



# Introduction to OGT, Inc.

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CAPE-OPEN Annual Meeting  
BP Offices, Sunbury on Thames, England  
October 2017





# Company Overview

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- Dr. Ralph Weiland – President
  - 30 years Chemical Engineering Professor / Researcher
  - 10 years in Mass Transfer with Koch-Glitsch
    - Lead Inventor of ULTRA-FRAC® trays
- Business Office
  - Located in Houston, Tx
  - Lead by Michael Jakobs, Business Director
- Technical Development Office
  - Located near Austin, Tx
  - Lead by Nate Hatcher, Technical Development Lead





# ProTreat®

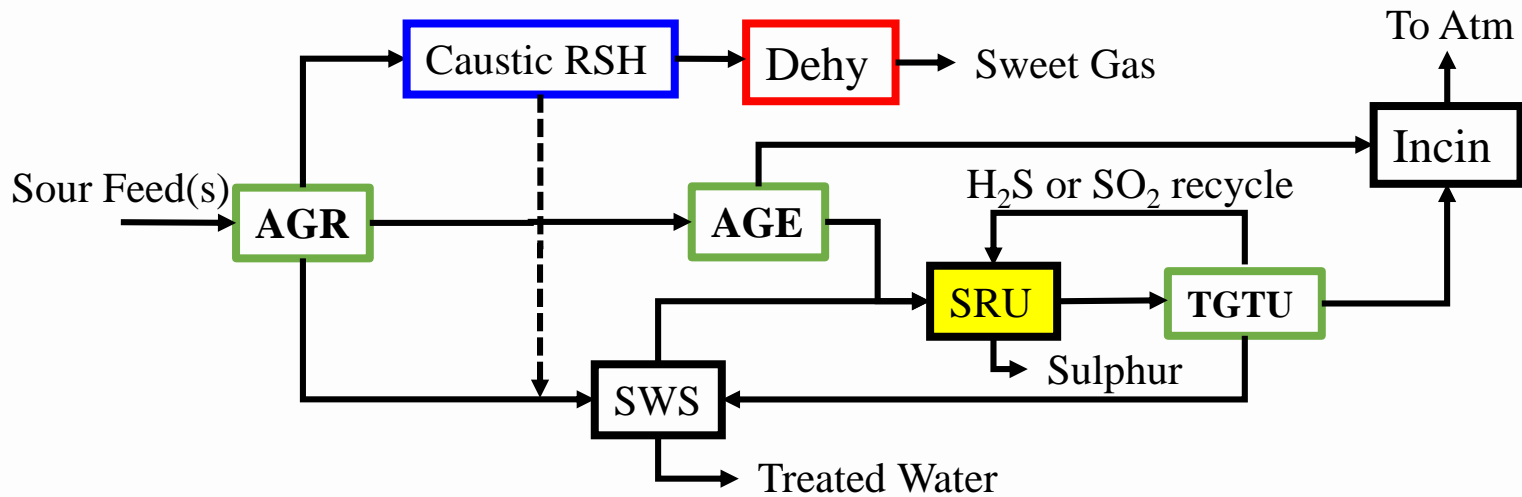
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- Simulation tool for:
  - Amine Treating
  - Physical Solvent Treating
  - Carbon Capture
  - Sour Water Strippers
  - Caustic Treating
  - Sulphur Recovery
  
- Fully flexible flowsheeting tool





# What can ProTreat Simulate?





# ProTreat<sup>®</sup> features

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- Rigorous electrolyte aqueous thermodynamics
- Many solvents available
  - Generic and licensed amines
  - Blend up to three amines
  - Physical solvents and Amino acids
  - NaOH, KOH, K<sub>2</sub>CO<sub>3</sub>
- Gas components
  - Acid gases including SO<sub>2</sub>, HCN
  - Ammonia
  - RSH, Phenol, Methanol
- Common Heat Stable Salt ions / Stripping Promoters
- Hydrocarbon / BTEX solubility in amines and physical solvents



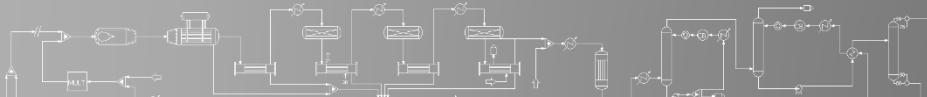


# ProTreat<sup>®</sup> features

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## Rigorous Sulphur Recovery Model

- Unique aspects of sulphur vapor / molten sulphur
  - Variable Mole Weight
  - H<sub>2</sub>S solubility
  - Viscosity / Heat Capacity inflections
- Reaction Furnace with kinetic ammonia / HC destruction
- WHB with rigorous heat transfer, chemical reactions
- Sulphur condenser
- Converter





# Our Customers – Operators / Suppliers





# Our Customers – Engineering / Consulting







# Our Customers – Engineering / Consulting



A Navratna Company



TOYO  
ENGINEERING





# How is the ProTreat Column Model Different?

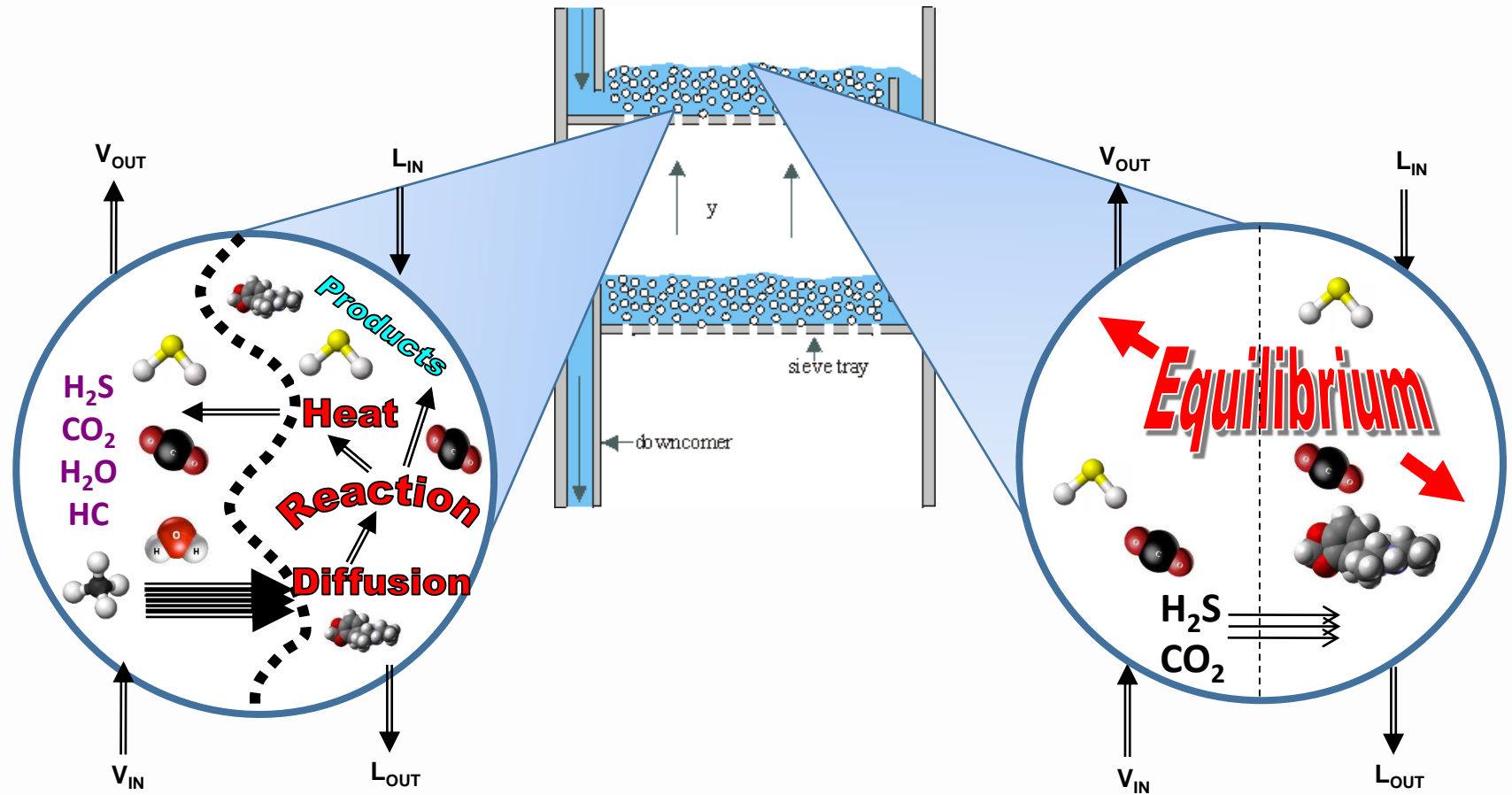
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- Column calculations are based on mass transfer and heat transfer rates
- Directly calculates tray-by-tray performance
- Predicts actual mass transfer
  - ProTreat does not calculate equilibrium then offset it by a user-supplied factor
  - No translation from ideal stages to real tower internals
  - No tray efficiency; no HETP
- User is not required to provide tuning factors
- Real trays, random packing, structured packing — hydraulics and mass transfer





# Mass Transfer Rate Model vs Ideal-Stage Model



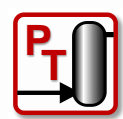


# ProTreat's COM Interface

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- Currently, a proprietary COM interface structure
- Allows 3<sup>rd</sup> party program to
  - Set input values
  - Execute the program to generate results
  - Retrieve calculated result values
- Current status
  - ProTreat flowsheet block available in VMG-Sim
  - Spreadsheet unit monitoring / flowsheet calc templates using VBA





# CAPE-OPEN plan

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- Use COM as the underlying technology to implement the CAPE-OPEN interface
- Implementation priorities
  - Unit Operations module, e.g. allow ProTreat columns to be integrated with other simulators
  - Thermo & Physical Props module, e.g. allow HTRI to use ProTreat for stream calculations within a heat exchanger
- Goal
  - Enable both Unit Operations and Thermo / Physical Props over the next year

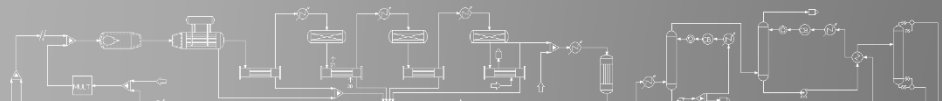




# Our next steps

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- Digest CAPE-OPEN documentation
- Begin implementing Unit Operations module
  
- Question:
- What technical support is available from CO-LaN?
  - Consultancy
  - Tech Support Single-Point-Of-Contact



The logo icon consists of a red square border containing a stylized grey vertical cylinder representing a gas treatment vessel. To the left of the cylinder, the letters 'P' and 'T' are stacked vertically in red. A black arrow points from the left into the cylinder.

# ProTreat<sup>®</sup>

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*The industry's most powerful gas treating and sulphur recovery simulation tool*

