

## Changes with *OGT Suite* Version 8.0

### ProTreat® & SulphurPro®

Release date: Oct 2022

#### 1. **Thermal Oxidizer and Stack**

- a. Two new SulphurPro blocks to model sulphur plant incineration
- b. Rigorous kinetics model for the destruction of reduced sulphur species, carbon monoxide, hydrogen and ammonia
- c. Rigorously model both natural draft and forced draft flowsheet configurations
- d. Calculate pressure drop through the vessels
- e. Stack hydraulic calculations with an insufficient stack draft warning.
- f. Component-wise emissions reporting in Stack

#### 2. **Sulphur Storage Block**

- a. New SulphurPro block to model sulphur storage and degassing
- b. Rigorous kinetics for H<sub>2</sub>S degassing from liquid sulphur
- c. Calculate H<sub>2</sub>S evolution with a specified sweep gas inlet
- d. Optional sparge gas inlet for modeling degassing
- e. Options for calculating Catalytic degassing include Morpholine, Ammonium Sulphite, Aromatic Amine, and Urea
- f. Account for oxidation of evolved H<sub>2</sub>S to SO<sub>2</sub> in the headspace
- g. Represent various configurations such as in-pit degasser, sulphur pit, above ground sulphur storage tank, etc.

#### 3. **CAPE-OPEN Property Package**

- a. ProTreat® is now a CAPE-OPEN compliant property package plug
- b. Create property packages from ProTreat with desired compounds and thermodynamic method
- c. Use within any CAPE-OPEN compliant thermodynamic socket

#### 4. **Hybrid Solvents Properties Update**

- a. Incorporated several new data sets for physical and transport properties for hybrid solvent mixtures of sulfolane and amines
- b. Better agreement with plant data

#### 5. **DIPA Heat Capacity Refit** – Refit of DIPA heat capacity to give better agreement with field data

## Changes With *ProTreat*® Version 7.0

Release date: May 2022

1. **New Physical Solvent – Sulfolane**
2. **New Thermodynamic Package – Hybrid Solvents**
  - a. Rigorously model mixtures of amines and physical solvents
  - b. Accurate VLE, physical properties and mass transfer characteristics for **blends of water, amines and sulfolane**
3. **Major COS Model Update**
  - a. Accurately predict the chemical absorption of COS in columns
  - b. Rigorous reaction chemistry and kinetics model
4. **DIPA Model Update**
  - a. Complete refit of DIPA VLE and physical properties
  - b. Incorporated latest published data
  - c. Improved mercaptans predictions
5. **New DOW Solvents**
  - a. UCARSOL™ HS-101
  - b. UCARSOL™ HS-102
  - c. UCARSOL™ HS-115
  - d. UCARSOL™ AP-802
  - e. UCARSOL™ LE-713
  - f. UCARSOL™ LE-777
  - g. UCARSOL™ LE-801
6. **New Hydrogenation Reactor Block**
  - a. **First-Gen Cobalt-Molybdenum Catalyst**
  - b. Rigorous reaction kinetic model
  - c. Hydrothermal ageing model
  - d. Catalyst poisoning model
  - e. Layer catalyst and inert sections in the same bed
  - f. Calculate pressure drop across the catalyst bed
  - g. Advanced vessel rating and sizing options
7. **Sulphur Converter Improvements**
  - a. New kinetic model for **Titania Catalyst**
  - b. Completely reworked interface
  - c. Layer Alumina and Titania catalyst along with inert support layers in the same bed
  - d. Calculate pressure drop across the catalyst bed
  - e. Advanced vessel rating and sizing options

8. **Burner Block Improvements**

- a. Advanced reaction control of H<sub>2</sub>S thermal splitting and partial oxidation to model **Acid Gas Fired Reheaters**
- b. Effluent temperature spec
- c. Solver variables implemented

9. **New Sulphur Conversion and Recovery Tables**

- a. Flexible calculation and display of Sulphur conversion and/or recovery on the Flowsheet
- b. User ability to create and customize any number of calculation 'stages'

10. **Callouts Improvements**

- a. Create and manage templates from new or existing callouts
- b. Change default callout configuration
- c. Add callouts by right-clicking on flowsheet streams

11. **Duplicator Block Reference Stream feature** – Seamless duplication of any flowsheet stream including internal streams

12. **Reworked Convergence Dashboard**

- a. Completely reworked flowsheet convergence dashboard to easily monitor and troubleshoot convergence
- b. Minimize, close, dock or float for convenience

13. **Docking/Floating capabilities** for the Flowsheet Window, Stream Table and Convergence Dashboard

14. **Graphics Update** – High resolution for flowsheet drawings and other windows

15. **Usability Improvements**

- a. Cntrl+Scroll to zoom flowsheet
- b. Hold Space key to multi-insert a selected palette block
- c. Drag to reorder components in the components manager
- d. Drag to reorder streams in the stream table
- e. File Auto-backup/Recovery system
- f. Unit manager changes in the PTR mode gets transferred to the PTD mode

16. Significant convergence speed improvement

# Changes With *ProTreat*<sup>®</sup> Version 6.6

Release date: May 2021

1. **Improved Piperazine-CO<sub>2</sub> kinetics**
  - a. Re-assessment of Piperazine-CO<sub>2</sub> reaction kinetics leads to faster predicted CO<sub>2</sub> pick up in mass transfer rate limited columns
2. **Data Call Outs for Streams**
  - a. Fully customizable to display any value in any units
  - b. This block can be added and specified in either the PTD (input file) or PTR (results file). If specified in the PTR, that information is carried back to the PTD.
  - c. Callouts appear in flow sheet exports and Custom Excel Reports
  - d. Values are automatically updated with each run
  - e. Can be placed anywhere on the flowsheet or minimized to save space
3. **Two new flowsheet blocks**
  - a. **Reheater** – Heat transfer rate-based block for sulfur plant simulation
    - i. Includes sizing, rating, and simple energy balance modes
    - ii. Rigorous heat transfer and pressure drop calculations
    - iii. Auto-generated utility streams
  - b. **Burner** – Simulates combustion operations
    - i. Auto-generated air and tempering steam streams
    - ii. Emission factors for estimating emissions
4. **Calculated Pressure Drop** feature available for most blocks
  - a. Pressure drop through blocks responds to changes in flow rate and density
  - b. **Import From Results** button makes data entry easier: First run the block with its design pressure drop, then on subsequent runs the pressure drop will scale up or down from that fixed point.
5. Improvements to **Column block**
  - a. **Liquid inventory** reported including trays, packing, and (optional) sump
  - b. **Downcomer backup** reported for trays if downcomer clearance is provided
  - c. More variables available to plot
  - d. Methods can be chosen for hydraulics and mass transfer in packed sections. Mass transfer methods include equilibrium stage.
  - e. Improved convergence stability
  - f. New **CO<sub>2</sub> Kinetics Adjustment Factor for Reactive Amine**
6. **Automatically renumber all streams** with a new command in the Edit menu
7. **New driverless HASP keys** available
8. Other improvements
  - a. New flowsheet components: C<sub>5</sub> – C<sub>8</sub> alkenes
  - b. Several **new units** added to Unit Manager
  - c. Several minor improvements to flowsheet drawing and stream auto-routing
  - d. Case Study can now manipulate amine concentration in Control Block
  - e. Gibbs Reactor includes fugacity effects at high pressure and more accurate flame temperatures
  - f. Updated VLE for butyl mercaptan in MDEA
  - g. Improved model for ICE3137 solvent

- h. New stream warning for salts of ammonia and CO<sub>2</sub> precipitating from aqueous liquids
- i. Minor bug fix for Sulfur Converter block related to catalyst loading spec