideal processes

$$S \geq rac{Q}{T}$$

Carnot heat cycle









# Thermodynamics Background

- Enthalpy and phases: latent heat
- Critical point
- Saturation dome
- Vapour quality
  - Saturated vapour: 100%
  - Saturated liquid: 0%







# Cycle components: heat exchanger

- Parallel flows or counter-flow
- Enthalpy in heat exchanger



$$egin{array}{ll} H_2 &= H_1 - dH \ H_4 &= H_3 + dH \end{array}$$



# Cycle components: two-fluid counterflow heat exchanger

• Temperatures - system of 1st-order linear differental equations

$$rac{d}{dx}egin{bmatrix} T_a \ T_b \end{bmatrix} = egin{bmatrix} -rac{C'}{C_a} & rac{C'}{C_a} \ -rac{C'}{C_b} & rac{C'}{C_b} \end{bmatrix} egin{bmatrix} T_a \ T_b \end{bmatrix}$$

Linear solution

$$egin{bmatrix} T_a \ T_b \end{bmatrix} = A_1ec v_1 e^{lpha_1 x} + A_2ec v_2 e^{lpha_2 x}$$

where  $A_1, A_2$  come from boundary conditions,  $\alpha_i$  is the eigenvalue and  $\vec{v}_i$  is the eigenvector







#### Cycle components: two-fluid counterflow heat exchanger

• Solve for total enthalpy transfer  $\Delta \dot{H}$ 

 $\Delta \dot{H} = C_H (T_3 - T_1)$ 

• If 
$$(C_a \neq C_b)$$
:  $C_H = \frac{C_a C_b (1 - e^{-\alpha})}{C_a e^{-\alpha} - C_b}$   
• If  $(C_a = C_b)$ :  $\frac{1}{C_H} = \frac{1}{C_a} + \frac{1}{C'L}$ 

where 
$$lpha = C'ig(rac{1}{C_a} - rac{1}{C_b}ig)$$

Solved for analytically

$\mathop{\downarrow} T_1$	$T_4$
x = 0	
$dx\downarrow$	
$C_a = C''$	$lackslash C_b$
x = L	
$\dagger T_2$	$T_3$





#### Cycle components: three-fluid counterflow heat exchanger

Same principle with heat equations







#### Cycle components: three-fluid counterflow heat exchanger

• Solving for total enthalpy transfer  $\Delta \dot{H}_1, \Delta \dot{H}_2$ 

 $v_{3i}$ 

where  $\alpha_i$  is the eigenvalue and  $\dot{v}_i =$  is the corresponding eigenvector

- Long analytic expressions with respect to two-fluid case, but still easy numerically
- In FREIA liquefier: two versions, one with cold flow x = L 2 reversed



 $T_{c2,f}$ 

 $C_{2}$ 

 $C_h$ 

 $T_{h,t}$ 

 $C'_1$ 

x = 0

dx .

 $C_1$ 





#### Cycle components: turboexpander

- Work against centrifugal force
- Ideal gas at sonic speeds and isentropic expansion

$$rac{T_1}{T_3} = 2$$
  $rac{P_1}{P_3} = 5.64$ 

• Gas must not liquefy







#### Cycle component: Joule-Thomson valve

- Joule-Thomson (JT) Effect
  - Temperature change of real gas
  - Irreversible, isenthalpic process:  $\Delta H=0$
  - JT coefficient

$$\mu_{JT} = \left(\frac{\partial T}{\partial P}\right)_{H=\text{const}}$$

Inversion temperature



JT Coefficient at 1 bar. Image courtesy: Hankwang, Wikimedia Commons





#### Cycle components - MATLAB implementation

• JT valve

%JT-valve - isenthalpic process h5 = py.CoolProp.CoolProp.PropsSI('H', 'P', P5, 'T', T5, 'Helium'); h6 = h5; T6 = py.CoolProp.CoolProp.PropsSI('T', 'P', P6, 'H', h6, 'Helium');

#### Turboexpander



%Turboexpander for helium h1 = py.CoolProp.CoolProp.PropsSI('H','P',P1,'T',T1,'Helium'); T3 = T1/2; P3 = P1/5.64; h3 = py.CoolProp.CoolProp.PropsSI('H','P',P3,'T',T3,'Helium'); deltah = h1 - h3; We = m\*deltah;





#### Liquefaction cycles – Linde-Hampson

- Most fundamental cycle, 1895
- 1<sup>st</sup> law of thermodynamics in dashed control volume

$$\dot{m}h_1=\dot{m}_fh_f+(\dot{m}-\dot{m}_f)h_5$$
  
Global yield:  $y\equiv rac{\dot{m}_f}{\dot{m}}=rac{h_1-h_5}{h_f-h_5}$ 

Around phase separator:  $\dot{m}h_2 = \dot{m}_f h_f + (\dot{m} - \dot{m}_f)h_4$ Local yield:  $y_l \equiv \frac{\dot{m}_f}{\dot{m}_3} = \frac{h_2 - h_4}{h_f - h_4}$ 







#### Liquefaction - MATLAB implementation

• Phase separator



if T6 < 5.1953 %critical point of He % enters liquid phase hliq= py.CoolProp.CoolProp.PropsSI('H','T',T6,'Q',0,'Helium'); %specific heat capacity hgas= py.CoolProp.CoolProp.PropsSI('H','T',T6,'Q',1,'Helium'); y1=(hgas-H5/Q5)/(hgas-hliq); y1=min(1,max(0,y1)); % local without HE Hgas = H6 - y1\*Q6\*hliq; | H7= Hgas;

• No matter the cycle complexity, local yield is always the same!







#### Liquefaction cycles - Claude

- Improvement, 1902
- Two more heat exchangers, isentropic expansion in turboexpander
- 1<sup>st</sup> law of thermodynamics

$$\dot{m}h_1=\dot{W}_e+(\dot{m}-\dot{m}_f)h_{11}+\dot{m}_fh_f$$

$$\dot{W}_e=\dot{m}_eh_2-\dot{m}_eh_e\qquad x=rac{m_e}{\dot{m}}$$

$$y\equiv rac{m_f}{\dot{m}}=rac{h_{11}-h_1}{h_{11}-h_f}+xrac{h_2-h_e}{h_{11}-h_f}$$





0.01

Cycle Point



#### Claude cycle simulations

Temperature and Gas Flow Profile: Claude









# Liquefaction cycles - Collins

- Collins 1946
- Similar to Claude, but two more heat exchangers and one more turboexpander
- Analogously: 1<sup>st</sup> law around control volume gives

$$y\equiv rac{\dot{m}_f}{\dot{m}}=rac{h_{17}-h_1}{h_{17}-h_f}+x_1rac{\Delta h_{e1}}{h_{17}-h_f}+x_2rac{\Delta h_{e2}}{h_{17}-h_f}$$

with

$$\dot{W}_{e_i}=\dot{m}_{e_i}\Delta h_{e_i}$$
  $x_i=rac{\dot{m}_{e_i}}{\dot{m}}$  ,  $i=1,2$ 







#### **Collins cycle simulations**











The FREIA Liquefier Schematic with Coldbox, with simplified drawing to the right



# The FREIA liquefier

- Similar to Collins
- Liquid nitrogen pre-cooling system: thermal reservoir
- Single mass flow through both TXP
- 1<sup>st</sup> law around control volume

$$y \equiv rac{\dot{m}_f}{\dot{m}} = rac{h_{16} - h_1}{h_{16} - h_f} + rac{\dot{m}_e}{\dot{m}} rac{\Delta h_1 + \Delta h_2}{h_{16} - h_f} \ - rac{\dot{m}_{pur}}{\dot{m}} rac{h_6 - h_{16}}{h_{16} - h_f} + rac{\dot{m}_N}{\dot{m}} rac{\Delta h_N}{h_{16} - h_f}$$

• Extra terms: purification and LN2 precooling

























Maximum yield of model and sanity check









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#### FREIA sensor values

• Some gas to cool purification system – periodic disruption in *P* and *T* 







#### Discussion and outlook

- Enthalpy conservation and physical processes
- Most cycle components linear
- Only liquefaction non-linear weak non-linearity for system as a whole
- Yield of model simulation close to yields of real liquefier
- Maximum yield of model mainly tune heat exchanger design parameter  $C^\prime$





#### Discussion and outlook

- Turboexpanders not entirely isentropic
- Specific heat capacity  $c_P$  not constant for last heat exchanger
- Importance of optimized liquefaction cycles
- More sensors for future comparisons?





# Thank you for your attention!

