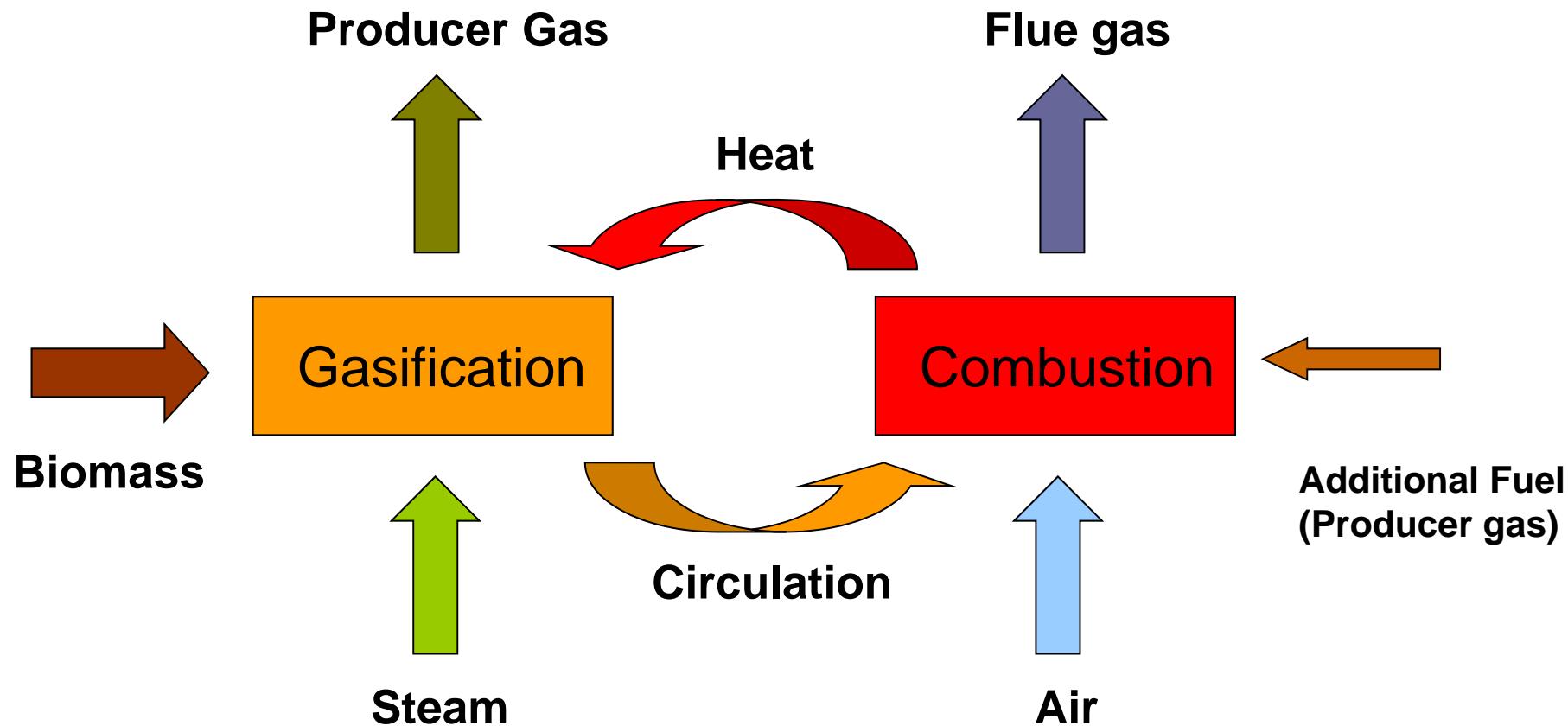


# **Gas Sampling, Measurement and Analyses at Güssing - DFB-Biomass gasification**

Dr. Reinhard Rauch

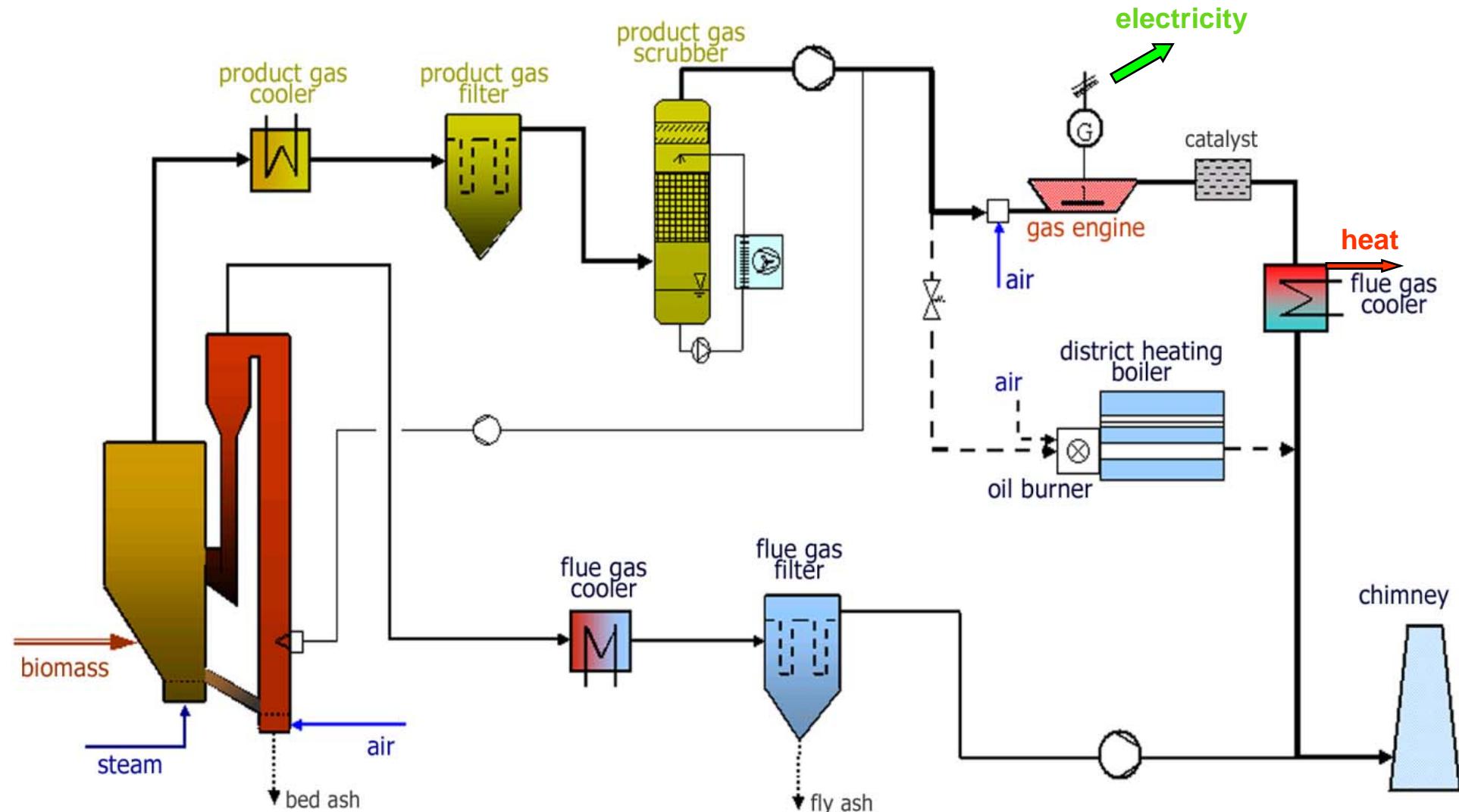
Vienna, University of Technology  
Bioenergy 2020+

# Gasification Concept of Dual Fluid



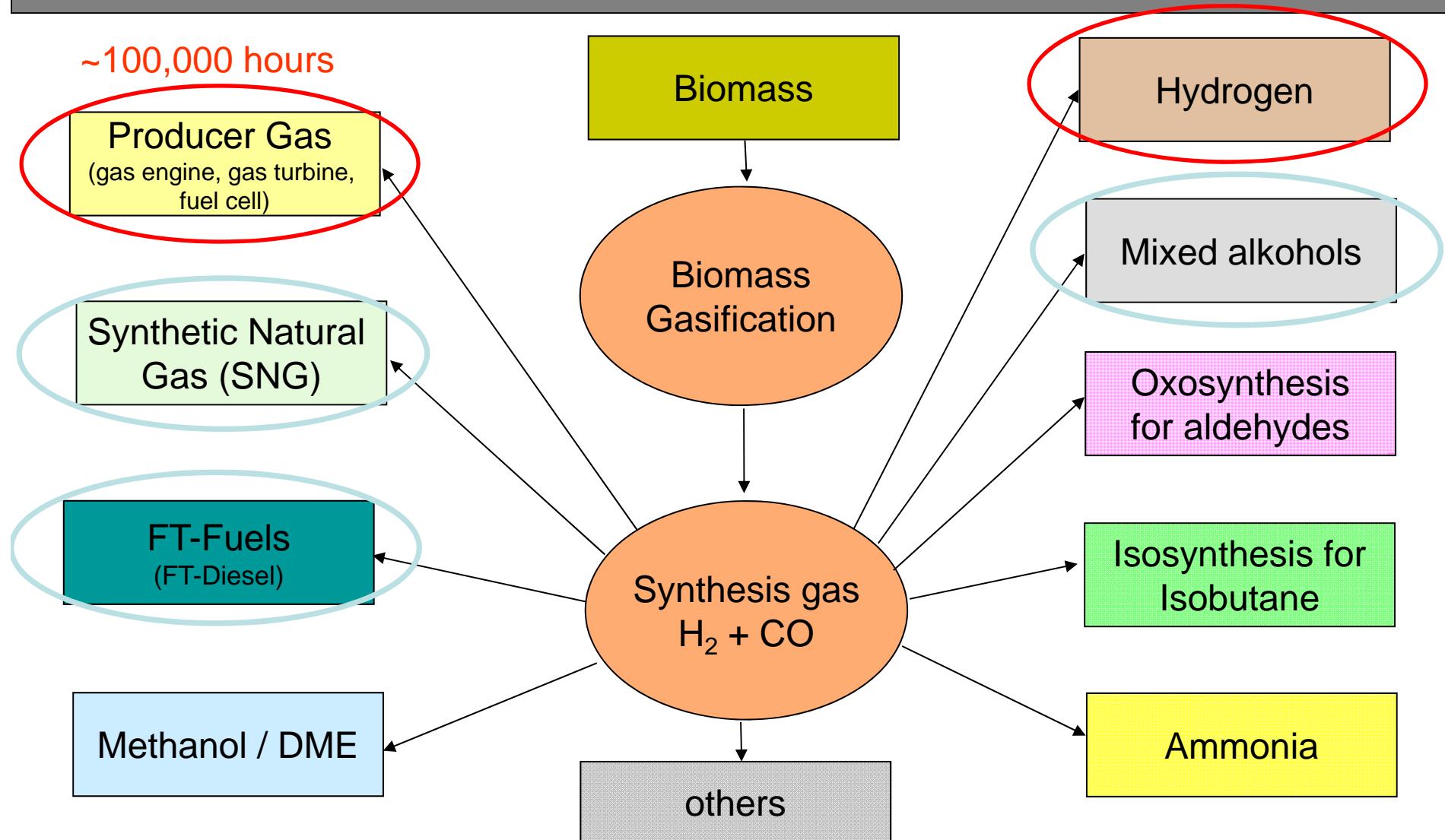
- Low temperature gasification (800-900°C)
- Atmospheric pressure
- Catalytic activity of bed material and ash influences the gas quality

# Flow chart of DFB for CHP bioenergy2020+



# The basic concept – “Green Chemistry”

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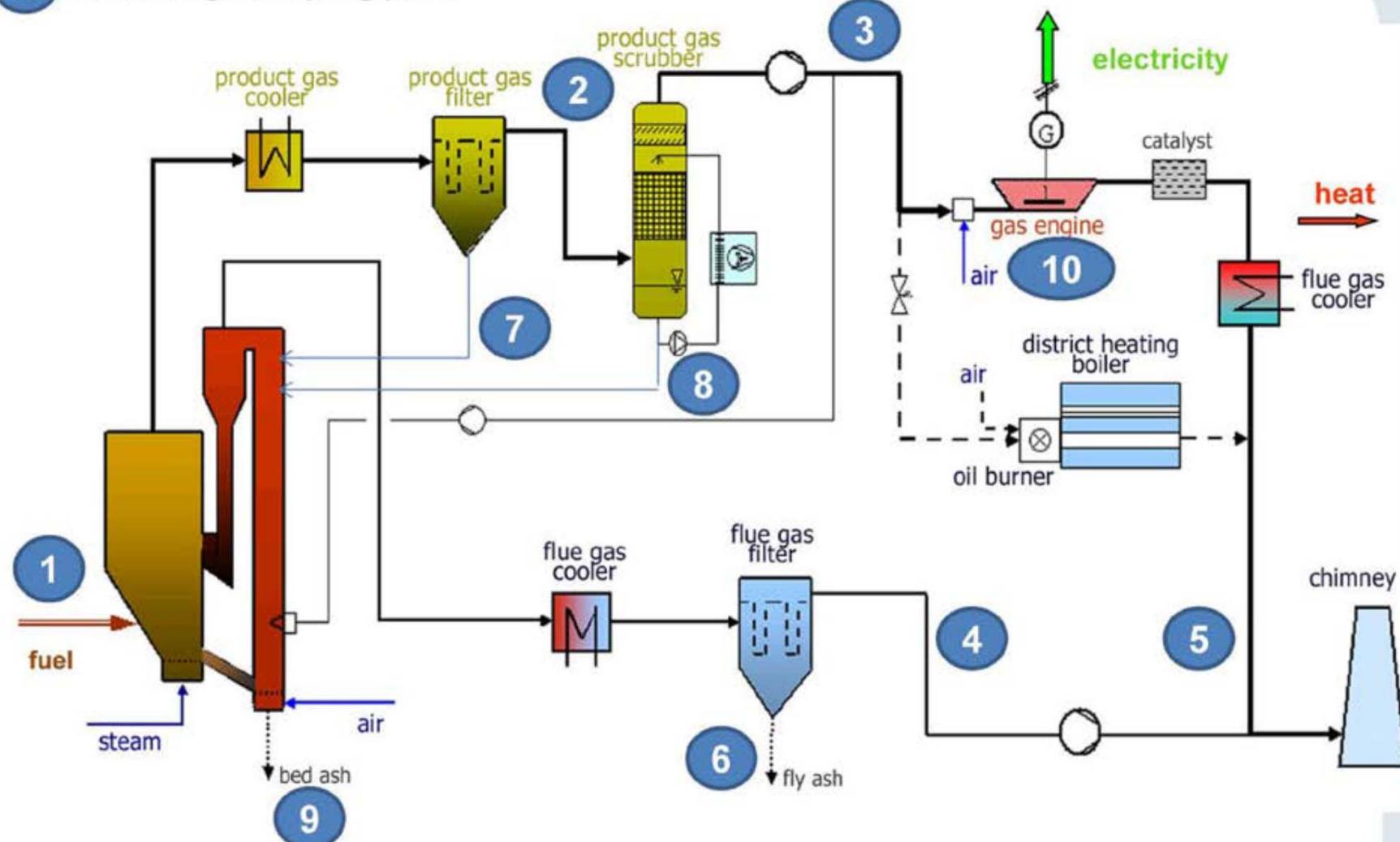


- Here fast results are necessary, so only online systems are used
- For normal operation only gas components are analysed (CO, CO<sub>2</sub>, O<sub>2</sub>), here typical online analysers are used (infrared, paramagnetic)
- Main difficulty is to clean the gas and to bring it into the analyser with good availability
- Analysis of the engine oil is also quite important as it shows the quality of the gas treatment
  
- For commissioning/optimisation of the process the following analysis are done:
  - All permanent gases in the producer gas (by GC)
  - Flue gas in combustor and gas engine (CO, NOx, SOx, PAH, ...)
  - Tar and particles in gas phase (gravimetric and GC-MS), here TU Vienna uses the standard method, only a different solvent (toluene) is used
  - Hydrocarbons / tar content in solvents (e.g. tar content in RME)
  - Bed material and ashes (activation of the bed material)

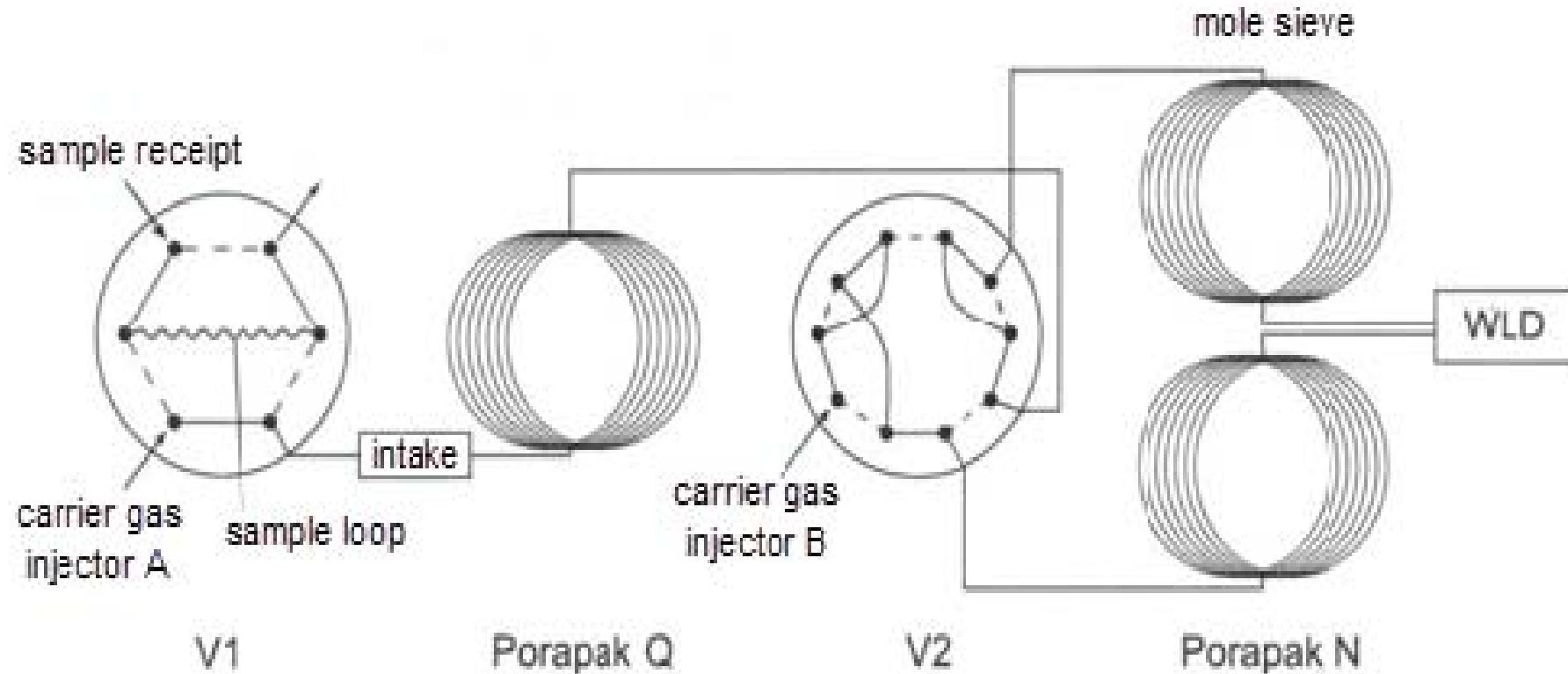
# Typical analytical points for full characterisation

n

Measuring / sampling point



# Typical GC configuration



# Gas Composition (after gasifier)

Main Components		
H <sub>2</sub>	% (dry)	35-45
CO	% (dry)	22-25
CH <sub>4</sub>	% (dry)	~10
CO <sub>2</sub>	% (dry)	20-25
H <sub>2</sub> O	%	30-50

Minor Components		
C <sub>2</sub> H <sub>4</sub>	% (dry)	2-3
C <sub>2</sub> H <sub>6</sub>	% (dry)	~0.5
C <sub>3</sub> H <sub>6</sub>	% (dry)	~0.4
O <sub>2</sub>	% (dry)	< 0,1
N <sub>2</sub>	% (dry)	0.3-3
C <sub>6</sub> H <sub>6</sub>	g/m <sup>3</sup>	5-10
C <sub>7</sub> H <sub>8</sub>	g/m <sup>3</sup>	1-2
C <sub>10</sub> H <sub>8</sub>	g/m <sup>3</sup>	~2
TARS (grav.)	g/m <sup>3</sup>	1-5

Possible poisons		
H <sub>2</sub> S	mgS/Nm <sup>3</sup>	~200
COS	mgS/Nm <sup>3</sup>	~10
Mercaptans	mgS/Nm <sup>3</sup>	~30
Thiophens	mgS/Nm <sup>3</sup>	~7
HCl	ppm	~3
NH <sub>3</sub>	ppm	500-1000
Dust	g/Nm <sup>3</sup>	10-100

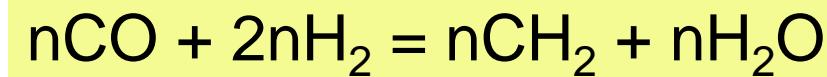
H<sub>2</sub>:CO = from 1.5:1 to 2:1

# Tar Composition (after gasifier)

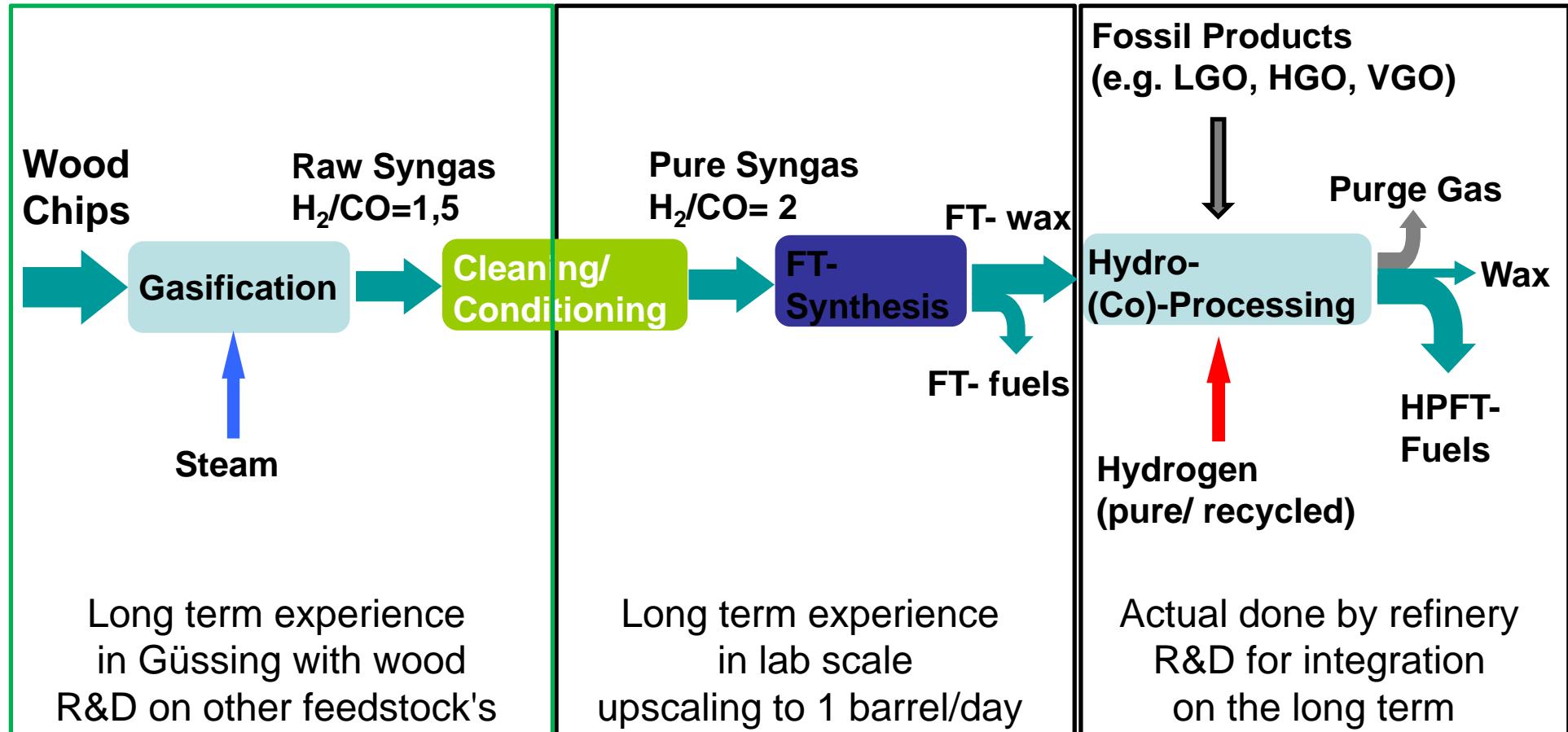
Phenylacetylen	mg/Nm <sup>3</sup>	122	141	Indol	mg/Nm <sup>3</sup>	58	112
Styrol	mg/Nm <sup>3</sup>	1360	1575	Biphenyl	mg/Nm <sup>3</sup>	222	290
Mesitylen	mg/Nm <sup>3</sup>	0	0	Isoeugenol	mg/Nm <sup>3</sup>	0	0
Phenol	mg/Nm <sup>3</sup>	345	515	Acenaphtylen	mg/Nm <sup>3</sup>	1769	2268
Benzofuran	mg/Nm <sup>3</sup>	243	301	Acenaphthen	mg/Nm <sup>3</sup>	56	70
1H-Inden	mg/Nm <sup>3</sup>	2621	2243	Dibenzofuran	mg/Nm <sup>3</sup>	145	176
2-Methylphenol	mg/Nm <sup>3</sup>	0	0	Flouren	mg/Nm <sup>3</sup>	435	633
4-Methylphenol	mg/Nm <sup>3</sup>	0	0	Dibenzothiophen	mg/Nm <sup>3</sup>	0	0
2-Methylbenzofuran	mg/Nm <sup>3</sup>	0	0	Anthracen	mg/Nm <sup>3</sup>	1209	1342
2,6-Dimethylphenol	mg/Nm <sup>3</sup>	0	0	Phenanthren	mg/Nm <sup>3</sup>	271	330
2,5 u. 2,4-				Carbazol	mg/Nm <sup>3</sup>	0	0
Dimethylphenol	mg/Nm <sup>3</sup>	0	0	4,5-Methylphenanthren	mg/Nm <sup>3</sup>	111	161
3,5-Dimethylphenol	mg/Nm <sup>3</sup>	0	0	9-Methylanthracen	mg/Nm <sup>3</sup>	0	0
2,3-Dimethylphenol	mg/Nm <sup>3</sup>	0	0	Flouranthen	mg/Nm <sup>3</sup>	284	329
3,4-Dimethylphenol	mg/Nm <sup>3</sup>	0	0	Pyren	mg/Nm <sup>3</sup>	301	358
2-Methoxy-4-				Benzo[a]anthracen	mg/Nm <sup>3</sup>	112	119
Methylphenol	mg/Nm <sup>3</sup>	0	0	Chrysen	mg/Nm <sup>3</sup>	180	199
Naphtalin	mg/Nm <sup>3</sup>	5954	4892	Benzo[b]flouranthen	mg/Nm <sup>3</sup>	0	31
1-Benzothiophen	mg/Nm <sup>3</sup>	0	0	Benzo[k]flouranthen	mg/Nm <sup>3</sup>	0	38
Chinolin	mg/Nm <sup>3</sup>	62	103	Benzo[a]pyren	mg/Nm <sup>3</sup>	8	84
2-Methylnaphthalin	mg/Nm <sup>3</sup>	457	534	Benzo[g,h,i]perylene	mg/Nm <sup>3</sup>	0	0
Isochinolin	mg/Nm <sup>3</sup>	0	0	Dibenz[a,h]anthracen	mg/Nm <sup>3</sup>	0	0
1-Methylnaphthalin	mg/Nm <sup>3</sup>	301	355	Indeno[1,2,3-cd]pyren	mg/Nm <sup>3</sup>	0	0
1-Indanon	mg/Nm <sup>3</sup>	0	0				
Eugenol	mg/Nm <sup>3</sup>	0	0				

Analytics in synthesis applications,  
as example FT is used:

*BioFiT*  
BIOMASS-TO-FISCHER-TROPSCH



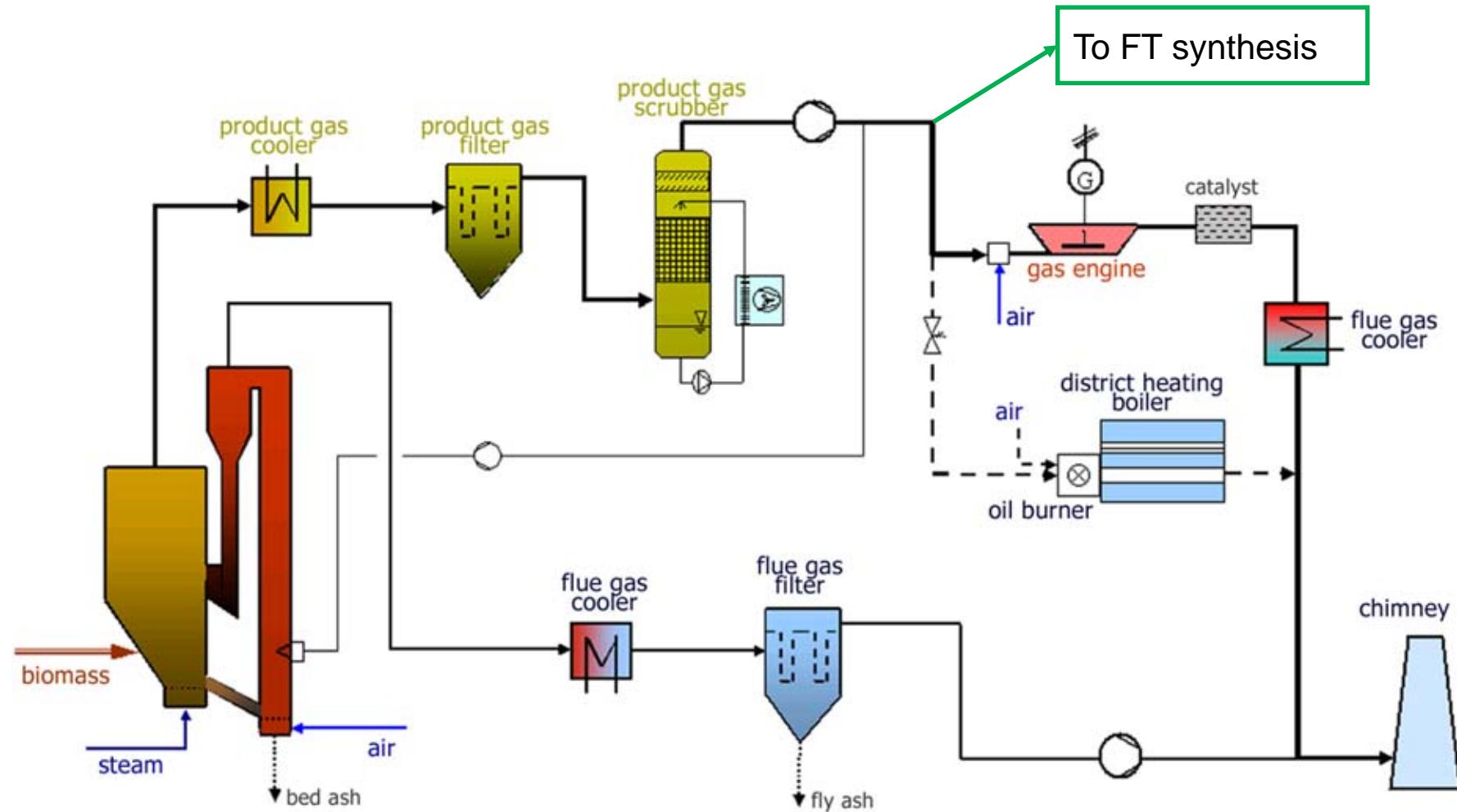
# Schema of FT

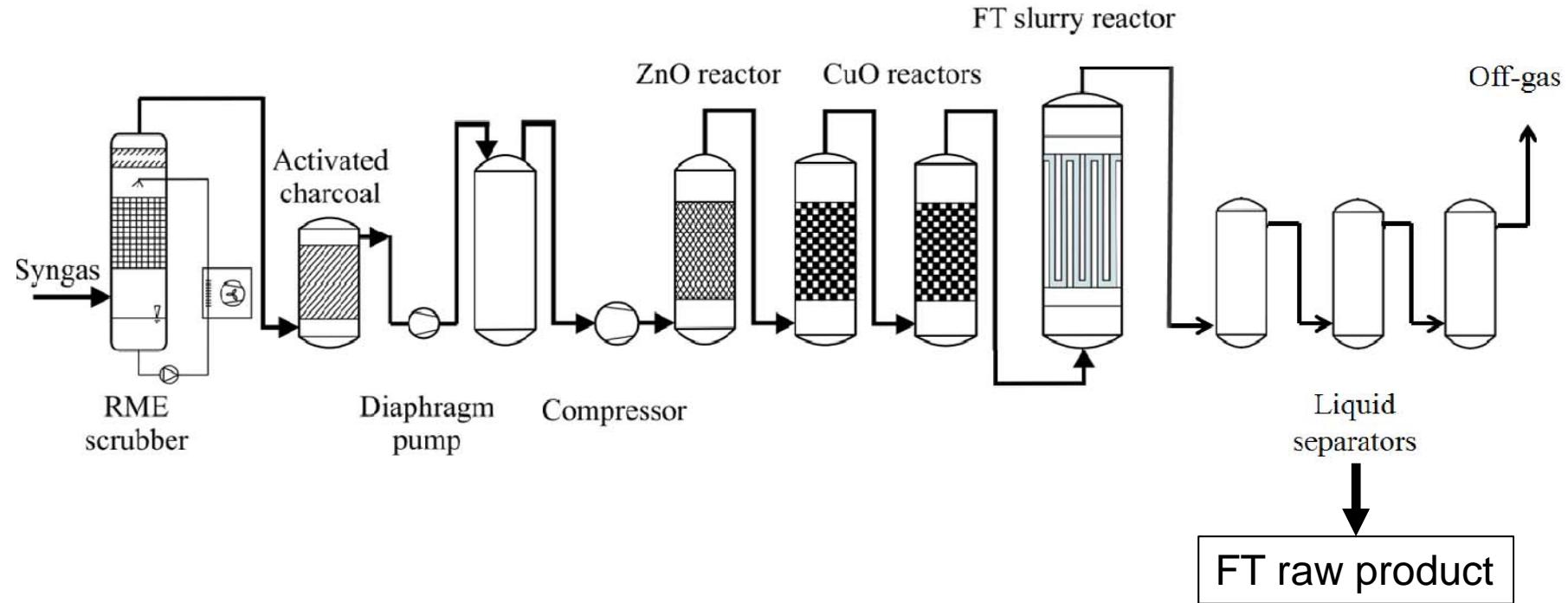


Same analytics as in CHP are used

Focus on catalyst poisons, like S, Cl or metals

Technology and analytics from refineries are used



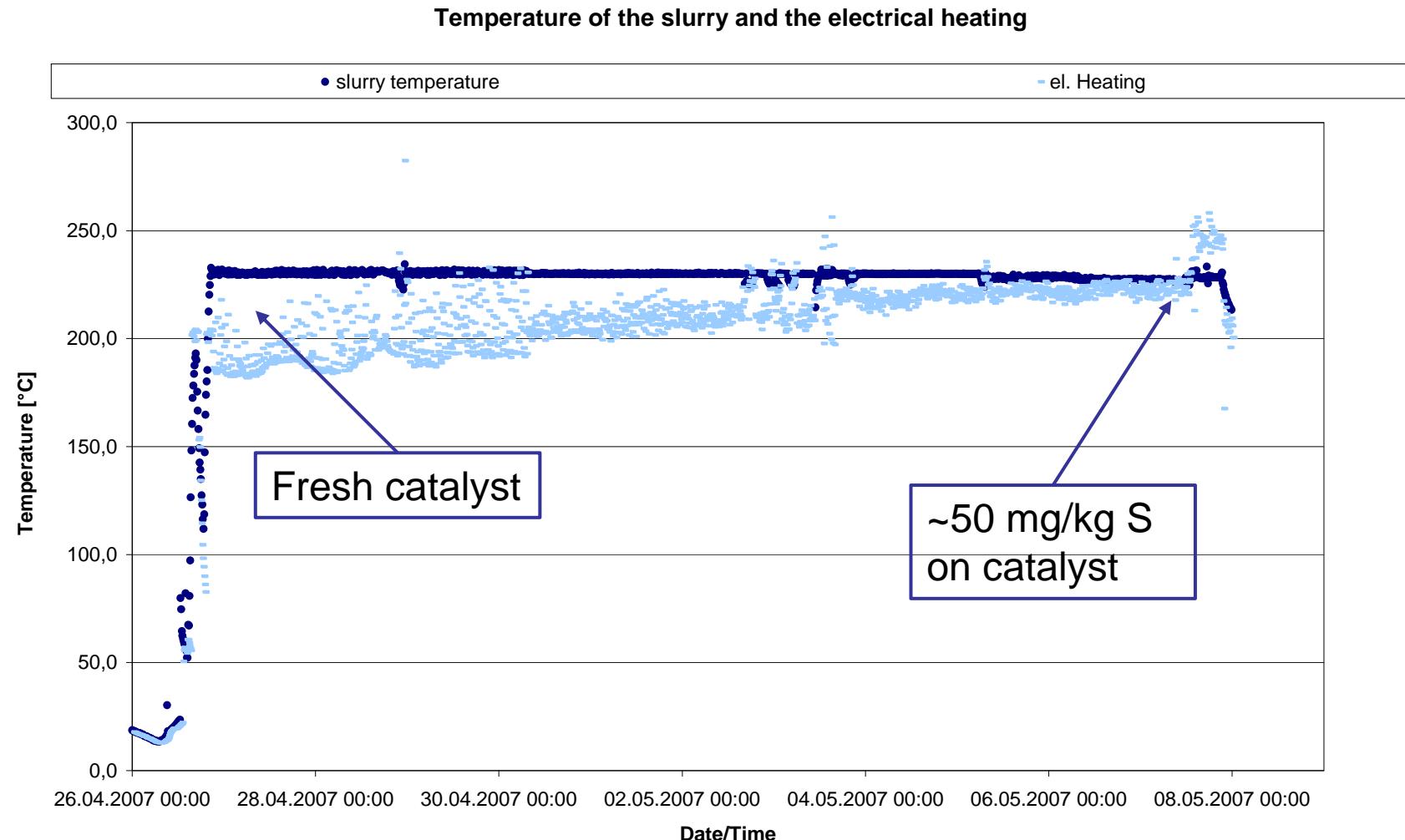


5-10kg/day of FT raw product

Slurry reactor, because of excellent heat transfer and easy scaling up

Gas treatment removes Sulphur to below 10ppb

Fully automatic



- Sulphur
  - In the gas phase by GC-SCD,
  - sampling is done by Tedlar bags
- Chlorine, Ammonia, HCN
  - Sampling in liquids (e.g. H<sub>2</sub>SO<sub>4</sub>),
  - analysis by IC
- BTX, Naphthalene
  - Sampling by Tedlar bags or online
  - Analysed by GC
- Metals, Carbonyls
  - Done by external organisations

# Conclusion

- Analytics for standard CHP are established and work well
- They are only too expansive and/or to much maintenance is necessary
- Species analysed for synthesis gas are different than CHP (focus more on inorganic)
- Analytics for synthesis applications are more difficult, as the range from **ppm** is changed to **ppb**
- In addition to the analytics long term tests in pilot scale are really necessary, because you do not know, if you miss one catalyst poison

## More Information

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More info at

<http://www.ieatask33.org>

<http://www.ficfb.at>

<http://www.vt.tuwien.ac.at>

<http://www.bioenergy2020.eu>