

«Hydrogen production via Steam Bio-methane Reforming»

Metacon Group

Δρ. Θωμάς Χαλκίδης, Διευθύνων Σύμβουλος, METACON SA, CTO Metacon Group

Personal information and studies

Personal information:

- Chalkidis I. Thomas
- Dr. Chemical Engineer
- Managing Director of METACON SA.

Education

1997: Diploma in Chemical Engineering with a focus on the Environment (University of Patras).

2002: PhD in Chemical Engineering in Environmental Catalysis (University of Patras)

2004: Master's Degree in Environment and Energy. (Univ. of Patras)

H Metacon SA

- ✓ It was founded in 2002 as a spin-off of the University of Patras for the commercialization of innovative processes and reform technologies developed at the university. Former name HELBIO SA.
- ✓ The core function of Metacon SA in the Metacon Group focuses on catalytic overhaul and R&D of reactors as well as the construction of hydrogen systems.
- ✓ The headquarters of Metacon SA remain in Patras and the university is still a talent pool for the company with many graduates among its employees, totaling 28 people.
- ✓ Metacon is active in the field of hydrogen and power generation, developing its own hydrogen production systems either for industrial use or integrated with fuel cells for Combined Heat and Power (CHP) solutions.
- ✓ Metacon S.A. has strong partnerships with Greek universities and institutions. The company is involved in many national and European funded projects with various national and European academic and industrial partners.
- ✓ Since 2023, Metacon S.A. has installed and operates a Management System in accordance with the requirements of ISO 9001:2015, ISO 14001:2015 and ISO 45001:2018 standards.



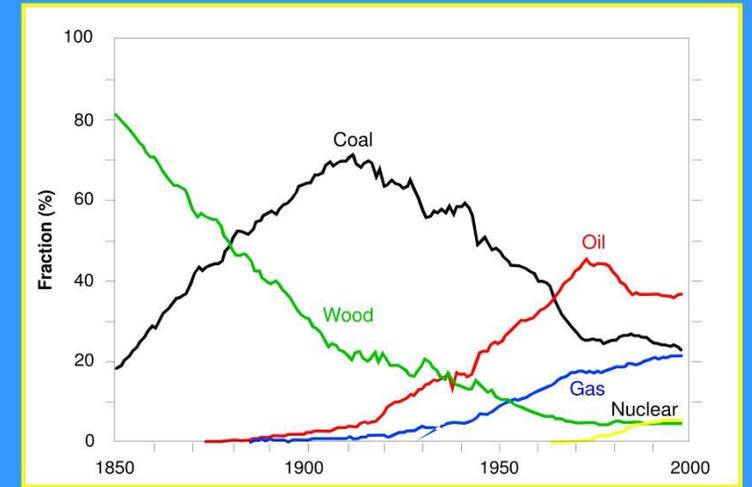
Source: Company information.
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Hydrogen: The energy carrier of the (near) future

- High energy content: 120 MJ/kg
- Zero emissions at the point of use
- Versatile applications: industry, transport, energy
- Hydrogen Economy Roadmap Gains Global Traction



World Primary Energy Substitution



86025 Energy Systems Analysis

Arnulf Grubler

| Fuel Type | Percent Hydrogen | Energy Content (Btu/lb) | Particulates (lb/Btu) | Carbon Dioxide (normal.) |
|-------------|------------------|-------------------------|-----------------------|--------------------------|
| Dry wood | 5 | 6,900 | 5.22 | 100 |
| Coal | 50 | 10,000 | 5.00 | 31 |
| Oil | 67 | 19,000 | 0.18 | 21 |
| Natural Gas | 80 | 22,500 | <0.01 | 15 |
| Hydrogen | 100 | 61,000 | 0.0 | 0 |

Source: Company information.

Hydrogen: The energy carrier of the (near) future

Current Hydrogen Production Landscape:

Total hydrogen production ~100 million tons in 2024
 ~96% from fossil fuels (natural gas, naphtha, coal)

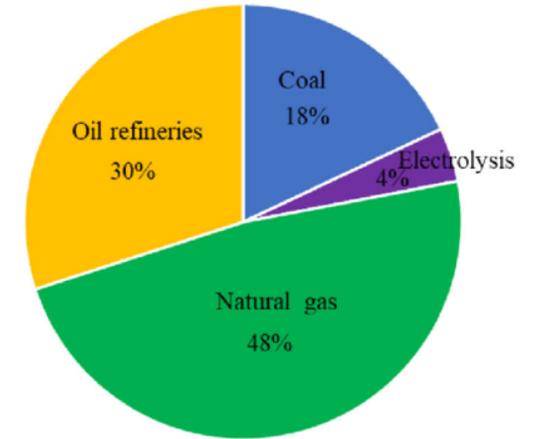
Dominant method: Catalytic Steam Methane Reforming (SMR)

Environmental impact: high CO₂ emissions

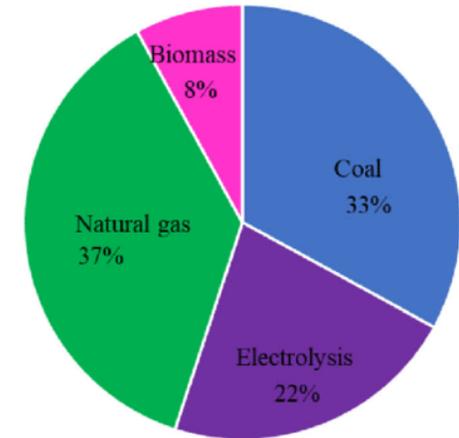
~9 kg CO₂/kg H₂ from natural gas

~20 kg CO₂/kg H₂ from carbon

Strong global push towards "blue" and "green" hydrogen



At present



Prediction Year 2050

Source: Company information.

Hydrogen production pathways

- Thermal: SMR, ATR, POX, Gasification, Methane Pyrolysis
- Electrolytic: Alkaline, PEM
- Biological : Dark or Photo fermentation
- Photolytic: Photoelectrochemical / photocatalytic water splitting (R&D)

CAPEX & OPEX technology comparison

| Technology | CAPEX | OPEX | Efficiency | Emissions | TRL | Scalability |
|-----------------------------|----------|--------------|----------------|-----------------------------|------------|-------------|
| SMR | Low | Low | High (thermal) | High | Commercial | Very High |
| SMR/ATR + CCS | Med | Low–Med | High (thermal) | Low–Med | Commercial | High |
| Gasification (coal/biomass) | Med–High | Med | High | High→Low (with CCS) | Commercial | High |
| Alkaline Electrolysis | Med | High (power) | 60–70% (LHV) | Near-zero (renewable power) | Commercial | High |
| PEM Electrolysis | High | High (power) | 60–70% (LHV) | Near-zero (renewable power) | Commercial | High |

SMR: Steam Methane Reforming
 ATR: Autothermal Reforming
 CCS: Carbon Capture & Sequestration

Source: Company information.

Case study:

Technology: Steam methane reforming for hydrogen production

Fuel origin: Biogenic methane (from upgraded Biogas)

Starting the design..

Inputs & Requirements:

- Feedstock Characteristics
- Water & Utility Requirements
- Hydrogen quantity and quality
- Catalyst Selection & Protection
- Energy Integration & Heat Management
- Carbon Management & Emissions
- Materials & Corrosion Management
- Plant Integration & Control Strategy
- Process Safety & Hazard Analysis
- Economic & Environmental Assessment

Modelling the process

***Boundary conditions**

Design Objectives:

Fuel: Biomethane (97 vol. % CH₄)*

Hydrogen output: 50 m³/h*

Purity: >99.97 vol. % (ISO 14687:2025)*

H₂ production cost <2 euros/kg*

Hydrogen production technology: Steam reforming reaction (H₂> 70 vol. %)

Molar Steam to carbon ratio: 3.5

CO minimization reaction(s): Single stage water Gas Shift reaction (CO: <1.0 vol. %)

Purification: Pressure Swing Adsorption (PSA)

H₂ Recovery rate >80%

Modelling the process-Desulfurization

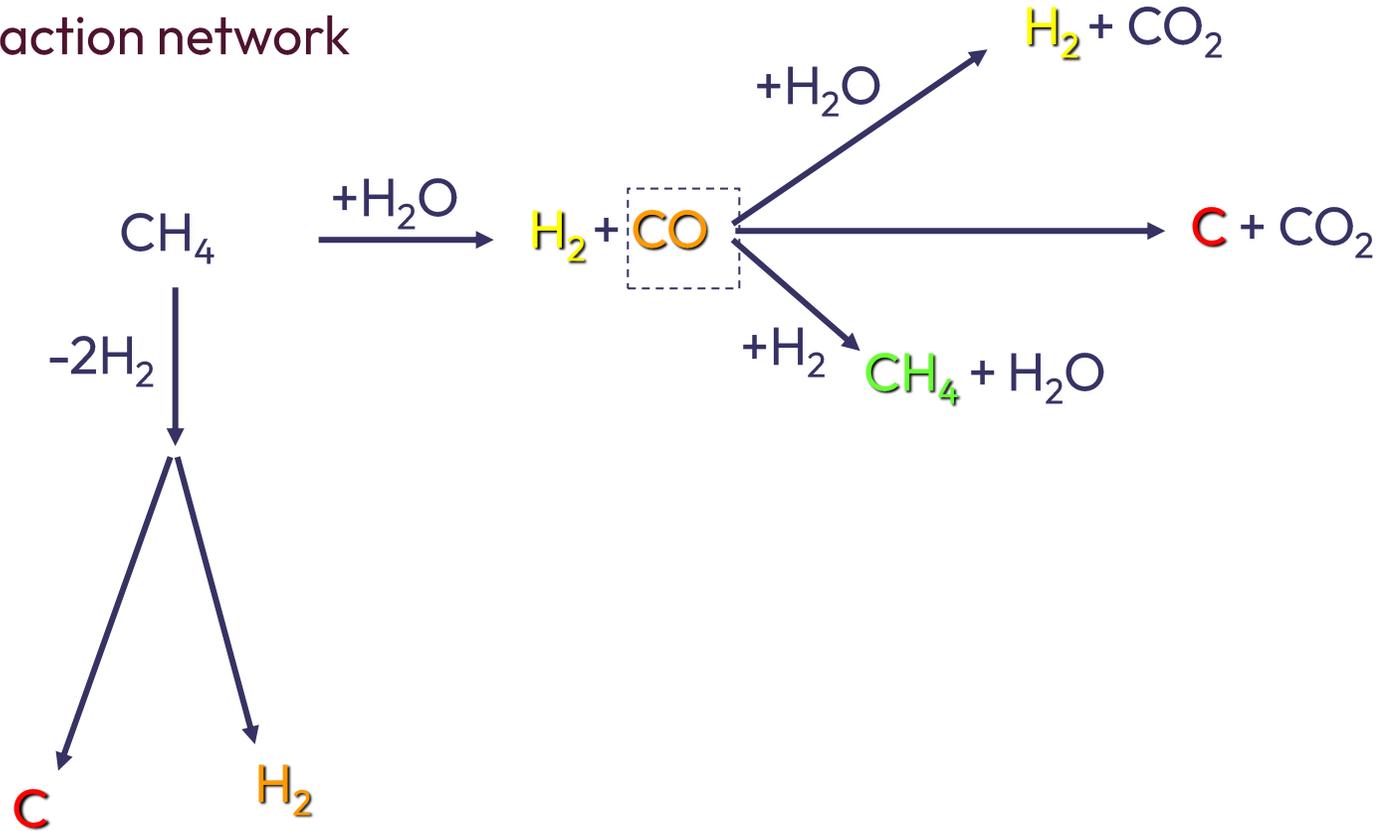
- **Fuels** may contain sulfur compounds (H_2S , mercaptans, odorants) and is essential to remove them to prevent poisoning of downstream catalysts
- Even a few ppm of sulfur can severely poison and deactivate them.

How is done...?

- a) If sulfur content is high, a hydrogenation step (hydrotreater with Co–Mo catalyst) converts organosulfur to H_2S , followed by ZnO bed to capture H_2S .
- b) For low sulfur fuels (odorant-level), a single ZnO guard bed or use of passive commercial adsorbent typically modified zeolites with additives like Cu, Zn, etc are often sufficient .
- c) You need the expected sulfur content (mg/m^3) and the uptake capacity of the adsorbent typically mg of S per g of adsorbent to calculate the amount of adsorbent needed for $X m^3$ of input fuel.

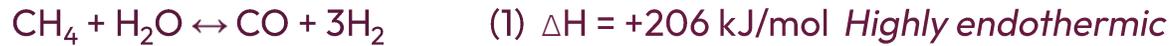
Modelling the process-Reformer

Steam/methane reaction network



Modelling the process-Reformer

Methane steam reforming reaction (700-800°C)



Side reaction(s) in reforming:



Combustion reactions (800-900°C)



Catalyst reforming: Ni/Al₂O₃ plus promoters/stabilizers

Catalyst combustion: Proprietary on Al₂O₃

Pressure: 10-11 barg (reforming)

Steam/Carbon: 3.5

Outlet temperature: 750-770 °C

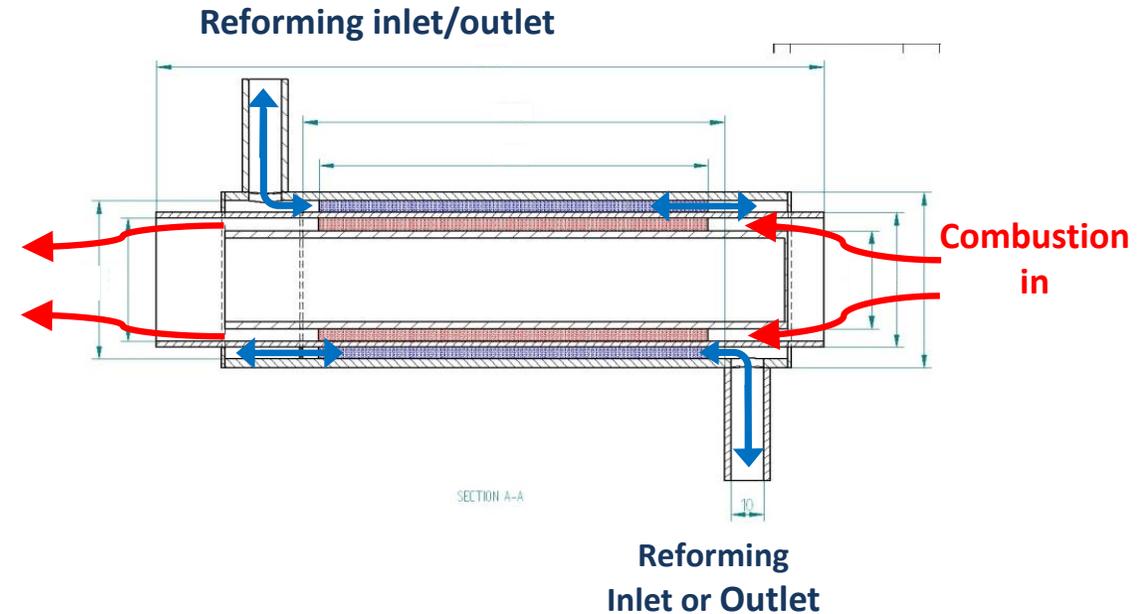


Reformer reactor

Modelling the process-Reformer

Reformer reactor:

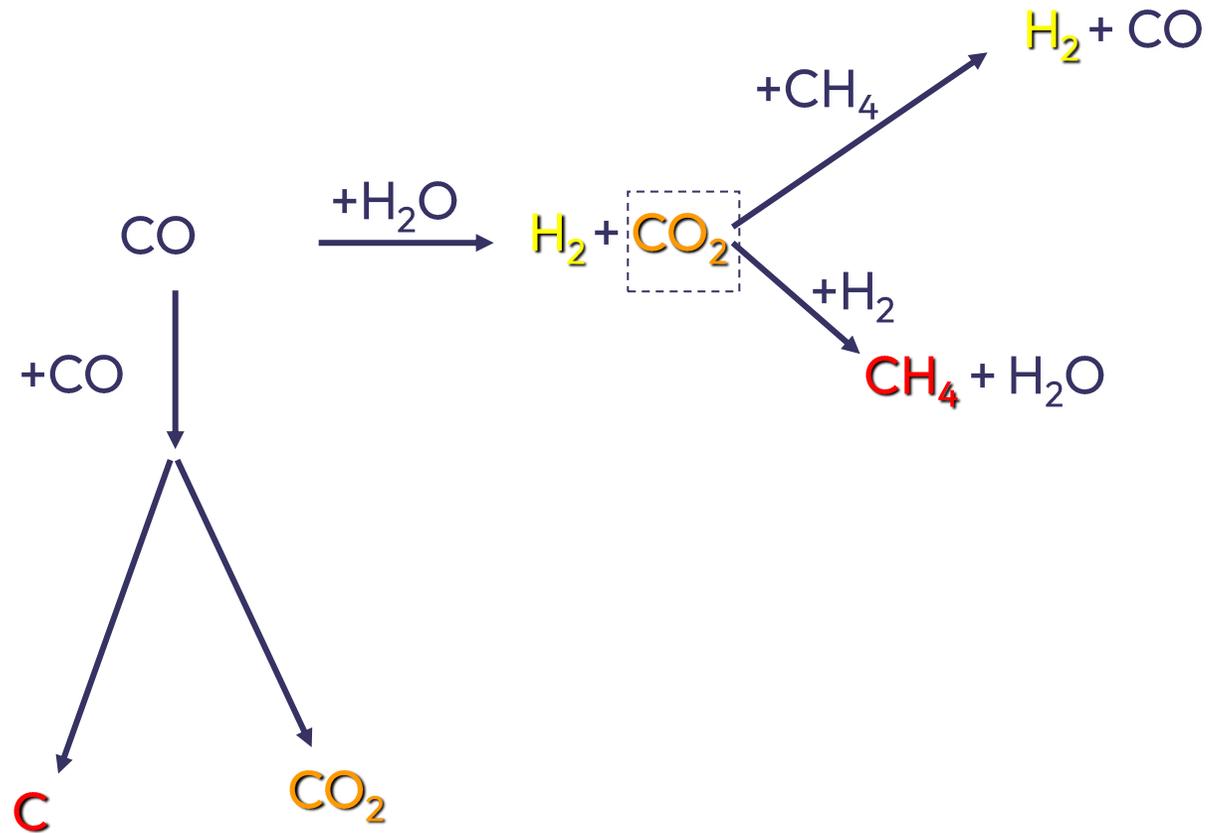
- The reformer consists of a double pipe reactor incorporating separate flow passages in which, both combustion and reforming reactions take place simultaneously.
- The chemical process that produces hydrogen, is the fuel steam reforming which is endothermic. A small portion of the fuel is combusted catalytically together with the reject gas to provide the necessary heat to the reforming reaction.



Suitable for:

- Methane steam reforming/combustion reaction

Modelling the process-WGS



Boudouard or CO disproportionation

Modelling the process-WGS

Water Gas Shift Reactor (WGSR)



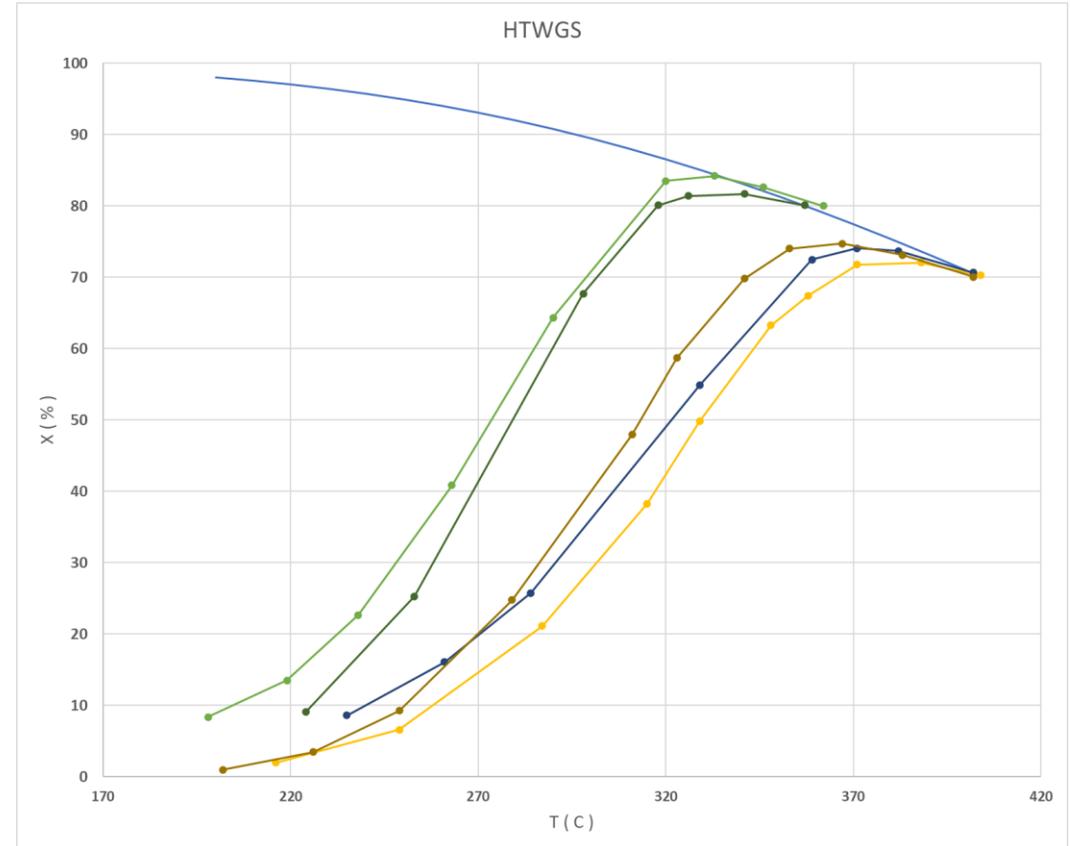
$T_{in} = 400 \text{ }^\circ\text{C}$

PGM catalyst

Heat exchanger type reactor

Typical inlet composition:

| | Vol. % |
|------------------|--------|
| CH ₄ | 2.8 |
| CO | 8.5 |
| H ₂ O | 33.1 |
| CO ₂ | 8.7 |
| H ₂ | 46.9 |



Source: Company information.

Modelling the process-PSA

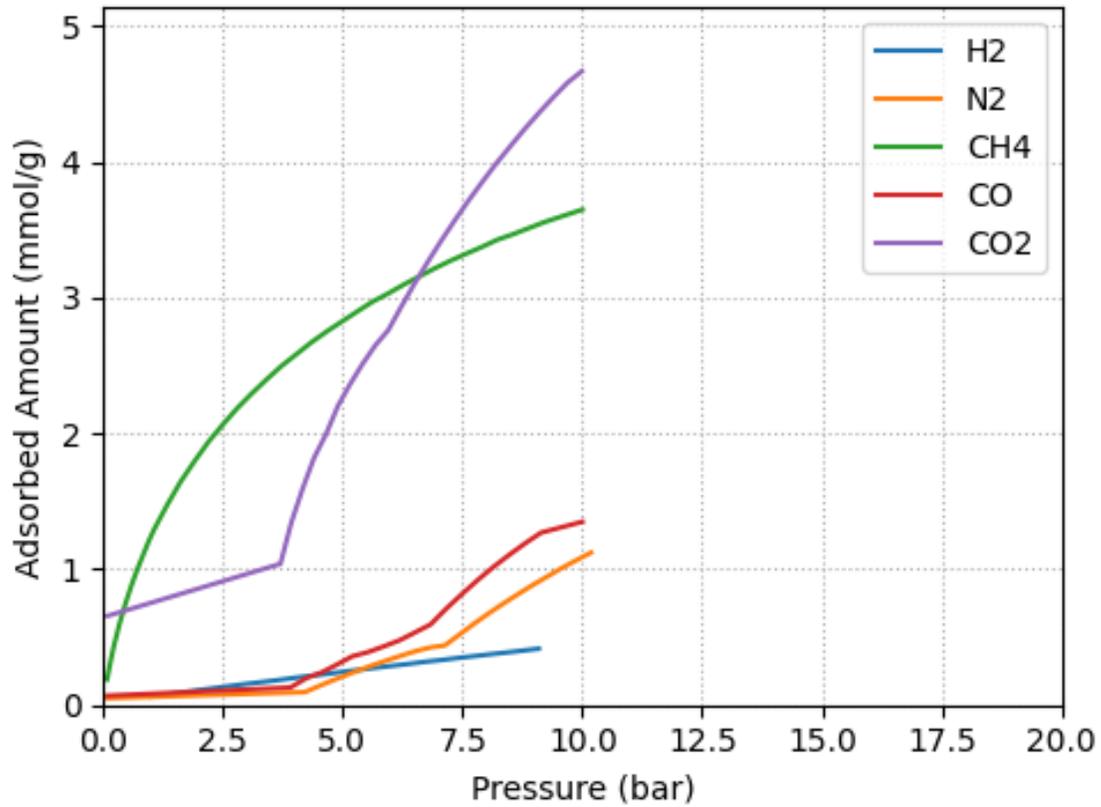
Some theory...

Pressure Swing Adsorption (PSA) is a cyclic gas-separation process that purifies a gas mixture by exploiting the fact that certain molecules are preferentially adsorbed onto solid adsorbents at high pressure and then released when the pressure is lowered.

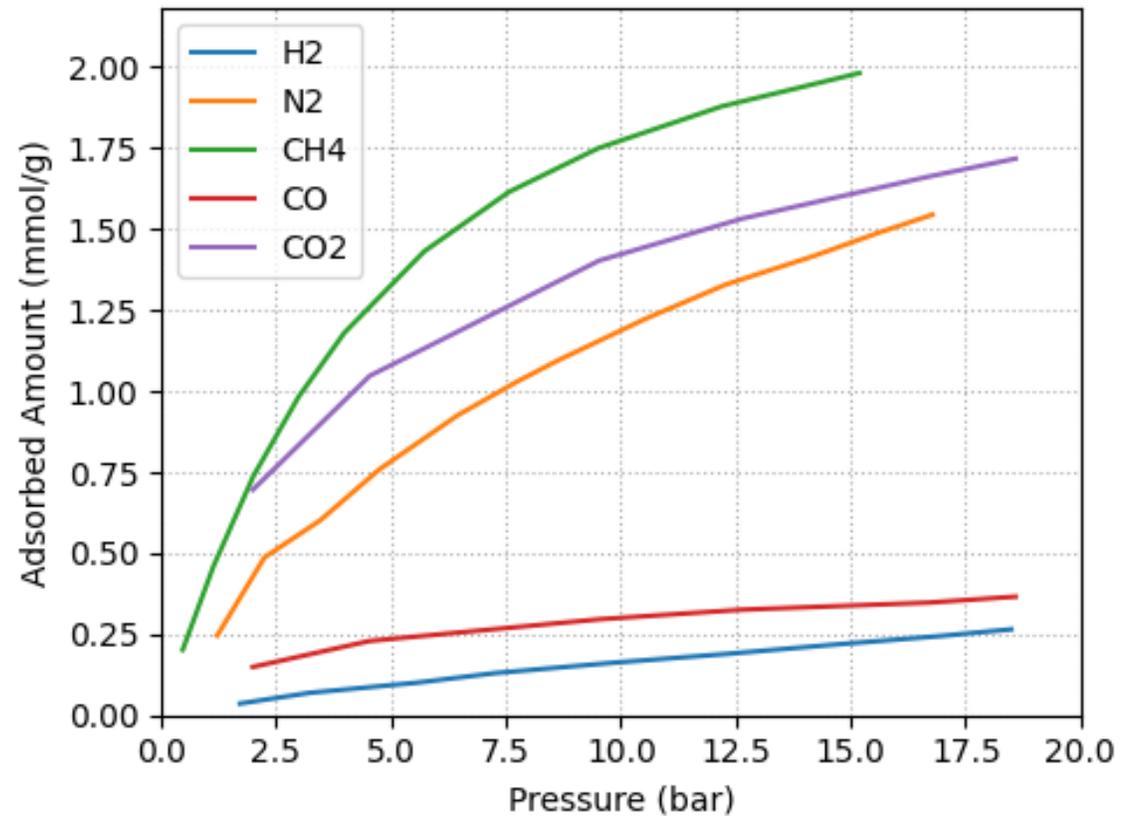
Adsorption isotherms are curves that relate the amount of gas adsorbed on a solid adsorbent to the gas pressure (or concentration) at constant temperature. They help characterize adsorbents, predict PSA performance, calculate surface area, and design layered beds with optimal selectivity.

Modelling the process-PSA

Activated Carbon Isotherms (25°C)

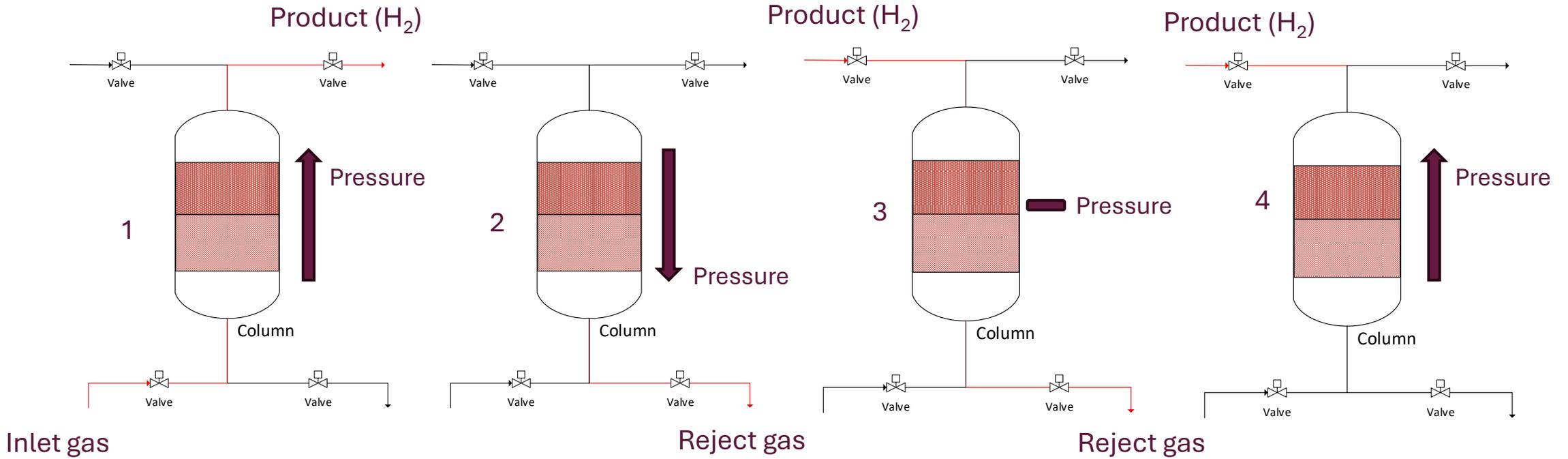


Zeolite 5A Isotherms (25°C)



Source: Company information.

Modelling the process-PSA

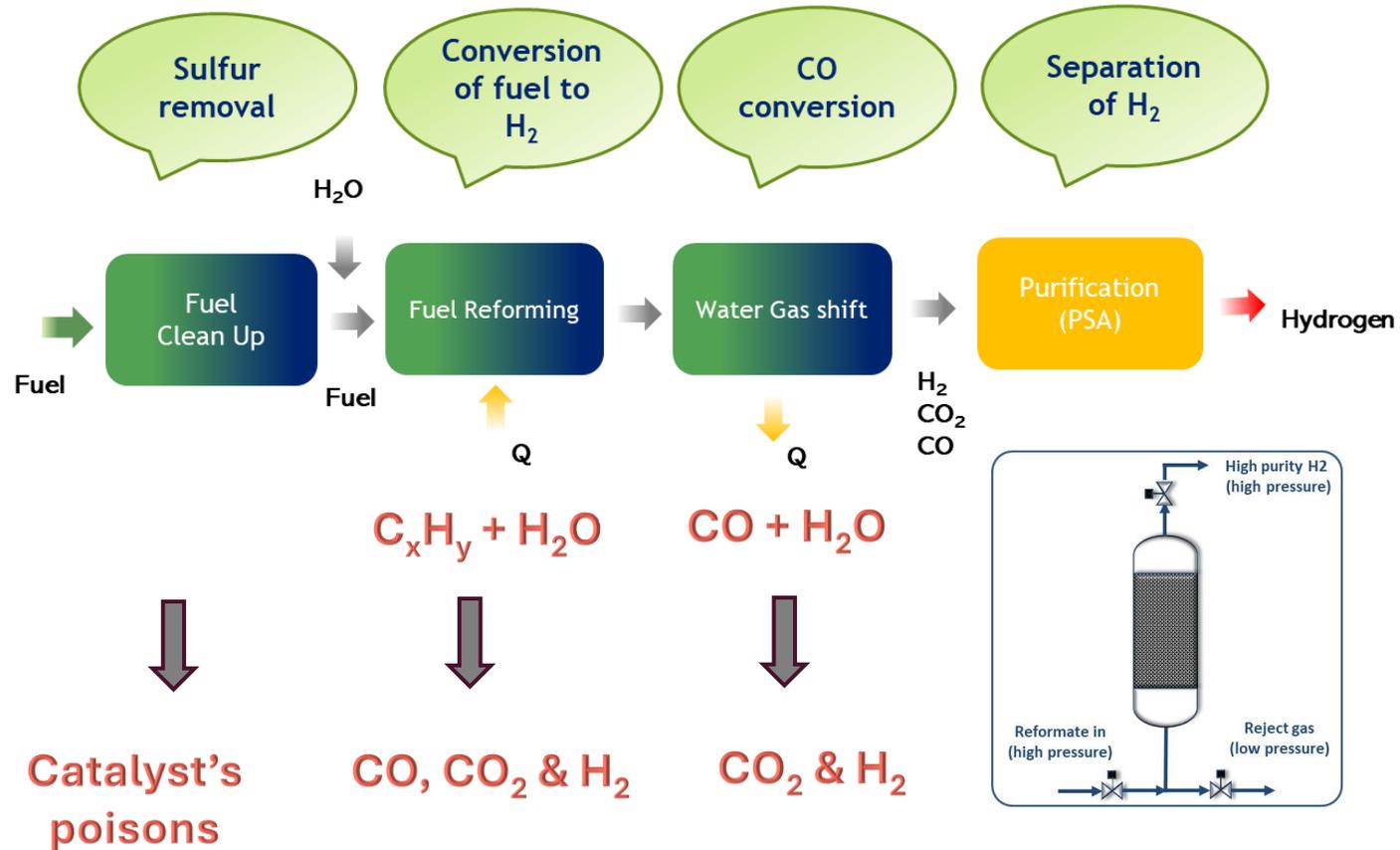


1. Adsorption/Equalization
2. Depressurization/Blowdown
3. Regeneration
4. Repressurization

Inlet gas composition (CH₄, CO, CO₂, H₂) : 4% 1% 20% 75%

Reject gas composition (CH₄, CO, CO₂, H₂): 12%, 3%, 48%, 37%

Modelling the process



Source: Company information.

Modelling the process

Unisim Design software

- **Selection:** of components and reactions
- **Thermodynamic Package:** Peng–Robinson is suitable for high-temperature syngas mixtures.
- **Reactor Models:** Equilibrium reactor for SMR and WGS, for CH_4 , H_2O , CO , CO_2 , H_2 at the given temperature (770 °C) and pressure (1200kPa).
- **Heat Integration:** We'll include heat exchangers to capture the energy exchange. We also model the furnace/combustor simply as a Gibbs reactor that takes fuel (PSA off-gas) and provides heat to the reformer reactor. By adjusting the fraction of fuel burned, we ensure the reformer outlet hits ~770 °C.
- **PSA Unit:** in UniSim we typically model it as a simple separator block that splits the stream into a hydrogen product (at the specified purity and recovery) and a tail gas.
- **Recycle Loops:** The simulation includes a recycle loop where the PSA tail gas is mixed with fresh fuel if needed and sent to the combustor. Convergence of this recycle loop in UniSim confirms that the system can reach a steady state where the energy from tail gas combustion is sufficient (or nearly sufficient) for the reformer needs. If not, we'd add a small fresh fuel makeup in the model.

Hints: By iterating in UniSim, we adjust parameters for optimal results, e.g. we might tweak the steam-to-carbon ratio in the model: increasing steam ratio generally increases CH_4 conversion (by Le Chatelier's principle) and reduces coking risk, but too much steam lowers overall efficiency. The simulation might show that going from $\text{S/C}=2.8$ to 3.5 raises H_2 yield by a few percentage points while increasing furnace duty by 5% – a tradeoff the designer must consider. UniSim also helps size equipment: the flow rates and energy duties from the model inform the sizing of burners, heat exchangers, and catalyst volume.

Source: Company information.

Modelling the process

Unisim Design software

Simulation Basis Manager

Current Fluid Packages

| | | |
|---------|--------|-------------------|
| Basis-1 | NC: 11 | PP: Peng-Robinson |
|---------|--------|-------------------|

View...
Add...
Delete
Copy
View Users...
Import...
Export...

Flowsheet - Fluid Pkg Associations

| Flowsheet | Fluid Pkg To Use |
|-------------|------------------|
| Case (Main) | Basis-1 |

Default Fluid Pkg: Basis-1

Fluid Pkg for New Sub-FlowSheets

Use Default Fluid Pkg Include Column
 Use Parent's Fluid Pkg

Components | **Fluid Pkgs** | Hypotheticals | Hypo Correlation Sets | Oil Manager | Reactions | Component Maps | User Properties

Enter PVT Environment... | Enter Regression Environment... | Return to Simulation Environment...

Simulation Basis Manager

Rxn Components

- Methane
- Ethane
- Propane
- n-Butane
- Hydrogen
- H2O
- CO2
- CO
- Oxygen
- Nitrogen
- I-Butane

Add Comps...

Reactions

- CO shift
- CH4_ref

View Rxn...
Add Rxn...
Delete Rxn
Copy Rxn...
Lock/Unlock...

Reaction Sets

- Reforming
- WGS

View Set...
Add Set...
Delete Set
Copy Set...
Import Set...
Export Set...
Add to FP

Assoc. Fluid Pkgs

- Basis-1

Components | Fluid Pkgs | Hypotheticals | Hypo Correlation Sets | Oil Manager | **Reactions** | Component Maps | User Properties

Enter PVT Environment... | Enter Regression Environment... | Return to Simulation Environment...

Reaction Set: Reforming

Name: Reforming

Set Info

Set Type: Equilibrium Ready Advanced...

Independent

| Active List | OK | Inactive List | Operations Attached |
|-------------|-------------------------------------|---------------|---------------------|
| CH4_ref | <input checked="" type="checkbox"/> | <empty> | Reformer |
| CO shift | <input checked="" type="checkbox"/> | | |
| <empty> | | | |

View Active... | View Inactive... | View Rxn Matrix...
Make Inactive -> | <- Make Active

Reaction Set: WGS

Name: WGS

Set Info

Set Type: Equilibrium Ready Advanced...

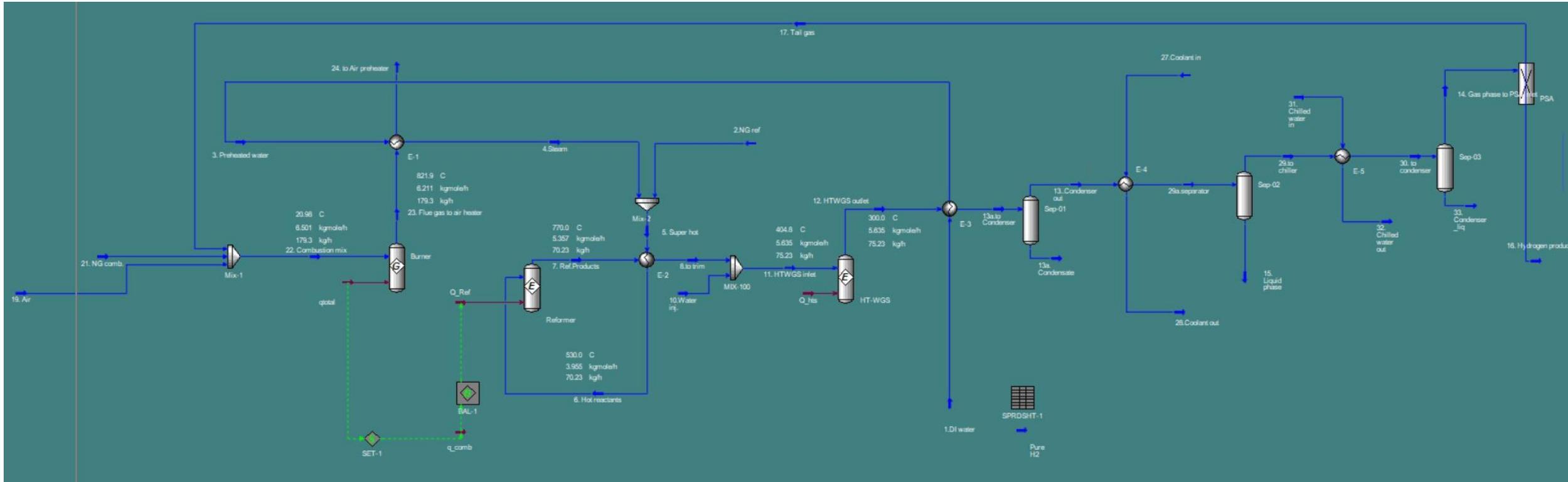
Independent

| Active List | OK | Inactive List | Operations Attached |
|-------------|-------------------------------------|---------------|---------------------|
| CO shift | <input checked="" type="checkbox"/> | <empty> | HT-WGS |
| <empty> | | | |

View Active... | View Inactive... | View Rxn Matrix...
Make Inactive -> | <- Make Active

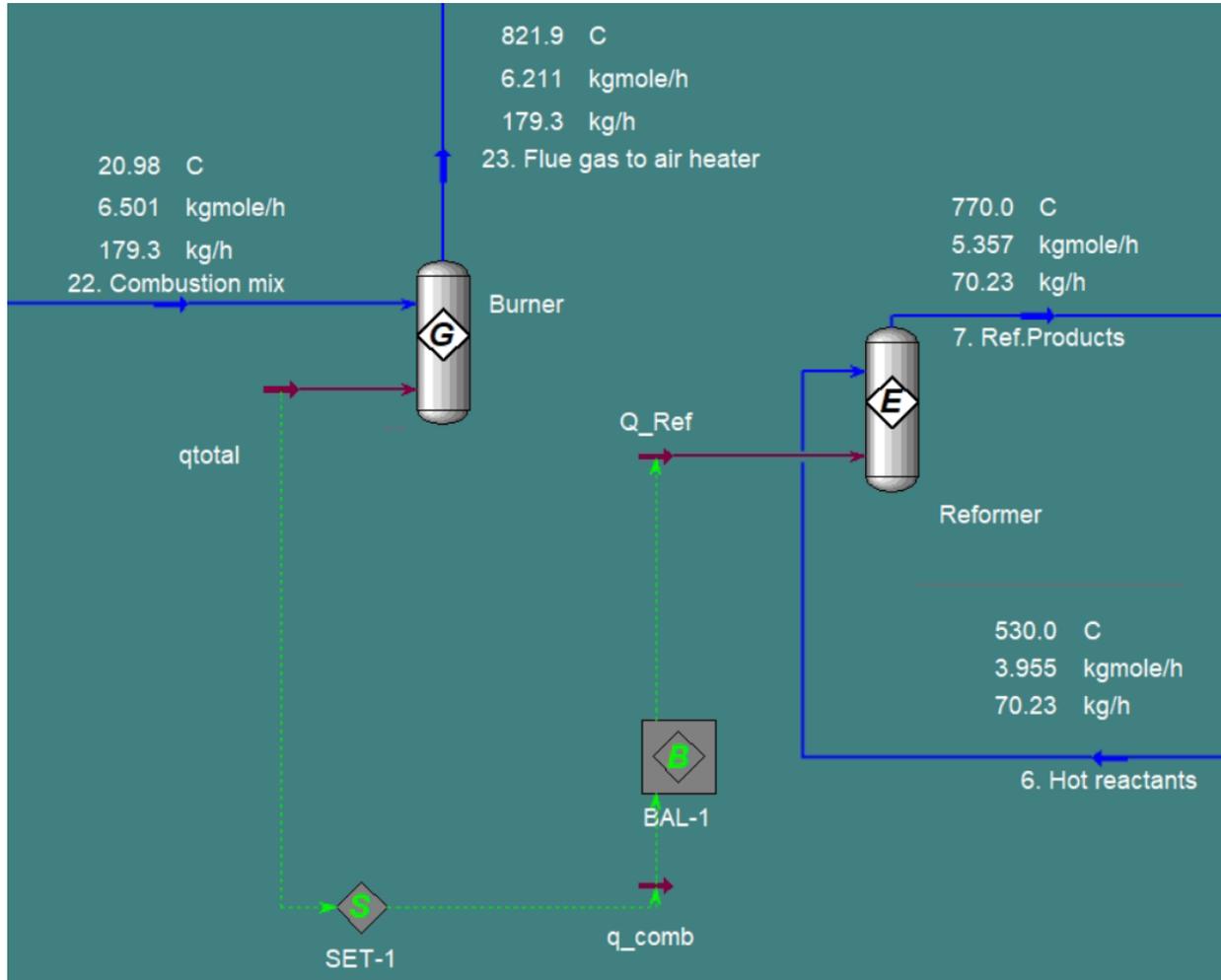
Modelling the process

Unisim Design software



Source: Company information.

Modelling the process



Reaction Set: Global Rxn Set-1

Name: Global Rxn Set-1

Set Info: Set Type: Equilibrium, Ready, Independent, Advanced...

| Active List | OK | Inactive List | Operations Attached |
|-------------|-------------------------------------|---------------|---------------------|
| CO shift | <input checked="" type="checkbox"/> | Comb | Reformer |
| CH4_ref | <input checked="" type="checkbox"/> | Meth | |
| <empty> | | CH4_comb | |

Buttons: View Active..., View Inactive..., View Rxn Matrix..., Make Inactive ->, <- Make Active

Reformer - Global Rxn Set-1

Reactions: Reaction Balance, Reaction Extents, Reaction Balance

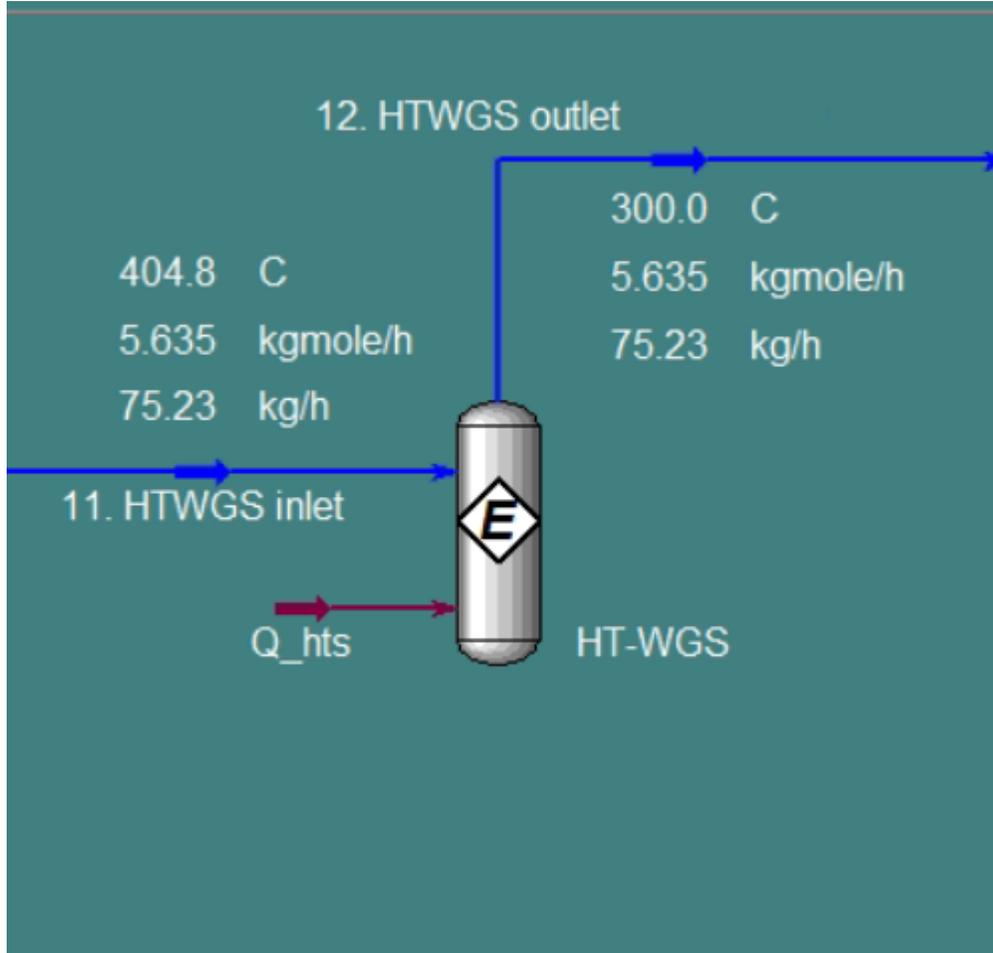
| Results | Act. % Cnv. | Base Comp | Eqm Const. | Rxn Extent |
|----------|-------------|-----------|------------|------------|
| CH4_ref | 80.32 % | Methane | 71.92 | 0.7012 |
| CO shift | 0.00 % | CO | 1.177 | 0.3328 |

Buttons: Design, Reactions, Rating, Worksheet, Dynamics, Delete, OK, Ignored

Unisim Design software

Source: Company information.

Modelling the process



Reaction Set: WGS

Name: **WGS**

Set Info

Set Type: Equilibrium Ready Advanced...
Independent

| Active List | OK | Inactive List | Operations Attached |
|-------------|-------------------------------------|---------------|---------------------|
| CO shift | <input checked="" type="checkbox"/> | <empty> | HT-WGS |
| <empty> | | | |

View Active... View Inactive... View Rxn Matrix...
Make Inactive -> <- Make Active

HT-WGS - WGS

Reactions

Details

Results

Reaction Balance

Reaction Extents Reaction Balance

| | Act. % Cnv. | Base Comp | Eqm Const. | Rxn Extent |
|----------|-------------|-----------|------------|------------|
| CO shift | 93.52 % | CO | 42.00 | 0.3446 |
| | | | | |
| | | | | |
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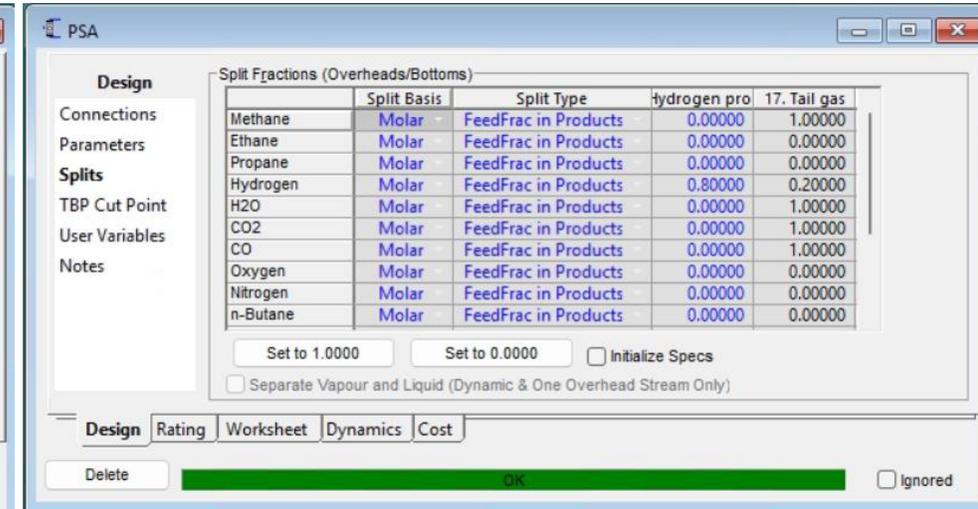
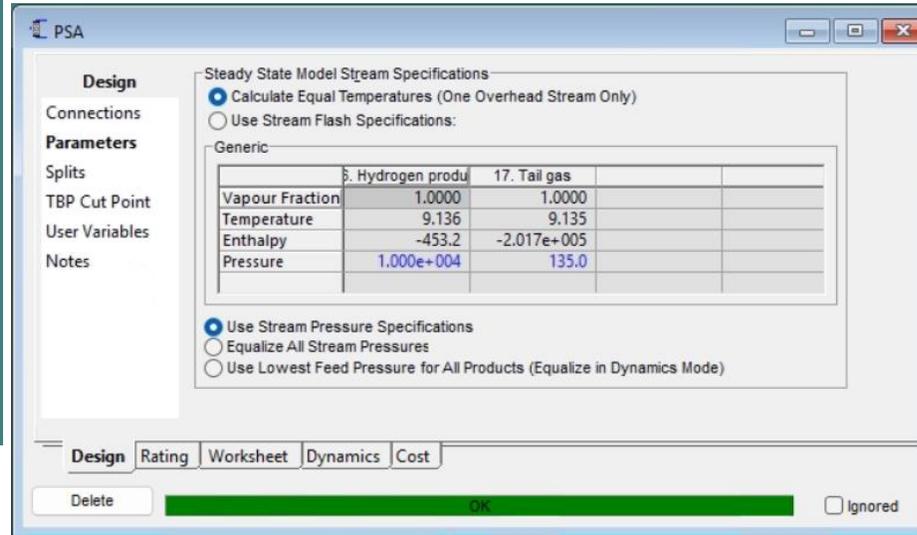
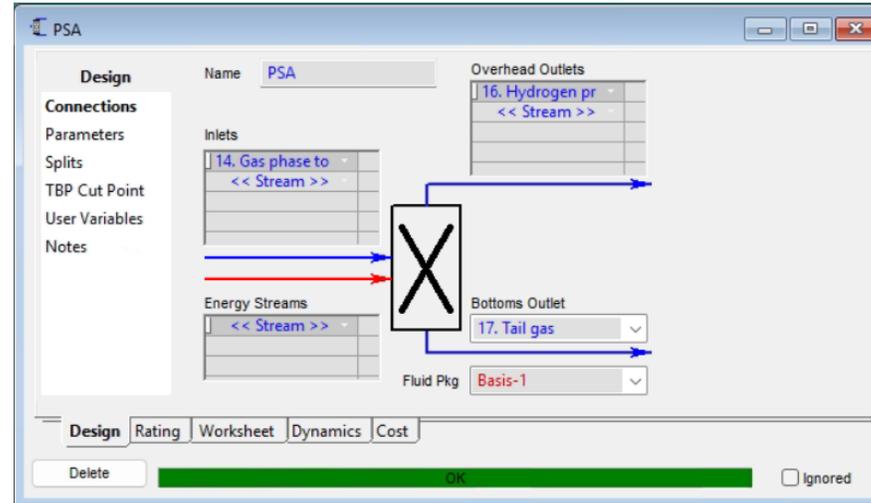
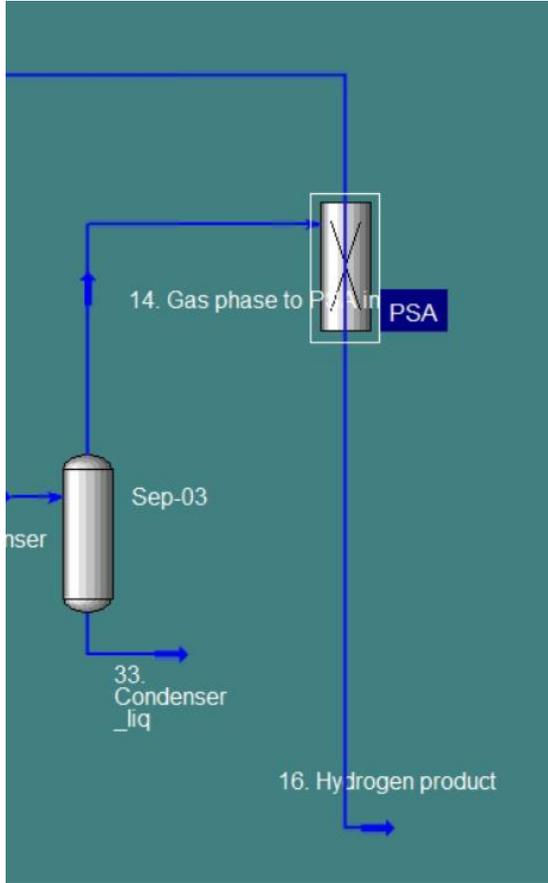
Design **Reactions** Rating Worksheet Dynamics

Delete **OK** Ignored

Unisim Design software

Source: Company information.

Modelling the process



Unisim Design software

Source: Company information.

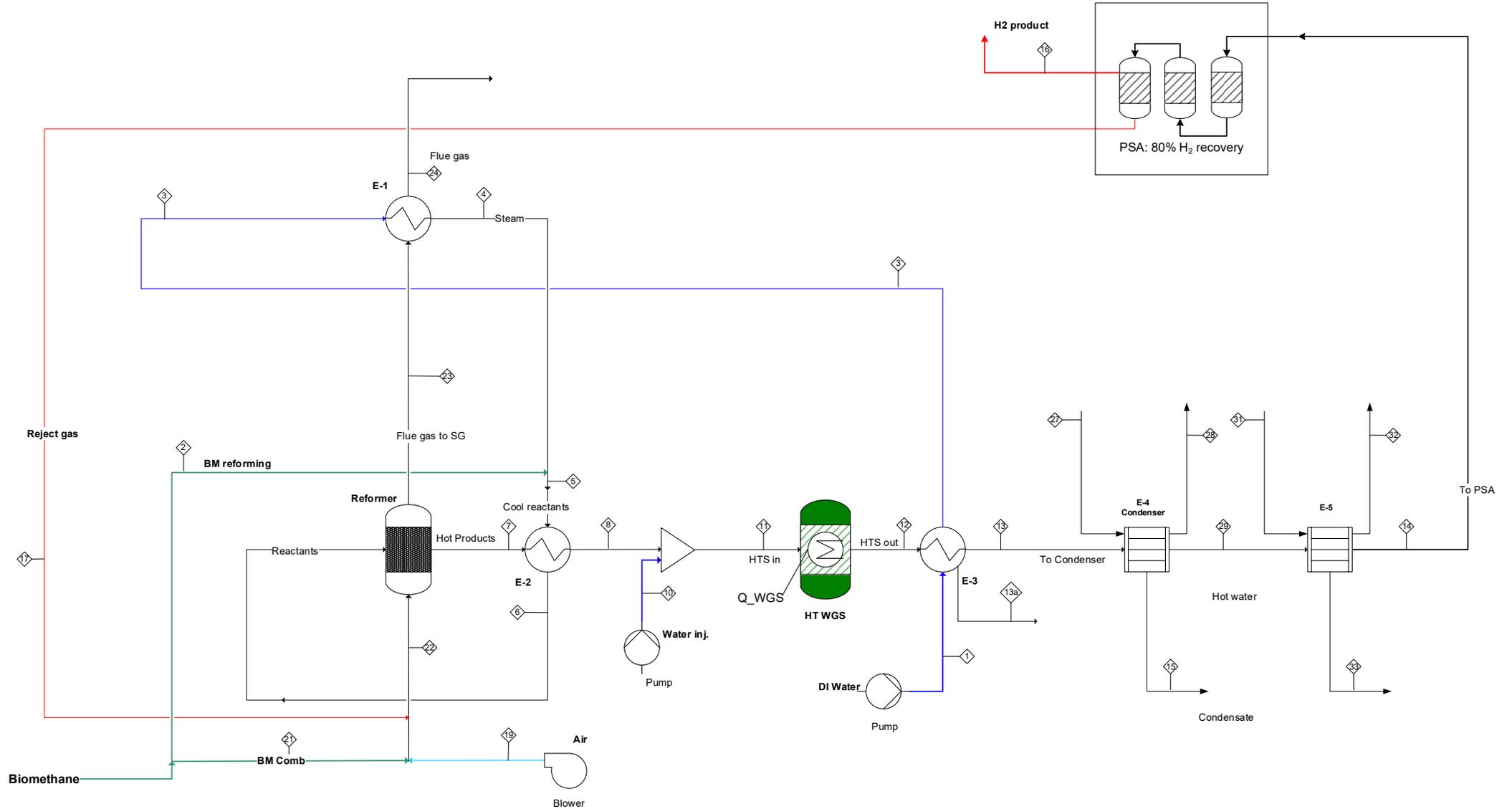
Process flow diagram and Mass and Energy balances

Process Flow Diagram (**PFD**) is a simplified sketch that uses symbols to identify instruments and vessels and to describe the primary flow path through a unit. It illustrates the general plant streams, major equipment and key control loops. They also provide detailed mass/energy balance data along with stream composition and physical properties.

A Piping and Instrumentation Diagram (**P&ID**) is a detailed diagram in the process industry which shows process equipment together with the instrumentation and control devices and control loops.

Process flow diagram and Mass and energy balances

PFD



Source: Company information.

Process flow diagram and Mass and energy balances

| Name | 1.DI water | 2.NG ref | 3. Preheated water | 4.Steam | 5. Super hot | 6. Hot reactants | 7. Ref.Products | 8.to trim | 10.Water inj. | 11. HTWGS inlet | 12. HTWGS outlet | 13..Condenser out | 13a. Condensate | 13a.to Condenser | 14. Gas phase to PSA inlet | 15. Liquid phase | 16. Hydrogen product | 17. Tail gas | 19. Air | 21. NG comb. | 22. Combustion mix | 23. Flue gas to air heater | 24. to Air preheater | 27.Coolant in | 28.Coolant out | 29.to chiller | 29a.separator | 30. to condenser | 31. Chilled water in | 32. Chilled water out | 33. Condenser _liq | |
|----------------------------------|--------------|--------------|--------------------|--------------|--------------|------------------|-----------------|--------------|---------------|-----------------|------------------|-------------------|-----------------|------------------|----------------------------|------------------|----------------------|--------------|--------------|--------------|--------------------|----------------------------|----------------------|---------------|----------------|---------------|---------------|------------------|----------------------|-----------------------|--------------------|------|
| Vapour Fraction | 0 | 1 | 0 | 1 | 1 | 1 | 1 | 1 | 0 | 1 | 1 | 1 | 0 | 1 | 1 | 0 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 0 | 0 | 0 | 1 | 0.65746987 | 0.994711816 | 0 | 0 | 0 |
| Temperature [C] | 5.0 | 25.0 | 101.5 | 250.0 | 192.8 | 530.0 | 770.0 | 489.7 | 25.0 | 404.8 | 300.0 | 180.0 | 180.0 | 180.0 | 10.0 | 35.0 | 9.1 | 9.1 | 25.0 | 25.0 | 21.0 | 821.9 | 199.1 | 25.0 | 32.1 | 35.0 | 35.0 | 10.0 | 5.0 | 7.4 | 10.0 | |
| Pressure [kPa] | 1200 | 1200 | 1195 | 1190 | 1190 | 1185 | 1175 | 1170 | 900 | 900 | 880 | 875 | 875 | 875 | 871 | 873 | 10000 | 135 | 130 | 120 | 120 | 115 | 250 | 248 | 873 | 873 | 871 | 150 | 148 | 871 | | |
| Molar Flow [kgmole/h] | 3.055 | 0.9 | 3.055 | 3.055 | 3.955 | 3.955 | 5.35745686 | 5.35745686 | 0.277544941 | 5.635001801 | 5.635001801 | 5.635001801 | 0 | 5.635001801 | 3.685252007 | 1.930157898 | 2.224819354 | 1.460432653 | 4.9 | 0.14 | 6.500432653 | 6.21039489 | 6.21039489 | 200 | 200 | 3.704843902 | 5.635001801 | 3.704843902 | 20 | 20 | 1.96E-02 | |
| Mass Flow [kg/h] | 55.04 | 15.19 | 55.04 | 55.04 | 70.23 | 70.23 | 70.23 | 70.23 | 5.00 | 75.23 | 75.23 | 75.23 | 0.00 | 75.23 | 40.07 | 34.80 | 4.49 | 35.59 | 141.37 | 2.36 | 179.32 | 179.32 | 3603.02 | 3603.02 | 40.43 | 75.23 | 40.43 | 360.30 | 360.30 | 0.35 | | |
| Std Ideal Liq Vol Flow [m3/h] | 5.51E-02 | 4.82E-02 | 5.51E-02 | 5.51E-02 | 0.103366262 | 0.103366262 | 0.148093575 | 0.148093575 | 5.01E-03 | 0.153103665 | 0.163126899 | 0.163126899 | 0.00E+00 | 0.163126899 | 0.127887621 | 3.49E-02 | 6.42E-02 | 6.37E-02 | 0.163419572 | 7.50E-03 | 0.234603905 | 0.214593144 | 0.214593144 | 3.610291139 | 3.610291139 | 0.128242076 | 0.163126899 | 0.128242076 | 0.361029114 | 0.361029114 | 3.54E-04 | |
| Heat Flow [kJ/h] | -8.77E+05 | -7.62E+04 | -8.54E+05 | -7.14E+05 | -790360.1155 | -734907.4212 | -544224.0992 | -599676.7935 | -7.92E+04 | -678886.1984 | -713101.9161 | -736110.4065 | 0.00E+00 | -736110.4065 | -295533.5966 | -5.49E+05 | -1.01E+03 | -2.95E+05 | -51.42454132 | -1.18E+04 | -306405.2618 | -518275.6196 | -657736.8589 | -57080835.7 | -56970047.11 | -297414.4661 | -846899.0041 | -301150.3762 | -5739221.199 | -5735485.288 | -5.62E+03 | |
| Molar Enthalpy [kJ/kgmole] | -286942.7506 | -84688.60244 | -279411.3298 | -233761.1696 | -199838.2087 | -185817.2999 | -101582.5444 | -111933.1073 | -285393.0777 | -120476.6604 | -126548.6581 | -130631.7961 | -272932.1279 | -130631.7961 | -80193.5922 | -284683.7238 | -453.1568133 | -201669.9667 | -10.49480435 | -84488.80493 | -47136.13357 | -83452.92509 | -105909.0236 | -285404.1785 | -284850.2355 | -80277.19221 | -150292.588 | -81285.57751 | -286961.0599 | -286774.2644 | -286688.9374 | |
| Mass Density [kg/m3] | 1022.532592 | 8.39586963 | 947.1839305 | 5.138606363 | 5.660394458 | 3.167204177 | 1.773701731 | 2.41864383 | 1007.604864 | 2.134870416 | 2.472074546 | 3.12444925 | 874.0272822 | 3.12444925 | 4.030445074 | 1000.293278 | 8.263750176 | 1.405706132 | 1.514050538 | 0.819442742 | 1.355076305 | 0.380464786 | 0.846093687 | 1007.409993 | 1002.028963 | 3.721727352 | 6.903570302 | 4.065859962 | 1022.234396 | 1020.468889 | 1019.253353 | |
| Viscosity [cP] | 1.501203667 | 0.011563874 | 0.274784532 | 1.80E-02 | 1.31E-02 | 0.024141402 | 2.89E-02 | 2.28E-02 | 0.890438924 | 2.11E-02 | 1.85E-02 | 1.55E-02 | 0.149297086 | 1.55E-02 | 1.06E-02 | 0.718491259 | 8.62E-03 | 1.40E-02 | 1.88E-02 | 0.011375772 | 1.74E-02 | 4.33E-02 | 2.32E-02 | 0.890438966 | 0.762112334 | 1.13E-02 | <empty> | <empty> | 1.50120411 | 1.398963045 | 1.234307796 | |
| Kinematic Viscosity [cSt] | 1.468123049 | 1.37732891 | 0.290106835 | 3.506335122 | 2.313355192 | 7.622306829 | 16.29091509 | 9.438452857 | 0.883718366 | 9.89871623 | 7.497920952 | 4.966517924 | 0.170815133 | 4.966517924 | 2.63958275 | 0.718280603 | 1.043586175 | 9.93900012 | 12.43975402 | 13.88232712 | 12.87270179 | 113.759624 | 27.43818002 | 0.883889352 | 0.760569167 | 3.03780995 | <empty> | <empty> | 1.468551749 | 1.3709022 | 1.210992137 | |
| Mass Heat Capacity [kJ/kg-C] | 4.320375272 | 2.207746649 | 4.398486522 | 2.085769321 | 2.169187249 | 2.529434606 | 2.904361348 | 2.724397895 | 4.312582363 | 2.624908036 | 2.589558528 | 2.509959702 | 4.739651144 | 2.509959702 | 4.309483028 | 14.43007457 | 1.390645682 | 1.01388398 | 2.141673312 | 1.104243414 | 1.333440476 | 1.152791538 | 4.313295725 | 4.313512028 | 2.841676249 | 3.520716786 | 2.844597639 | 4.321321489 | 4.319700237 | 4.31037425 | | |
| Thermal Conductivity [W/m-K] | 0.577988737 | 3.40E-02 | 0.681297073 | 3.96E-02 | 3.98E-02 | 8.11E-02 | 0.182587683 | 0.14149498 | 0.611014986 | 0.123561088 | 0.117031895 | 9.56E-02 | 0.676602971 | 9.56E-02 | 9.93E-02 | 0.625068154 | 0.174575923 | 4.65E-02 | 2.59E-02 | 3.31E-02 | 3.01E-02 | 7.44E-02 | 3.52E-02 | 0.611014416 | 0.621195361 | 0.10586831 | <empty> | <empty> | 0.577987829 | 0.5823102 | 0.587545327 | |
| Heat Of Vaporization [kJ/kgmole] | 36219.42503 | 7196.154314 | 36231.56517 | 36243.85816 | 51031.48058 | 51034.57617 | 38296.6798 | 38295.95405 | 37012.16558 | 40059.85141 | 38214.82581 | 38215.16291 | 73137.60954 | 38215.16291 | 13130.12865 | 39894.84076 | <empty> | 18603.41064 | 5823.080866 | 9400.811649 | 15666.08183 | 30611.1443 | 30624.11055 | 39667.7551 | 39681.19165 | 14288.46714 | 38215.14429 | 14288.65698 | 40465.18873 | 40484.70169 | 42920.27284 | |
| Comp Mole Frac (Methane) | 0 | 0.97 | 0 | 0 | 0.220733249 | 0.220733249 | 3.21E-02 | 3.21E-02 | 0 | 3.05E-02 | 3.05E-02 | 3.05E-02 | 2.56E-06 | 3.05E-02 | 4.66E-02 | 3.92E-09 | 0 | 0.117616901 | 0 | 0.97 | 4.73E-02 | 9.68E-38 | 9.68E-38 | 0 | 0 | 4.64E-02 | 3.05E-02 | 4.64E-02 | 0 | 0 | 8.31E-10 | |
| Comp Mole Frac (Ethane) | 0 | 0 | 0 | 0.00E+00 | 0.00E+00 | 0 | 0.00E+00 | 0.00E+00 | 0 | 0 | 0.00E+00 | 0.00E+00 | 0 | 0 | 0 | 0 | 0.00E+00 | 0.00E+00 | 0.00E+00 | 0 | 0.00E+00 | 1.98E-47 | 1.98E-47 | 0 | 0 | 0.00E+00 | 0 | 0 | 0 | 0 | 0 | |
| Comp Mole Frac (Propane) | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 1.70E-48 | 1.70E-48 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | |
| Comp Mole Frac (n-Butane) | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 3.11E-87 | 3.11E-87 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | |
| Comp Mole Frac (Hydrogen) | 0 | 0.00E+00 | 0 | 0.00E+00 | 0.00E+00 | 0.00E+00 | 0.454782778 | 0.454782778 | 0 | 0.432383023 | 0.493530067 | 4.94E-01 | 3.64E-04 | 0.493530067 | 7.55E-01 | 9.60E-06 | 1 | 3.81E-01 | 0.00E+00 | 0.00E+00 | 8.56E-02 | 1.37E-09 | 1.37E-09 | 0 | 0 | 0.750645467 | 0.493530067 | 0.750645467 | 0 | 0 | 4.53E-06 | |
| Comp Mole Frac (H2O) | 1 | 0 | 1 | 1 | 0.772439949 | 0.772439949 | 0.377227069 | 0.377227069 | 1 | 0.407900968 | 0.346753924 | 0.346753924 | 0.999104364 | 0.346753924 | 1.48E-03 | 0.999366637 | 0 | 3.73E-03 | 0 | 0 | 8.39E-04 | 1.89E-01 | 1.89E-01 | 1 | 1 | 6.75E-03 | 0.346753924 | 6.75E-03 | 1 | 1 | 0.998849627 | |
| Comp Mole Frac (CO2) | 0.00 | 0.03 | 0.00 | 0.00 | 0.01 | 0.01 | 0.07 | 0.07 | 0.00 | 0.06 | 0.12 | 0.12 | 0.00 | 0.12 | 0.19 | 0.00 | 0.00 | 0.48 | 0.00 | 0.03 | 0.11 | 0.17 | 0.17 | 0.00 | 0.00 | 0.19 | 0.12 | 0.19 | 0.00 | 0.00 | 0.00 | |
| Comp Mole Frac (CO) | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.07 | 0.07 | 0.00 | 0.07 | 0.00 | 0.00 | 0.00 | 0.00 | 0.01 | 0.00 | 0.00 | 0.02 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.01 | 0.00 | 0.01 | 0.00 | 0.00 | 0.00 | |
| Comp Mole Frac (Oxygen) | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.21 | 0.00 | 0.16 | 0.02 | 0.02 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 |
| Comp Mole Frac (Nitrogen) | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.79 | 0.00 | 0.60 | 0.62 | 0.62 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | |
| Comp Mole Frac (i-Butane) | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | 0.00 | |

Source: Company information.

Process flow diagram and Mass and energy balances

| Name | 1.DI water | 2.NG ref | 3. Preheated water | 4.Steam | 6. Hot reactants | 7. Ref.Products | 10.Water inj. | 11. HTWGS inlet | 12. HTWGS outle | Hydrogen produ | 17. Tail gas | 19. Air | 21. NG comb. |
|-----------------------------------|--------------|--------------|--------------------|--------------|------------------|-----------------|---------------|-----------------|-----------------|----------------|--------------|--------------|--------------|
| Vapour Fraction | 0 | 1 | 0 | 1 | 1 | 1 | 0 | 1 | 1 | 1 | 1 | 1 | 1 |
| Temperature [C] | 5.0 | 25.0 | 101.5 | 250.0 | 530.0 | 770.0 | 25.0 | 404.8 | 300.0 | 9.1 | 9.1 | 25.0 | 25.0 |
| Pressure [kPa] | 1200 | 1200 | 1195 | 1190 | 1185 | 1175 | 900 | 900 | 880 | 10000 | 135 | 130 | 120 |
| Molar Flow [kgmole/h] | 3.055 | 0.9 | 3.055 | 3.055 | 3.955 | 5.35745686 | 0.277544941 | 5.635001801 | 5.635001801 | 2.224819354 | 1.460432653 | 4.9 | 0.14 |
| Mass Flow [kg/h] | 55.04 | 15.19 | 55.04 | 55.04 | 70.23 | 70.23 | 5.00 | 75.23 | 75.23 | 4.49 | 35.59 | 141.37 | 2.36 |
| Std Ideal Liq Vol Flow [m3/h] | 5.51E-02 | 4.82E-02 | 5.51E-02 | 5.51E-02 | 0.103366262 | 0.148093575 | 5.01E-03 | 0.153103665 | 0.163126899 | 6.42E-02 | 6.37E-02 | 0.163419572 | 7.50E-03 |
| Heat Flow [kJ/h] | -8.77E+05 | -7.62E+04 | -8.54E+05 | -7.14E+05 | -734907.4212 | -544224.0992 | -7.92E+04 | -678886.1984 | -713101.9161 | -1.01E+03 | -2.95E+05 | -51.42454132 | -1.18E+04 |
| Molar Enthalpy [kJ/kgmole] | -286942.7506 | -84688.60244 | -279411.3298 | -233761.1696 | -185817.2999 | -101582.5444 | -285393.0777 | -120476.6604 | -126548.6581 | -453.1568133 | -201669.9667 | -10.49480435 | -84488.80493 |
| Mass Density [kg/m3] | 1022.532592 | 8.39586963 | 947.1839305 | 5.138606363 | 3.167204177 | 1.773701731 | 1007.604864 | 2.134870416 | 2.472074546 | 8.263750176 | 1.405706132 | 1.514050538 | 0.819442742 |
| Viscosity [cP] | 1.501203667 | 0.011563874 | 0.274784532 | 1.80E-02 | 0.024141402 | 2.89E-02 | 0.890438924 | 2.11E-02 | 1.85E-02 | 8.62E-03 | 1.40E-02 | 1.88E-02 | 0.011375772 |
| Kinematic Viscosity [cSt] | 1.468123049 | 1.37732891 | 0.290106835 | 3.506335122 | 7.622306829 | 16.29091509 | 0.883718366 | 9.89871623 | 7.497920952 | 1.043586175 | 9.93900012 | 12.43975402 | 13.88232712 |
| Mass Heat Capacity [kJ/kg-C] | 4.320375272 | 2.207746649 | 4.398486522 | 2.065769321 | 2.529434606 | 2.904361348 | 4.312582363 | 2.624908036 | 2.589558528 | 14.43007457 | 1.390645682 | 1.01358398 | 2.141673312 |
| Thermal Conductivity [W/m-K] | 0.577988737 | 3.40E-02 | 0.681297073 | 3.96E-02 | 8.11E-02 | 0.182587683 | 0.611014986 | 0.123561088 | 0.117031895 | 0.174575923 | 4.65E-02 | 2.59E-02 | 3.31E-02 |
| Heat Of Vapourization [kJ/kgmole] | 36219.42503 | 7196.154314 | 36231.56517 | 36243.85816 | 51034.57617 | 38296.6798 | 37012.16558 | 40059.85141 | 38214.82581 | <empty> | 18603.41064 | 5823.080866 | 9400.811649 |
| Comp Mole Frac (Methane) | 0.0% | 97.0% | 0.0% | 0.0% | 22.1% | 3.2% | 0.0% | 3.0% | 3.0% | 0.0% | 11.8% | 0.0% | 97.0% |
| Comp Mole Frac (Ethane) | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% |
| Comp Mole Frac (Propane) | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% |
| Comp Mole Frac (n-Butane) | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% |
| Comp Mole Frac (Hydrogen) | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 45.5% | 0.0% | 43.2% | 49.4% | 100.0% | 38.1% | 0.0% | 0.0% |
| Comp Mole Frac (H2O) | 100.0% | 0.0% | 100.0% | 100.0% | 77.2% | 37.7% | 100.0% | 40.8% | 34.7% | 0.0% | 0.4% | 0.0% | 0.0% |
| Comp Mole Frac (CO2) | 0.0% | 3.0% | 0.0% | 0.0% | 0.7% | 6.7% | 0.0% | 6.4% | 12.5% | 0.0% | 48.1% | 0.0% | 3.0% |
| Comp Mole Frac (CO) | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 6.9% | 0.0% | 6.5% | 0.4% | 0.0% | 1.6% | 0.0% | 0.0% |
| Comp Mole Frac (Oxygen) | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 21.0% | 0.0% |
| Comp Mole Frac (Nitrogen) | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 79.0% | 0.0% |
| Comp Mole Frac (i-Butane) | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% | 0.0% |

Source: Company information.

Initial equipment sizing: BoP

| Name | 1. DI water | 19. Air |
|-----------------------------------|--------------|--------------|
| Vapour Fraction | 0 | 1 |
| Temperature [C] | 5.0 | 25.0 |
| Pressure [kPa] | 1200 | 130 |
| Molar Flow [kgmole/h] | 3.055 | 4.9 |
| Mass Flow [kg/h] | 55.04 | 141.37 |
| Std Ideal Liq Vol Flow [m3/h] | 5.51E-02 | 0.163419572 |
| Heat Flow [kJ/h] | -8.77E+05 | -51.42454132 |
| Molar Enthalpy [kJ/kgmole] | -286942.7506 | -10.49480435 |
| Mass Density [kg/m3] | 1022.532592 | 1.514050538 |
| Viscosity [cP] | 1.501203667 | 1.88E-02 |
| Kinematic Viscosity [cSt] | 1.468123049 | 12.43975402 |
| Mass Heat Capacity [kJ/kg-C] | 4.320375272 | 1.01358398 |
| Thermal Conductivity [W/m-K] | 0.577988737 | 2.59E-02 |
| Heat Of Vapourization [kJ/kgmole] | 36219.42503 | 5823.080866 |
| Comp Mole Frac (Methane) | 0% | 0% |
| Comp Mole Frac (Ethane) | 0% | 0% |
| Comp Mole Frac (Propane) | 0% | 0% |
| Comp Mole Frac (n-Butane) | 0% | 0% |
| Comp Mole Frac (Hydrogen) | 0% | 0% |
| Comp Mole Frac (H2O) | 100% | 0% |
| Comp Mole Frac (CO2) | 0% | 0% |
| Comp Mole Frac (CO) | 0% | 0% |
| Comp Mole Frac (Oxygen) | 0% | 21% |
| Comp Mole Frac (Nitrogen) | 0% | 79% |
| Comp Mole Frac (i-Butane) | 0% | 0% |

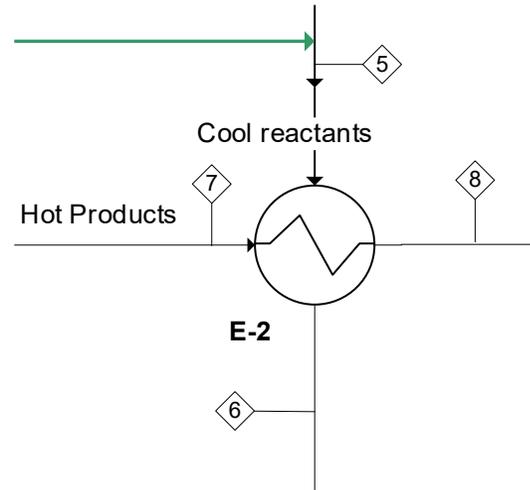
Example of equipment for preliminary sizing:

- Water pump preliminary sizing
- Air blower preliminary sizing

Source: Company information.

Initial equipment sizing-HEx

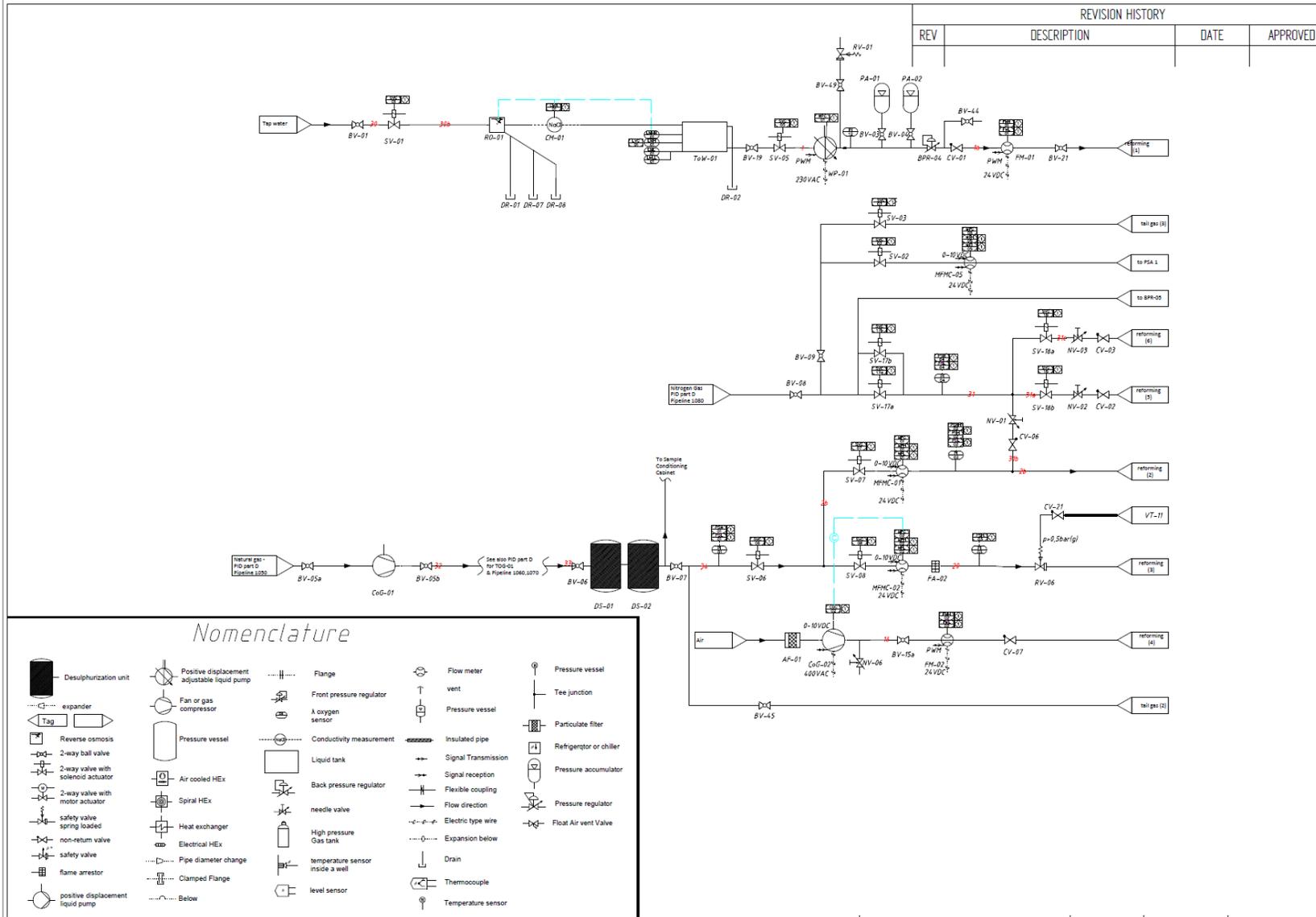
- **Heat exchanger type**
(shell-and-tube, plate, air-cooled, spiral, etc.)
- **Duty type:** heating, cooling, condensation, vaporization.
- **Hot & cold fluids:** identify both streams and their roles.
- **Operating conditions:**
Flow rates, inlet/outlet temperatures, pressures, Phase states (liquid, gas, two-phase), allowable pressure drops (tube side & shell side). [
- **Constraints:** space limits, materials, corrosion, fouling tendency.



| Name | 5. Super hot | 6. Hot reactants | 7. Ref.Products | 8.to trim |
|-----------------------------------|--------------|------------------|-----------------|--------------|
| Vapour Fraction | 1 | 1 | 1 | 1 |
| Temperature [C] | 192.8 | 530.0 | 770.0 | 489.7 |
| Pressure [kPa] | 1190 | 1185 | 1175 | 1170 |
| Molar Flow [kgmole/h] | 3.955 | 3.955 | 5.35745686 | 5.35745686 |
| Mass Flow [kg/h] | 70.23 | 70.23 | 70.23 | 70.23 |
| Std Ideal Liq Vol Flow [m3/h] | 0.103366262 | 0.103366262 | 0.148093575 | 0.148093575 |
| Heat Flow [kJ/h] | -790360.1155 | -734907.4212 | -544224.0992 | -599676.7935 |
| Molar Enthalpy [kJ/kgmole] | -199838.2087 | -185817.2999 | -101582.5444 | -111933.1073 |
| Mass Density [kg/m3] | 5.660394458 | 3.167204177 | 1.773701731 | 2.41864383 |
| Viscosity [cP] | 1.31E-02 | 0.024141402 | 2.89E-02 | 2.28E-02 |
| Kinematic Viscosity [cSt] | 2.313355192 | 7.622306829 | 16.29091509 | 9.438452857 |
| Mass Heat Capacity [kJ/kg-C] | 2.169187249 | 2.529434606 | 2.904361348 | 2.724397895 |
| Thermal Conductivity [W/m-K] | 3.98E-02 | 8.11E-02 | 0.182587683 | 0.14149498 |
| Heat Of Vapourization [kJ/kgmole] | 51031.48058 | 51034.57617 | 38296.6798 | 38295.95405 |
| Comp Mole Frac (Methane) | 0.220733249 | 0.220733249 | 3.21E-02 | 3.21E-02 |
| Comp Mole Frac (Ethane) | 0.00E+00 | 0 | 0.00E+00 | 0.00E+00 |
| Comp Mole Frac (Propane) | 0 | 0 | 0 | 0 |
| Comp Mole Frac (n-Butane) | 0 | 0 | 0 | 0 |
| Comp Mole Frac (Hydrogen) | 0.00E+00 | 0.00E+00 | 0.454782778 | 0.454782778 |
| Comp Mole Frac (H2O) | 0.772439949 | 0.772439949 | 0.377227069 | 0.377227069 |
| Comp Mole Frac (CO2) | 0.01 | 0.01 | 0.07 | 0.07 |
| Comp Mole Frac (CO) | 0.00 | 0.00 | 0.07 | 0.07 |
| Comp Mole Frac (Oxygen) | 0.00 | 0.00 | 0.00 | 0.00 |
| Comp Mole Frac (Nitrogen) | 0.00 | 0.00 | 0.00 | 0.00 |
| Comp Mole Frac (i-Butane) | 0.00 | 0.00 | 0.00 | 0.00 |

Source: Company information.

Piping and Instrumentation diagram



Source: Company information.

Hazardous Operation – Risk assessment

| Aspect | HAZOP | Risk Analysis |
|-----------------|------------------------------------------|-------------------------------------------------|
| Goal | Identify deviations & operability issues | Evaluate and prioritize risks |
| Timing | Detailed design stage | Any project stage |
| Inputs | P&ID, control logic, process data | Varies (concept design to detailed engineering) |
| Approach | Guidewords, structured brainstorming | Risk matrices, models (QRA), expert judgement |
| Output | Actionable deviations & safeguards | Risk ranking, mitigation strategy |
| Scope | Narrow, process-specific | Broad, system-level or project-level |

HAZOP example

| Node | 1b Reformer reforming feed: From BV-01 and BV-06 (through SV-07) to FM-01 and RV-02 | | | | | | | | |
|---------------|-------------------------------------------------------------------------------------|-----------|--------------------------------------------|-----------------------------------------------------------------------------------------------------------------|--------------------------------|-----------------------|----------|-----------|----------------------------------------------------------------------------------|
| Drawing | xx-xxx-xxx-xx | | | | | | | | |
| Parameter | Nr. | Deviation | Cause | Consequences | Safeguards | Additional safeguards | Severity | Frequency | Recommendations/ Actions |
| Flow | 1 | High | WP-01 malfunction | High water flow. Low quality steam. Low reformer temperature | FMH01 | TSL25, TSL23 | Low | moderate | |
| | 1 | | FM-01 instrument malfunction | | TSL25 | TSL23 | Low | Low | |
| | 2 | | MFMC-01 instrument malfunction | High Natural gas flowrate in reforming side. Low reformer temperature, Low quality hydrogen, | MFH01 | | Low | Low | |
| | 2 | | SV-16 fails to shut-off | | Nitrogen in the reforming side | PSH04 | | Low | Low |
| | 3 | Low | WP-01 malfunction/failure | Low or no water flow. High reformer temperature, Catalyst coking, High CO at WGS reactor effluent | FML01 | TSH25, TSH23 | Low | Moderate | |
| | 3 | | FM-01 instrument malfunction | | TSH25 | TSH23 | Low | Low | |
| | 3 | | ToW-01 low water level | | LML01 | | Low | Low | |
| | 3 | | SV-05 fails open | | FML01 | TSH25, TSH23 | Low | Low | |
| | 3 | | Air bubbles in tubing to WP-01 | FML01 | TSH25, TSH23 | Low | Low | | |
| | 4 | | MFMC-01 instrument malfunction | Low Natural gas flowrate in reforming side. High reformer temperature, Low hydrogen flowrate | MFL01 | PSL02 | Low | Low | |
| | 4 | | SV-07 fails open | | MFL01 | PSL02 | Low | Low | |
| | 4 | | SV-06 fails open | | MFL01 | PSL02 | Low | Low | |
| | Temperature | 5 | High | None | | | | | |
| 6 | | Low | None | | | | | | |
| Pressure | 7 | High | Downstream blokage, | Limited or no reactants flow. Reforming temperature high. Possible reformer shell or combustion tube rupture | PSH02 | FML01, MFL01 | High | Low | |
| | 8 | High | High flowrate of NG | Reforming temperature low. Catalyst coking. Possible reformer shell or combustion tube rupture | MFH01 | | High | Low | |
| | 9 | Low | Natural gas compressor failure. | High reformer temperature. Low or no hydrogen production. | PSL01 | MFL01 | Low | Moderate | |
| | 10 | Low | No or low water flowrate | High reformer temperature. Low hydrogen production. Catalyst coking | FML01 | | Low | Moderate | |
| | 11 | Low | Natural gas compressor failure. | Low combustion temperature or no combustion | PSL01 | MFL01 | Low | Moderate | |
| Level | 12 | High | SV-04 fails close | High water level. Possible water overflow | LMH01 | LMHH01 | Low | Low | |
| | 13 | Low | SV-01 fails open | Low water level in ToW-01 | LML01 | LMLL01 | Low | Low | |
| | 13 | | SV-02 fails open | | LML01 | LMLL01 | Low | Low | |
| | 13 | | SV-03 fails open | | LML01 | LMLL01 | Low | Low | |
| 13 | Prefilter PF-01 partial/full blockage | LML01 | LMLL01 | Low | High | Check/clean filter | | | |
| Concentration | 14 | High | None | | | | | | |
| | 15 | Low | None | | | | | | |
| Contamination | 16 | High | Reverse osmosis device(s) need replacement | High water conductivity. Fouling effects of the water heater and steam generator | CMH01 | | Low | Moderate | |
| | 17 | High | Sulfur traps uptake capacity reached | Reforming Catalyst deactivation. Reduced hydrogen production. Contamination of WGS catalyst, PSA and fuel cell. | None | None | Low | Low | Use of a safety factor during DS-01 unit desing. Use of precious metals catalyst |
| | 18 | Low | None | | | | | | |

Source: Company information.

Equipment list

| Instrument List | | | | | |
|-----------------|----------------------|-----------------|-----------|--------------|-------|
| Displayed Text | Description | Connection Size | Service | Manufacturer | Model |
| MFC-1 | Mass Flow Controller | 1/4" | | -W | |
| MFC-2 | Mass Flow Controller | 1/4" | | -W | |
| TC-01 | TC | K-type | 0-1000 oC | co | |
| TC-02 | TC | K-type | 0-1000 oC | co | |
| TC-03 | TC | K-type | 0-1000 oC | co | |
| TC-04 | TC | K-type | 0-1000 oC | co | |
| TC-05 | TC | K-type | 0-1000 oC | co | |
| TC-06 | TC | K-type | 0-1000 oC | co | |
| TC-07 | TC | K-type | 0-1000 oC | co | |
| TC-08 | TC | K-type | 0-1000 oC | co | |
| TC-09 | TC | K-type | 0-1000 oC | co | |
| TC-10 | TC | K-type | 0-1000 oC | co | |
| TC-11 | TC | K-type | 0-1000 oC | co | |
| TC-12 | TC | K-type | 0-1000 oC | co | |
| PI-02 | Pressure transducer | 1/4" | 0-4 barg | IO | |
| LV-1 | Level meter | | On/Off | no | |

| Valve List | | | | | |
|----------------|--------------------------------|-----------|---------------|--------------|-------|
| Displayed Text | Description | Line Size | Valve Class | Manufacturer | Model |
| CV-1 | Check Valve | 1/4" | SS 304 | | |
| CV-2 | Check Valve | 1/4" | Spring loaded | | |
| V-1 | On/Off Valve (Normally Closed) | 1/4" | Brass | | |
| V-2 | On/Off Valve (Normally Closed) | 1/4" | Brass | | |
| V-3 | On/Off Valve (Normally Closed) | 1/4" | Brass | | |
| V-4 | On/Off Valve (Normally Closed) | 1/4" | Brass | | |
| V-5 | On/Off Valve (Normally Open) | 1/4" | Brass | | |
| V-6 | On/Off Valve (Normally Closed) | 1/4" | Brass | | |
| V-7 | On/Off Valve (Normally Open) | 1/2" | Brass | | |
| V-8 | On/Off Valve (Normally Open) | 1/2" | Brass | | |
| V-9 | On/Off Valve (Normally Open) | 1/2" | Brass | | |

| Equipment List | | | |
|-----------------|-------------------------------------|--------------|----------|
| Displayed Text | Description | Manufacturer | Material |
| BL-1 | Air Pump | | Aluminum |
| DI water tank | Distilled water tank | | Plastic |
| DS | Desulfurization | | SS 304 |
| E-1 | Steam Gen | | SS 304 |
| E-2 | Heat Exchanger (Recupurator) | | SS 304 |
| E-3 | Air cooled HEX | | SS 304 |
| E-5 (Condenser) | Heat Exchanger (Condenser) | | SS 304 |
| E-6 | Air Preheater | | SS 316 |
| EHL-1 | Heat Exchanger (External heat load) | | Aluminum |
| Preheater BL-2 | Air Pump | | Aluminum |
| REFORMER | Reformer | | SS 304 |
| RP-1 | Recirculation pump | | Plastic |
| SG-WGS | Reactor Vessel | | SS 304 |
| WP-1 | Water Pump | | Plastic |

Source: Company information.

Final equipment sizing

| Summary Table – SMR Equipment to Be Sized | |
|--------------------------------------------------|-------------------------------------------------------|
| Section | Equipment (valves, pressure meter, level meter, etc.) |
| Feed | Filters, desulfurizers |
| Reforming | SMR reactor |
| Shift | HTS reactor |
| Heat Recovery | Heat Exchangers |
| Separation | Knock Out drums |
| Purification | PSA |
| Product | Compressors, storage |
| Utilities | Water, fuel, nitrogen |
| Safety | PSV, flare |

Source: Company information.

Example of solenoid valve sizing

1. Process Data (Mandatory)

Fluid type (water, steam, gas, nitrogen, fuel gas)
 Flow rate (normal & max)
 Upstream pressure (P_1)
 Downstream pressure (P_2)
 $\Delta P = P_1 - P_2$
 Fluid temperature
 Fluid density / specific gravity
 Viscosity (important for liquids)

2. Valve Functional Requirements

Normally Open (NO) or Normally Closed (NC)
 Fail position (fail open / fail closed)
 On-off or modulating (solenoids are normally on-off)
 Opening time requirement (fast / standard)

3. Mechanical & Design Inputs

Line size
 Design pressure & temperature
 Valve body material
 Seal material (EPDM, Viton, PTFE, etc.)
 End connections (threaded, flanged)

4. Electrical Inputs

Voltage (e.g. 24 VDC, 110 VAC, 230 VAC)
 Area classification (safe / ATEX / IECEx)
 IP rating

Kv Equation for Liquid Flow

$$Q = K_v \sqrt{\frac{\Delta P}{SG}}$$

Where:

Q = liquid flow rate (m^3/h)

Kv = valve flow coefficient

ΔP = pressure drop across valve (bara)

SG = specific gravity (water = 1.0)

Rearranged (to size Kv):

$$K_v = \frac{Q}{\sqrt{\Delta P / SG}}$$

Boring section...😊

Economic and Environmental Considerations

No engineering design is complete without evaluating its economics and environmental impact. We consider these aspects for hydrogen via steam reforming of biomethane.

Why?

Environment: Regulations are becoming stringer and processes must comply otherwise can be rejected from permitting.

Projects need investment: We must justify it by either cost reduction or new revenue stream creation.

Economic considerations

- **Cost-effective** hydrogen production with mature, proven technology.
- **Lower** €/kg H₂ compared to electrolysis; competitive as blue hydrogen with CCS.
- High overall efficiency (**>85% in CHP setups**) through integrated heat recovery.
- Uses existing biomethane infrastructure, **reducing logistics costs**.
- PSA delivers high-purity H₂ (>99.97%), enabling **premium applications**.
- Catalyst and feed-gas quality influence OPEX; biomethane requires less sulfur removal devices.
- Modular design **reduces installation costs** and supports scalable deployment.
- **Low NO_x emissions** with catalytic combustion lower environmental compliance costs.

Environmental Impact and Sustainability

- **Low-carbon hydrogen pathway:** Biogenic methane originates from recent biomass, significantly reducing lifecycle greenhouse gas emissions compared to fossil SMR.
- **Potential for net-negative emissions:** When combined with carbon capture and storage (CCS), SMR on biogenic methane can achieve near-zero or carbon-negative CO₂ emissions.
- **Circular-economy benefits:** Utilizes renewable, waste-derived feedstocks (e.g. biogas, landfill gas), reducing methane emissions from waste streams.
- **Controlled local environmental impact:** Comparable air emissions to conventional SMR; NO_x and sulfur emissions are effectively mitigated through proven design and controls.
- **High efficiency and scalability:** Leverages mature SMR technology with extensive heat recovery and compatibility with existing infrastructure.
- **Feedstock sustainability dependent:** Environmental performance depends on certified, responsibly sourced biogenic methane, with waste-based sources offering the highest benefit.

Conclusions



THANK YOU FOR YOUR ATTENTION

Questions...?

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Conclusions

- Process design is rigorous process that pertains to selection, sequencing and specifications of the units (equipment) required for the conversion of raw materials into finished products.
- Process design is considered the "pinnacle" of Chemical Engineering, as it combines all the individual knowledge (thermodynamics, kinetics, mass and energy transfer) to create a functional industrial unit.
- Chemical engineering design software are powerful tools for process design, but we need to pay extra attention for delivering concrete results.
- PFD (Process Flow Diagram) and P&ID (Piping & Instrumentation Diagram) are the two most basic tools in the design and operation of industrial facilities, functioning complementary at different stages of a project.
- The HAZOP (Hazard and Operability Study) is a systematic and structured risk assessment technique, which aims to identify potential safety and operability problems in an industrial process.
- Environmental and economic considerations of the process determines the “go or no-go” decision for a process/project.