

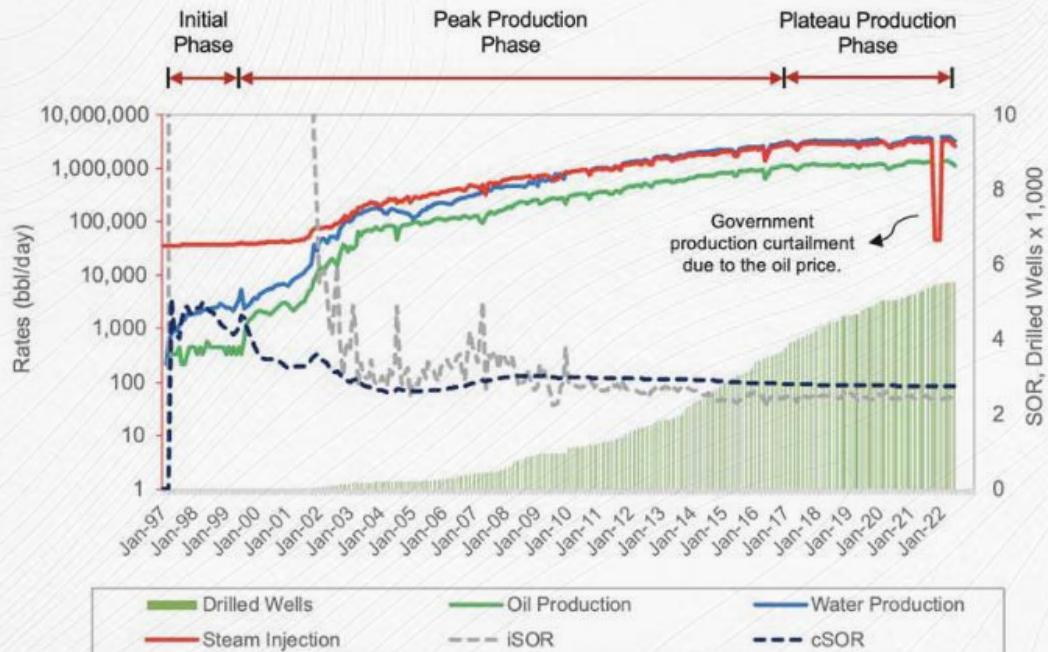
# 基于溶剂的稠油提高采收率模拟技术

## Enabling the Future of Solvent-Based EOR

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# Current Status of SAGD in Canada

Comprises 25% of the Canadian oil production – 1.44 million bbl/day



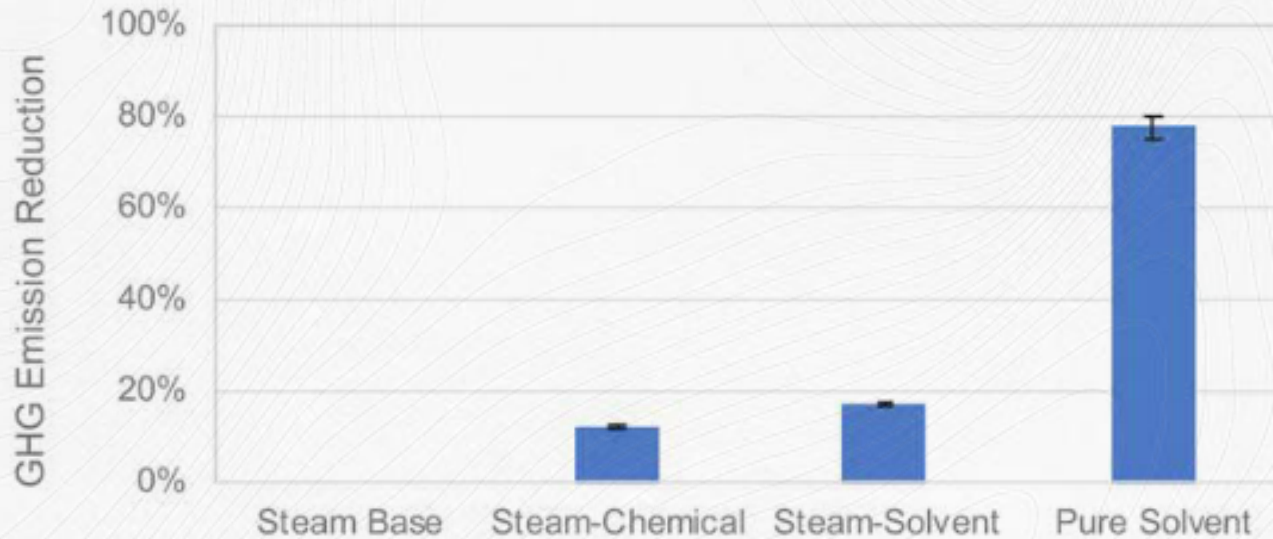
30 years of SAGD performance

22 SAGD projects  
300 pads and 2,700 well pairs



Active SAGD projects distribution\*

# Great Potential for Thermal Technologies



**SAGD/post-SAGD improvements in reducing GHG emissions  
(adopted from Canadian Energy Research Institute 2020)**

## Steam - Chemical:

- Alkalis
- Surfactants
- Ionic liquids
- Solvents
  - dimethyl ether
  - diethyl ether
  - additives
- Nanoparticles
- Chelating agents
- Deep eutectic solvents
- Biodiesel
- Switchable-hydrophilicity solvent

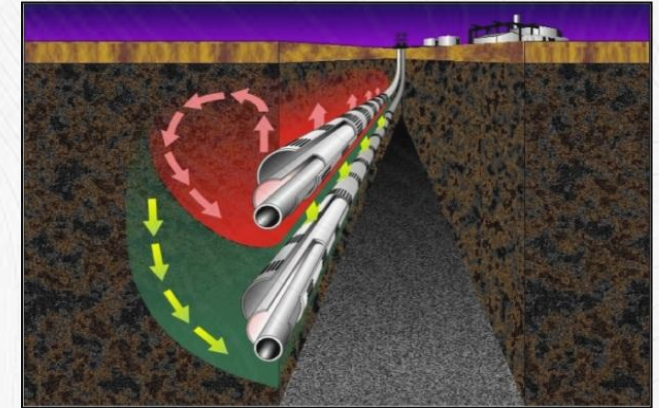
## Solvents

- NCG (methane)
- CG (propane)
- Liquid (C5+)

# Steam Processes

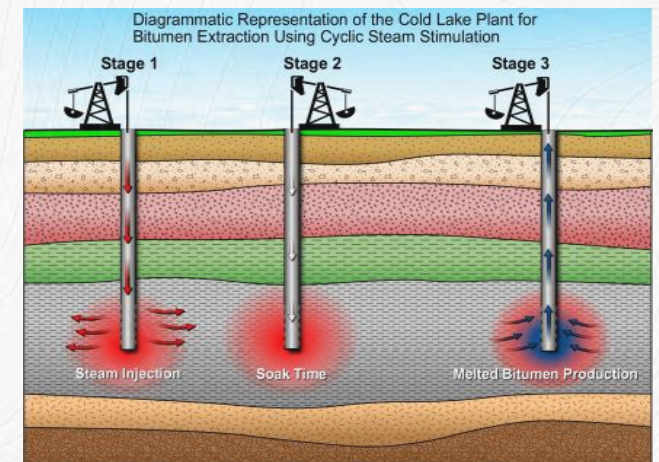
- High temperature steam
- Heavy oil viscosity reduction by temperature
- Heat loss to the surrounding formations
- Up to 70% recovery
- High operating cost

## Steam Assisted Gravity Drainage (SAGD)



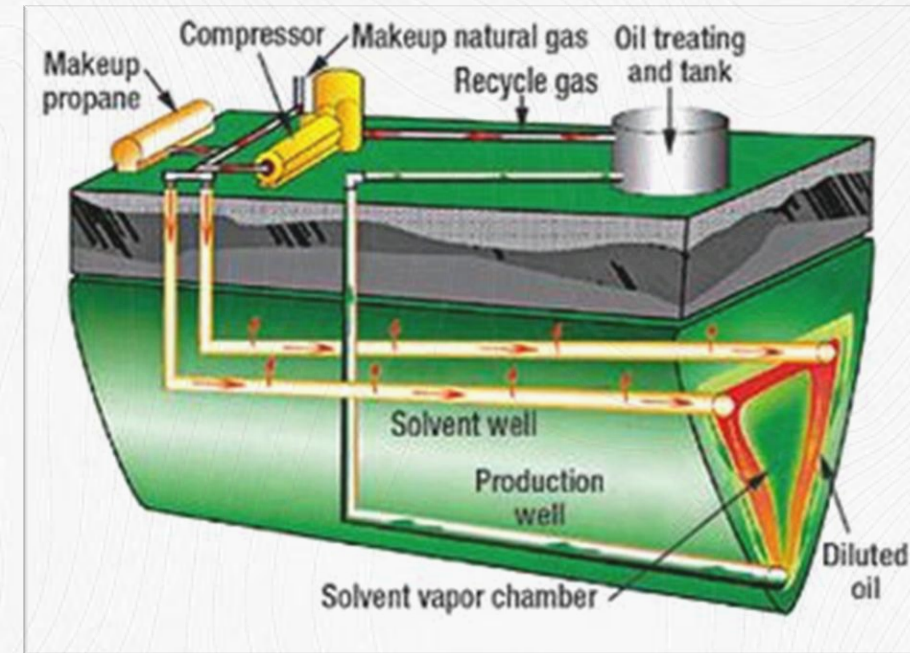
[www.oilsands.ualberta.ca](http://www.oilsands.ualberta.ca)

## Cyclic Steam Stimulation CSS



# Solvent Processes

- Solvent dissolution reduces oil viscosity
- Solvents in vapor phase decrease trapping
- Solvents mostly stay with the oil
- Diffusion and dispersion of solvent control the dilution and mixing process
- Mass-transfer is often slower than heat transfer



<http://bittooth.blogspot.ca>

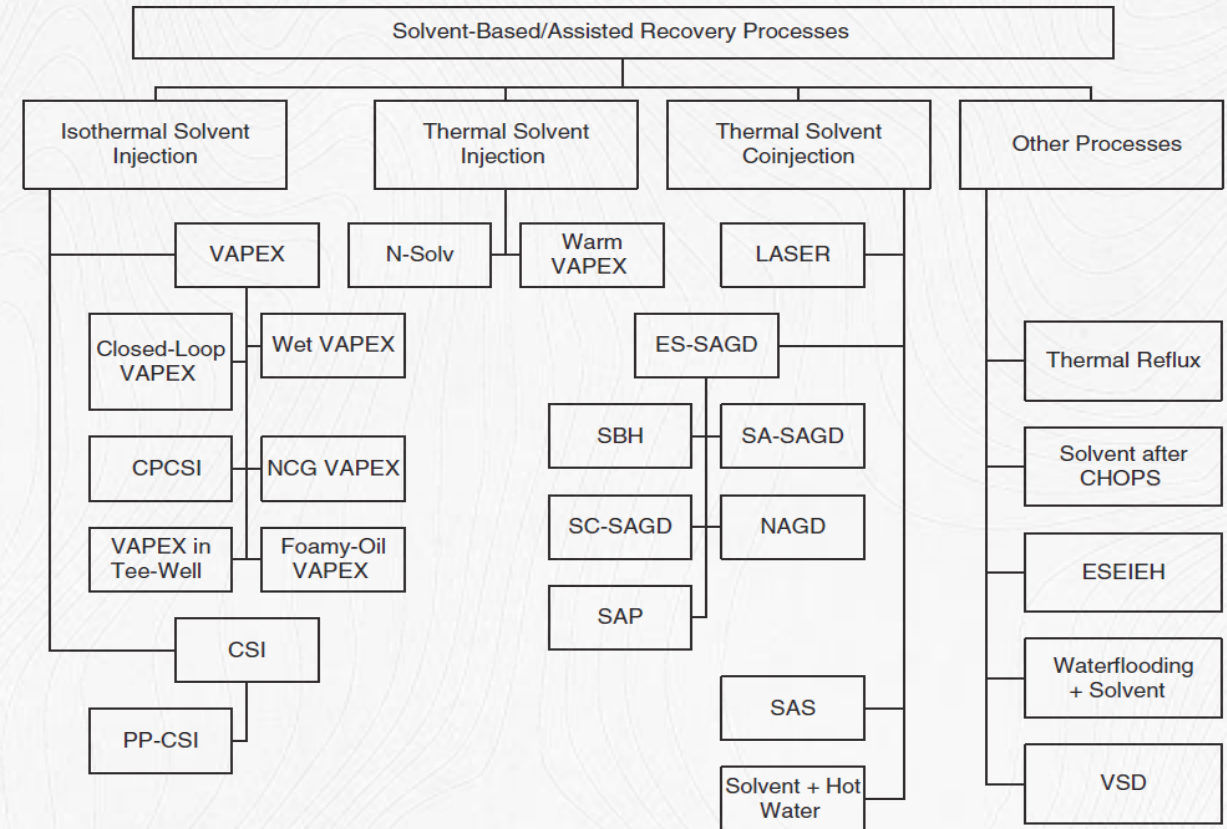
# Solvent-Based Recovery Processes

## Hybrid Processes

- Addition of heat to VAPEX
- Addition of solvent to SAGD/CSS/Steamflood

## Examples

- Heated solvent injection (N-Solv)
  - Solvent is injected at the dew point of the reservoir condition
- Expanding solvent SAGD (ES-SAGD)
- Solvent aided SAGD (SA-SAGD)
- Steam and gas push (SAGP)
- VAPEX initiated by SAGD (SAVEX)
- Steam alternating solvent (SAS)
- Liquid addition to steam for ER (LASER)



# Key Development Areas

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## Solvent Process Wizard

- For Quicker Model Setups
- Boundary leak-off considerations



## Post-Processing Workflows

- For arriving to insights sooner



## Advanced Process Dynamics

- Viscosity Mixing
- Rate-Dependent Dissolution/Exsolution
- Asphaltene Deposition
- Comp-Dependent Rel Perms



## Advanced Well Dynamics

- Extensibility of well controls
- Coupling to a strong wellbore model



## Modelling More than 3-Phases

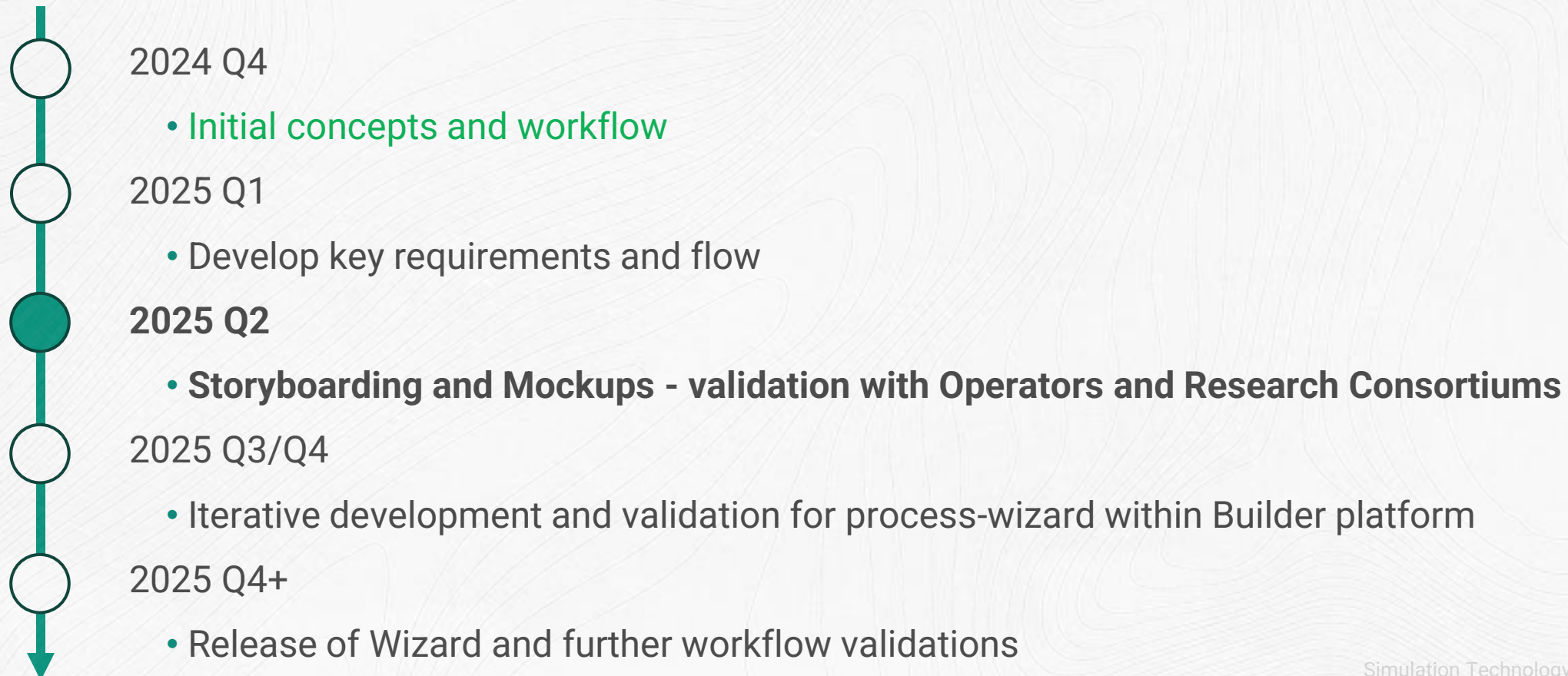
- Multiple oleic phases

- 
- > Workstream Development Plans
  - > Current Status of Developments
  - > What is Coming Next

# Solvent Process Wizard

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Creation of a comprehensive guided task for providing step-by-step input support for solvent studies. Intended to include the key physics of consideration and intelligent default values



# Solvent Process Wizard

- **Streamlined Workflow**
  - Structured, step-by-step process
- **Reduced Complexity**
  - Automates key setup tasks
- **Improved Consistency & Accuracy**
  - Standardization in model setup
- **Friendly Design**
  - Guided inputs, tooltips, and error handling to support both expert and new users

Solvent Wizard - Reservoir & Fluid Model Setup

01 Fluids

02 Physics

03 Relative Permeability

04 Field Strategy

05 I/O Control

### Solvent Components

Select the hydrocarbon components that should be treated as solvents in the simulation. For each well pair, you can enable output options for Solvent/Oil, Solvent/Water, and Steam/Oil ratios. Both rate-based and cumulative outputs can be enabled simultaneously. At least one output type must be selected.

COMPONENT	TREAT AS SOLVENT
CH4	<input checked="" type="checkbox"/>
IC4toNC5	<input checked="" type="checkbox"/>
FC6toFC8	<input type="checkbox"/>
FC9toFC11	<input type="checkbox"/>
Bitumen	<input type="checkbox"/>

### Well Pair Outputs

Select injector... Select producer... + Add Well Pair

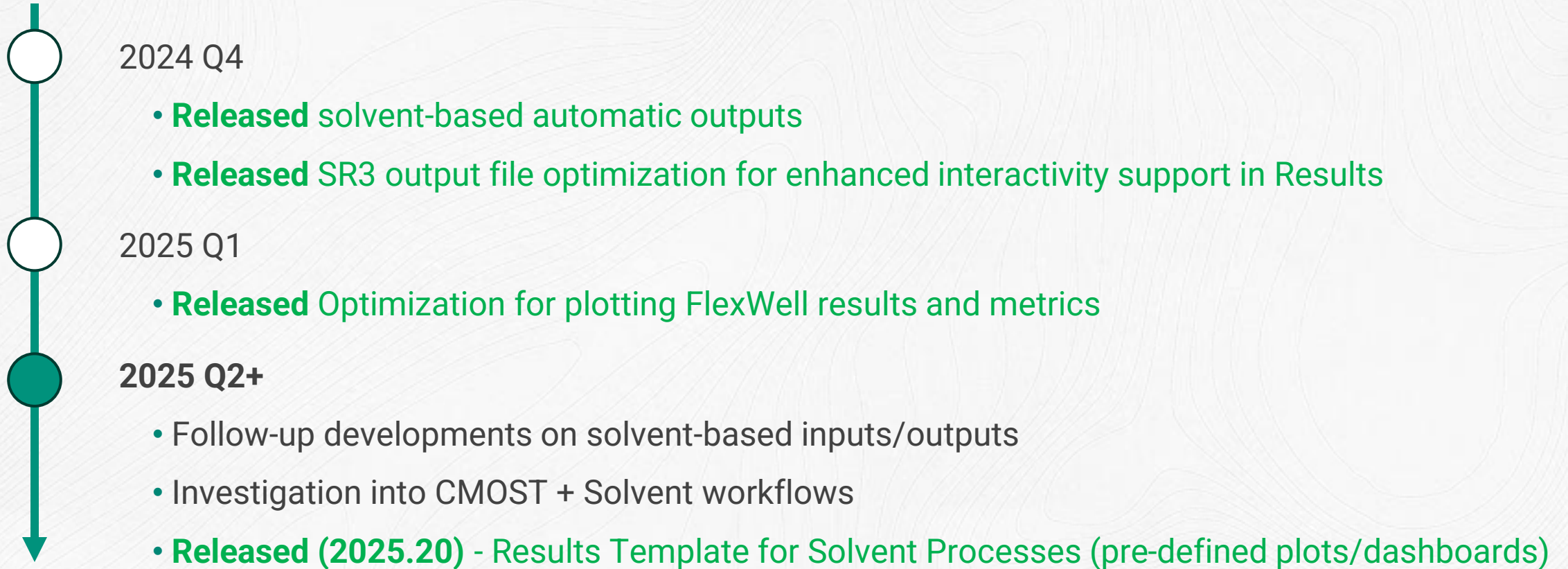
WELL PAIR	SOLVENT/OIL OUTPUT	SOLVENT/WATER OUTPUT	STEAM/OIL OUTPUT	STEAM/OIL OUTPUT (COMP)	ACTIONS
Injector: INJ_SAGD_1 Producer: PROD_SAGD_1	<input checked="" type="checkbox"/> Enable <input checked="" type="checkbox"/> Rate	<input checked="" type="checkbox"/> Enable <input checked="" type="checkbox"/> Rate	<input type="checkbox"/> Enable	<input type="checkbox"/> Enable	<input checked="" type="button"/> Remove

← Back Finish

# Post-Processing Workflows

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Ability to 'Tag' Solvent components and auto-output key metrics for analysis



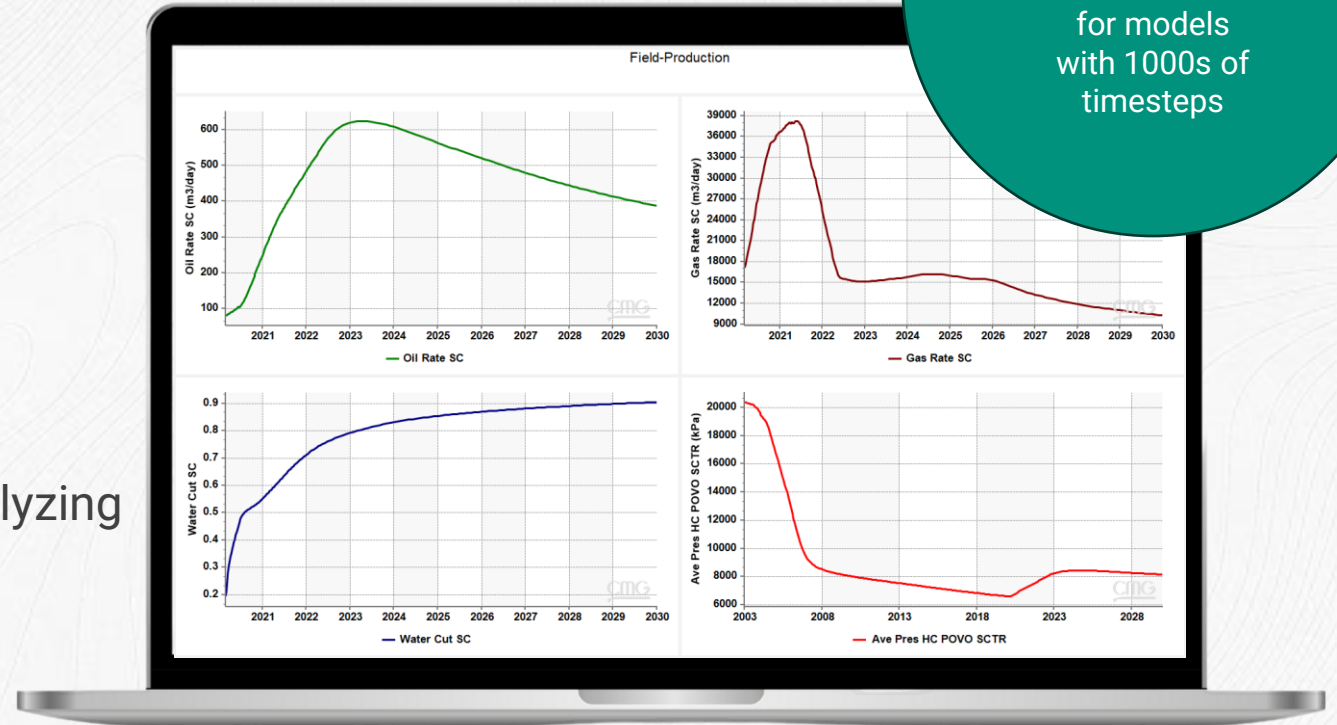
# Performance Enhancements via SR3 Optimization



## SR3 Optimized for Results performance:

- Optimize SR3 files in CMG Launcher
- Load simulation result files (SR3) significantly faster in RESULTS
- Plot timeseries instantaneously
- Profile plotting speed drastically improved
  - **Under 3s**
- Works with all three simulators IMEX, GEM and STARS
- Spend less time waiting and more time analyzing

Up to 60x  
**faster**  
for models  
with 1000s of  
timesteps



# Greatly reduce the time spent in Results dashboards and formula setups

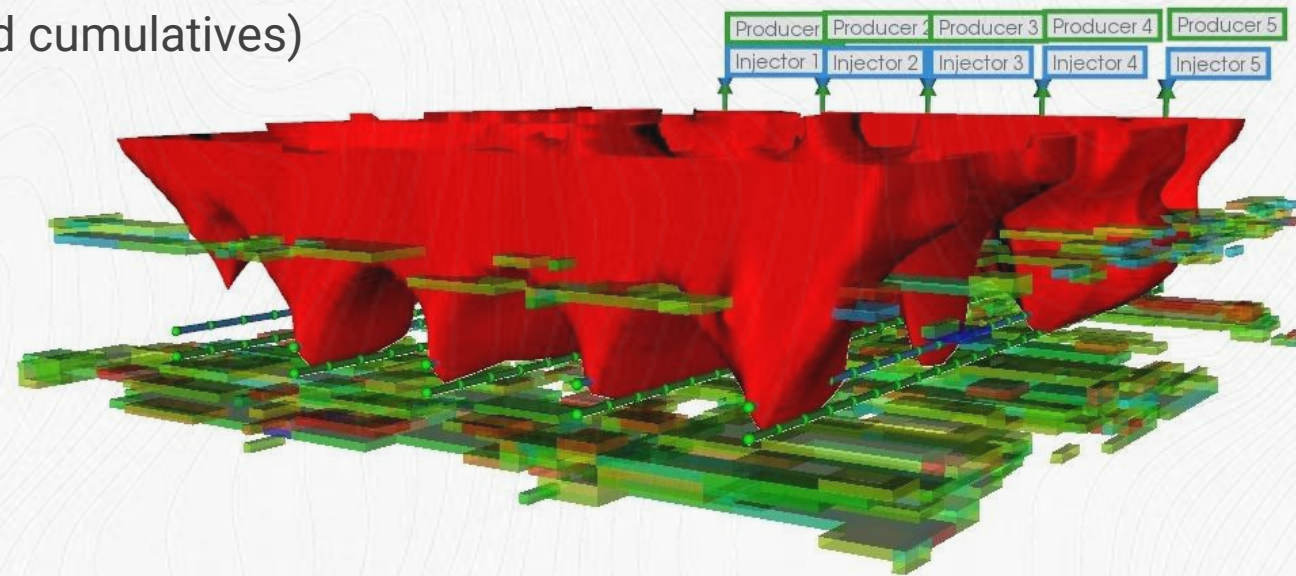


Tag Components as 'Solvents' for output purposes – auto-outputs:

- Solvent split from oil (i.e. non-solvent oil rates and cumulatives)
- Solvent inj and prod, rates and cumulatives
- Solvent-Oil Ratio, inst. and cumulative
- Solvent-Water Ratio, inst. and cumulative
- Solvent Recovery

Outputs available per-sector and well pair

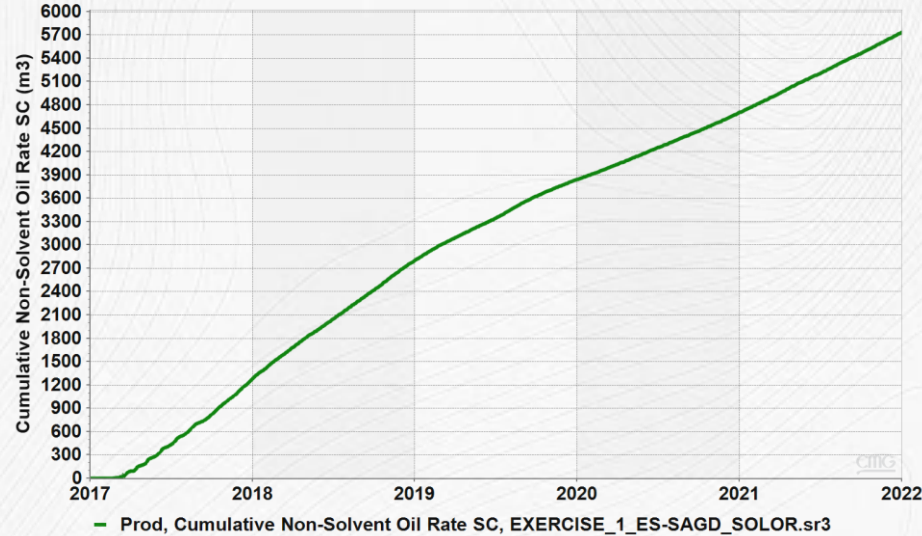
Reduce challenges in defining uncertainty & optimization workflows in **CMOST**.



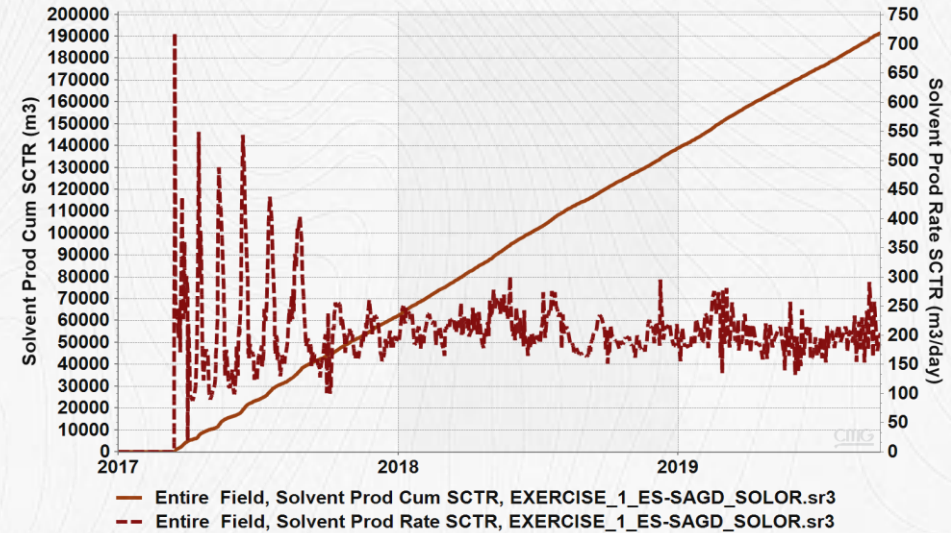
# Greatly reduce the time spent in Results dashboard and formula setups



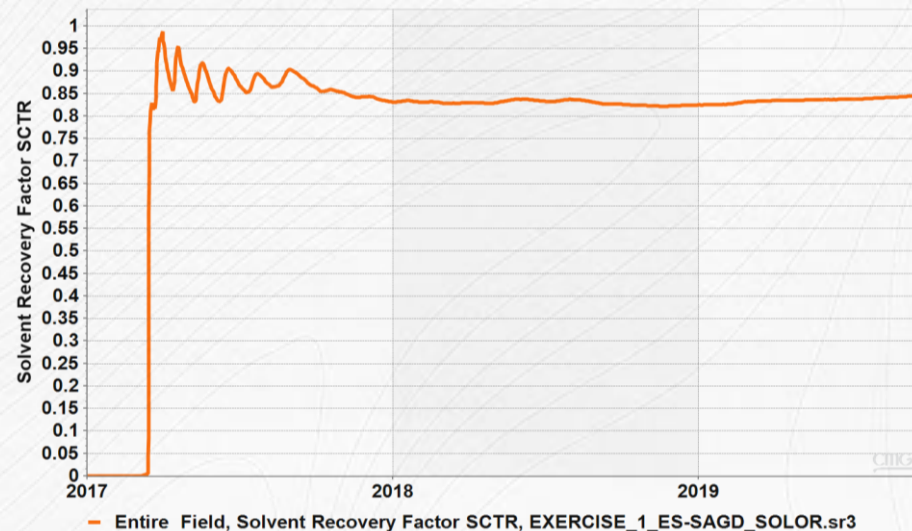
Net Non-Solvent Cumulative Oil Production



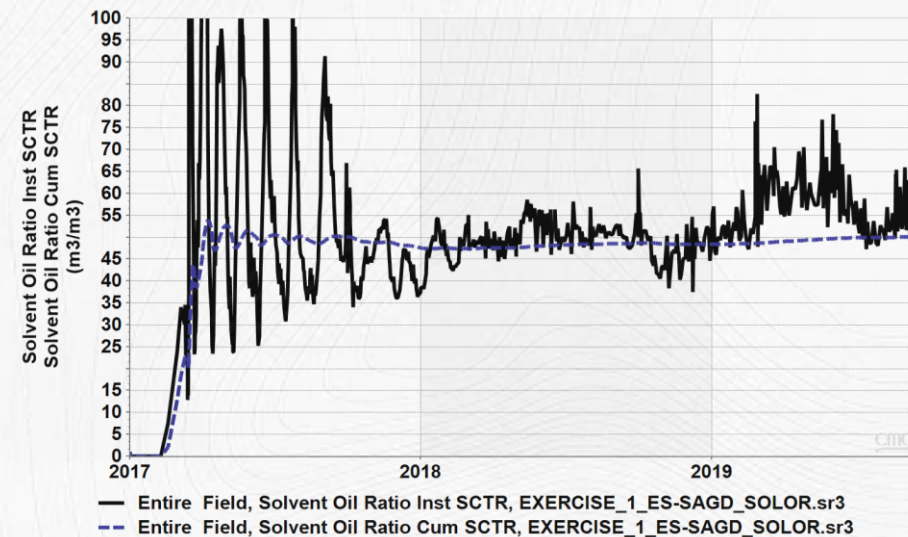
Produced Solvent Rate and Cumulative



Dynamic Solvent Recovery Factor



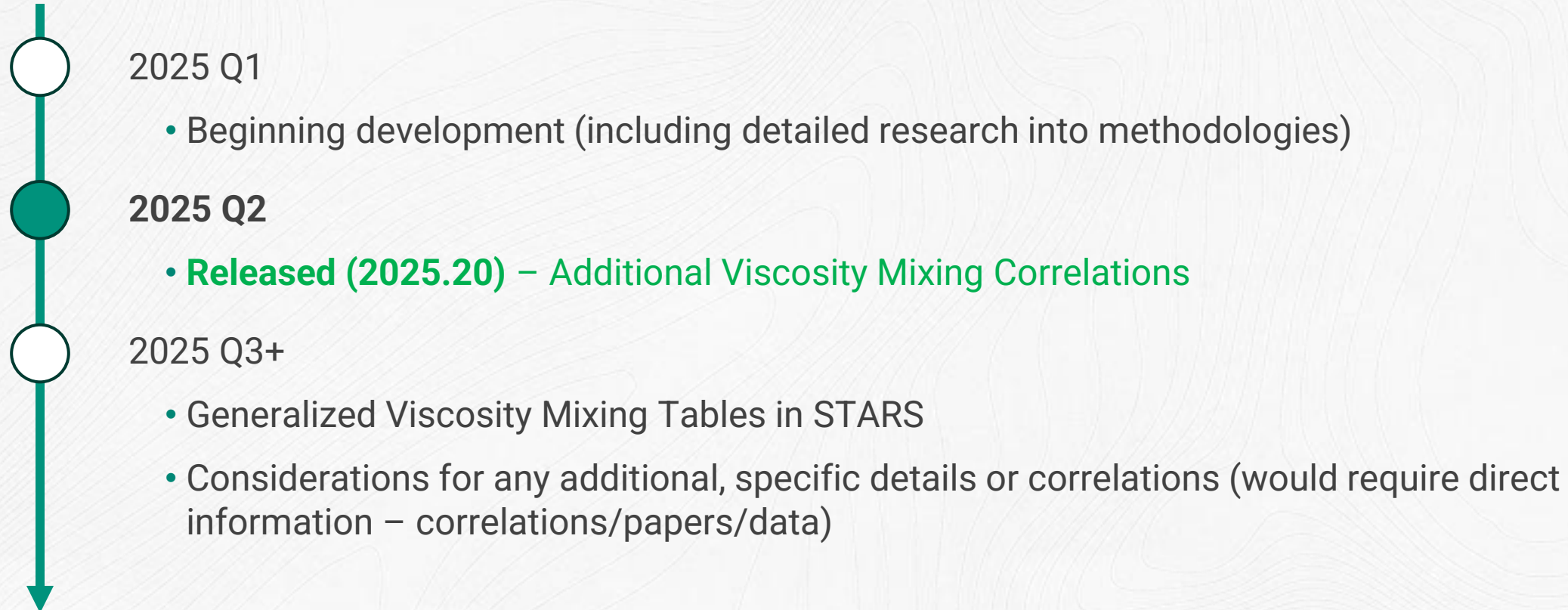
Instantaneous and Cumulative Solvent-Oil Ratio





# Process Dynamics – Viscosity Mixing

Flexible input for a Generalized Viscosity Mixing Table Format into STARS. Different tables for each phase; pressure, temperature and composition effects.





# Process Dynamics – Viscosity Mixing

## Prior to 2025.10

- Linear
- Non-Linear
  - Log-Log
  - Mole or Mass-Fraction based

## New in 2025.20

- Bingham
- Cragoe
- Mehrotra
- Chirinos
- Power-Law
- Arrhenius
- Expanded-Fluid (EF)

Linear	$\mu_{mix} = \sum_{i=1} x_i \mu_i$
Power-law	$\mu_{mix}^n = \sum_{i=1} x_i \mu_i^n$
Arrhenius	$\ln(\mu_{mix}) = \sum_{i=1} x_i \ln(\mu_i)$
Double-double log	$\log(\log(\mu_{mix} + 1)) = \sum_{i=1} x_i \log(\log(\mu_i + 1))$
Kendel	$\frac{1}{3} \ln(\mu_{mix}) = \sum_{i=1} \frac{1}{3} x_i \ln(\mu_i)$
Bingham	$\frac{1}{\mu_{mix}} = \sum_{i=1} \frac{x_i}{\mu_i}$
Crago	$\frac{1000 \ln(20)}{\ln\left(\frac{\mu_{mix}}{5 \times 10^{-4}}\right)} = \sum_{i=1} x_i \frac{1000 \ln(20)}{\ln\left(\frac{\mu_i}{5 \times 10^{-4}}\right)}$
Forth-root	$\mu_{mix}^{-0.25} = \sum_{i=1} x_i \mu_i^{-0.25}$
Wilkie	$\mu_{mix} = \frac{\sum_{i=1} x_i \varphi_i \mu_i}{\sum_{i=1} x_i \varphi_i}$
Mehrorta	$\log(\mu_{mix} + 0.7) = \sum_{i=1} x_i \frac{M_i}{M_{mix}} \log(\mu_i + 0.7)$
Refutas	$A \log(\log(\mu_{mix} + 0.8)) + B = \sum_{i=1} (A x_i \log(\log(\mu_i + 0.8)) + B)$
Chirinos	$\log\left(\log\left(\frac{\mu_{mix}}{\rho_{mix}} + 0.7\right)\right) = \sum_{i=1} x_i \log\left(\log\left(\frac{\mu_i}{\rho_i} + 0.7\right)\right)$

# Process Dynamics – Viscosity Mixing

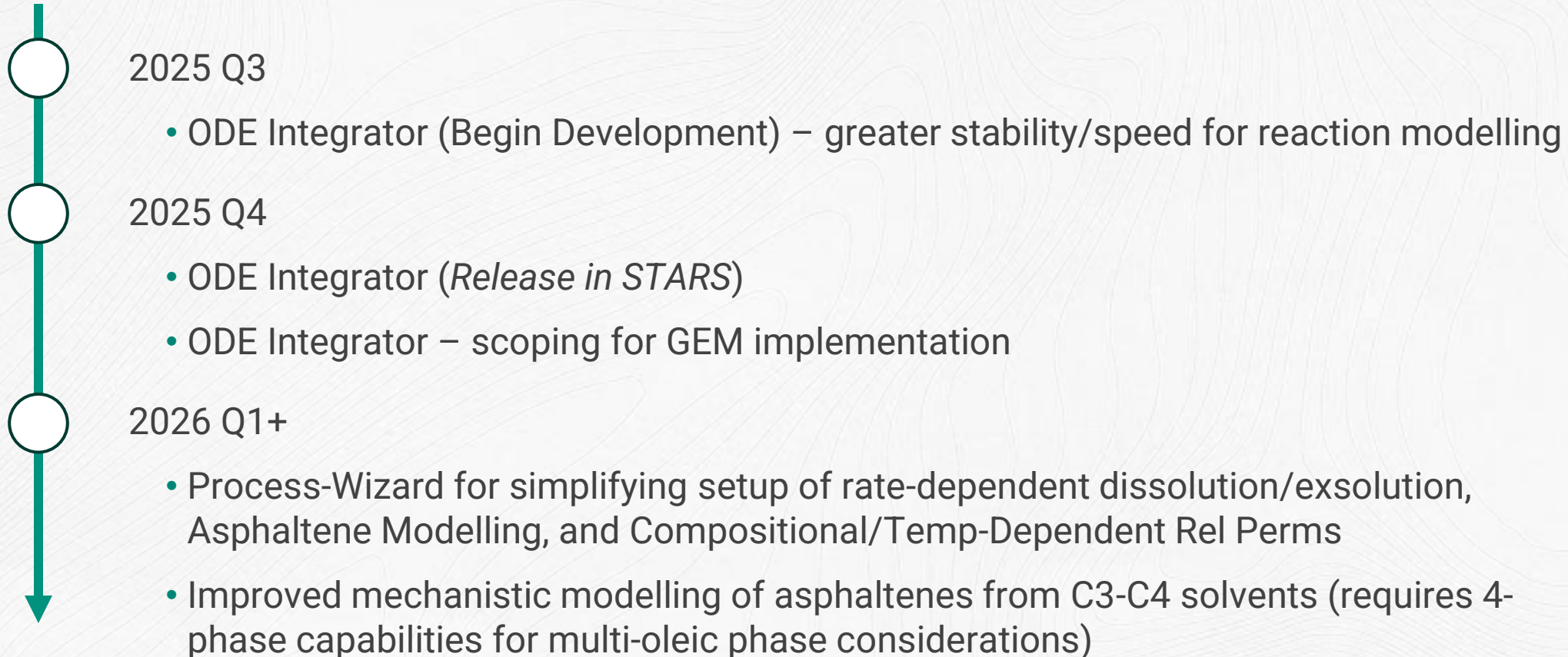


GENVISTABLE			OIL							
PRESS			2000	3000	4000					
TEMP			50	100						
MFRAC			‘C3’	0.1	0.2	0.3	0.4	0.5	0.6	0.7
**	P	T	FR (C3)	0.1	0.2	0.3	0.4	0.5	0.6	0.7
	1	1		590	470	390	270	230	190	180
	1	2		550	450	370	260	210	170	140
	2	1		600	490	400	282	241	199	187
	2	2		559	458	380	267	217	182	149

# Process Dynamics – Rate Dependent Dissolution/Exsolution, Asphaltene Modelling, & Compositional-Dependent Rel Perms

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Validation of current capabilities to known data and enhancements for easier setup and more stable simulations.





# Advanced Well Dynamics

Enhancing extensibility of simulators to work with complex well controls and scripts.  
Consideration of strong wellbore modelling module.



2024

- Development of Drift-Flux formulation for FlexWell
- PyControl Enhancements to support external scripts

2025 Q1

- **Released** - Drift-Flux formulation in FlexWell, Phase 1 (Co-Current)
- **Released** - Pump with Segregation Enabled in FlexWell

2025 Q2+

- **Released (2025.20)** Drift-Flux, Phase 2 (Counter-Current, Liquid Level); Emulsion-Dispersion
- Downhole Separator with FlexWell
- Further stability of FlexWell to provide greater performance
  - New linear solver, compute velocity implicitly, parallelization of FlexWell

# Original Formulation

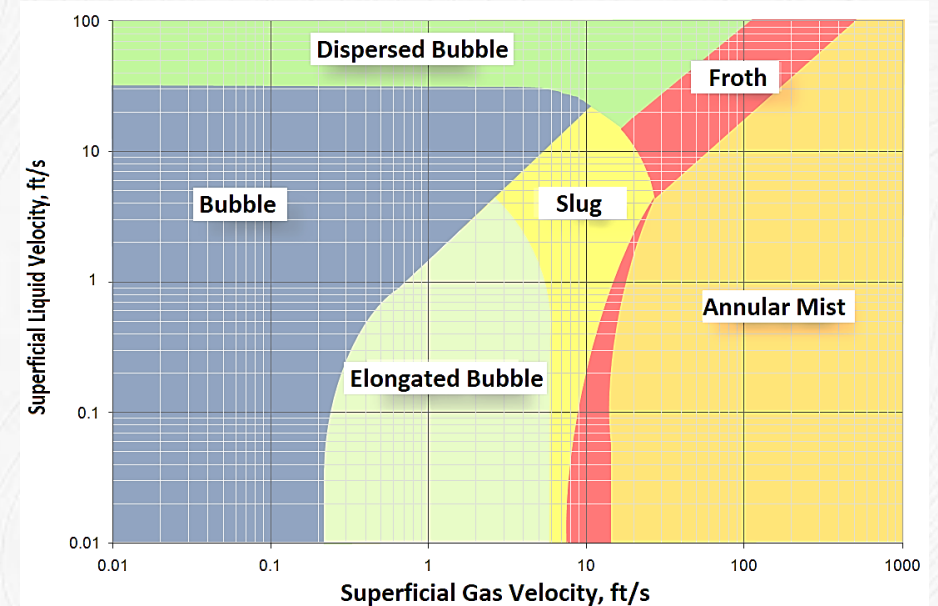
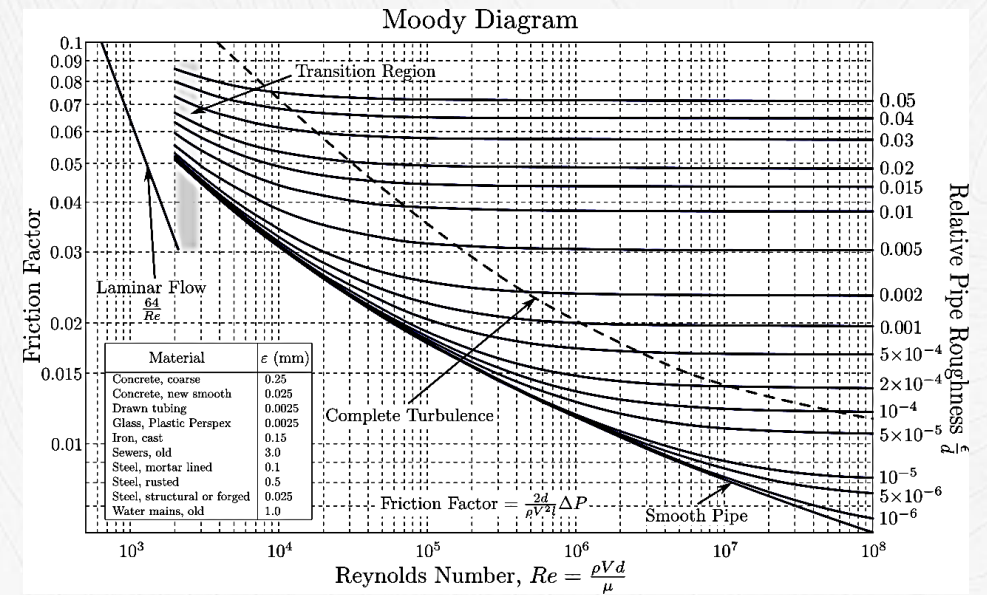
$$f_D = \begin{cases} \frac{64}{Re} & \text{Laminar} \\ \left( \frac{-0.5}{\log_{10} \left( \frac{(\epsilon/D)}{3.7} + \frac{2.51}{Re\sqrt{f_D}} \right)} \right)^2 & \text{Turbulent} \end{cases}$$

$$\Delta\Phi = \frac{\rho f_{cor} f_D v |v|}{2} L/D = K_{cor} v \quad (\text{per phase})$$

- Multi-phase momentum Eqs.
- Flow regime calculation
- High fidelity pressure drop
- Discontinuous correlations

$f_D$

$f_{cor}$



# Updated Formulation

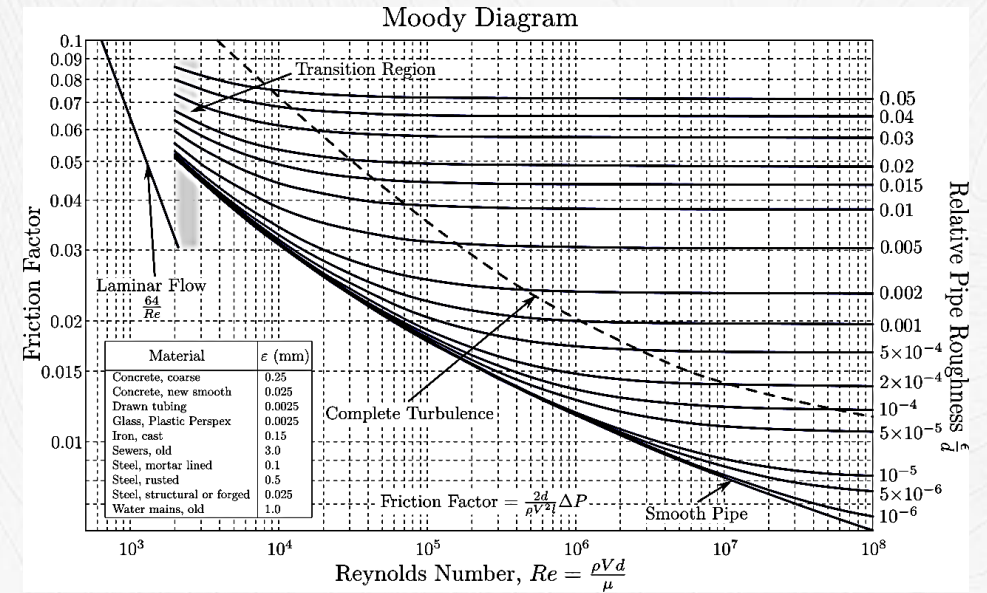
$$f_D = \begin{cases} \frac{64}{Re} & \text{Laminar} \\ \left( \frac{-0.5}{\log_{10} \left( \frac{(\varepsilon/D)}{3.7} + \frac{2.51}{Re\sqrt{f_D}} \right)} \right)^2 & \text{Turbulent} \end{cases}$$

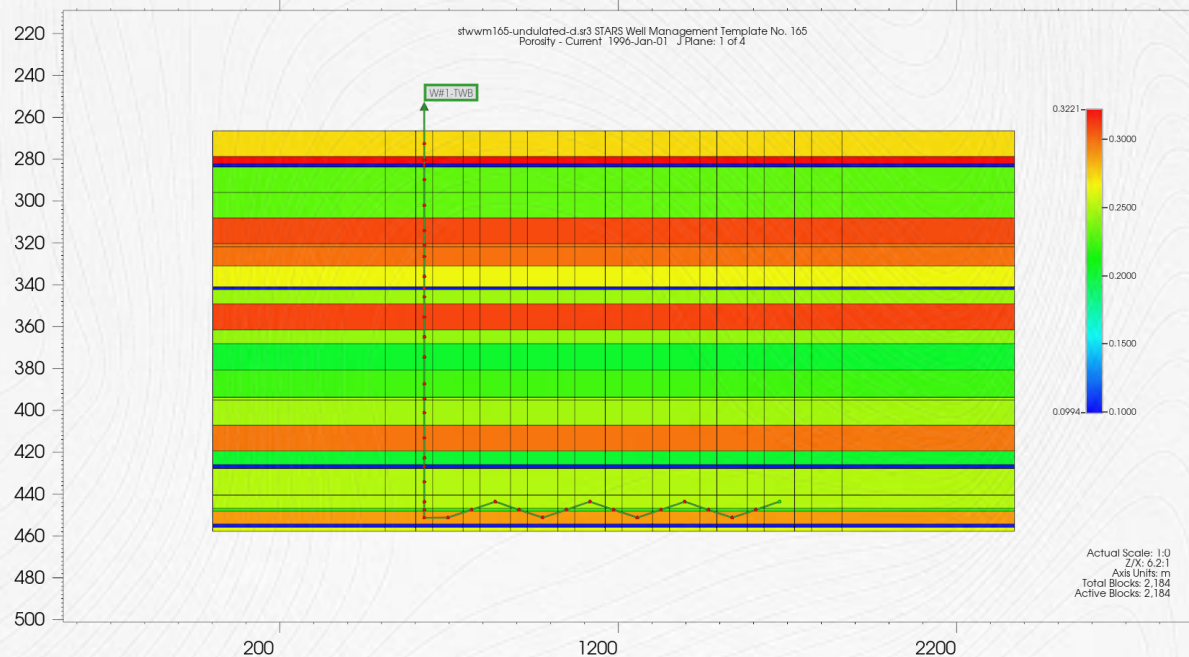
$f_D$

$$\Delta\Phi = \frac{\rho f_D v |v|}{2} L/D = K_{cor} v$$

$$v_g = C_0 v + v_d \quad v_l = \frac{(v - s_g v_d)}{s_l}$$

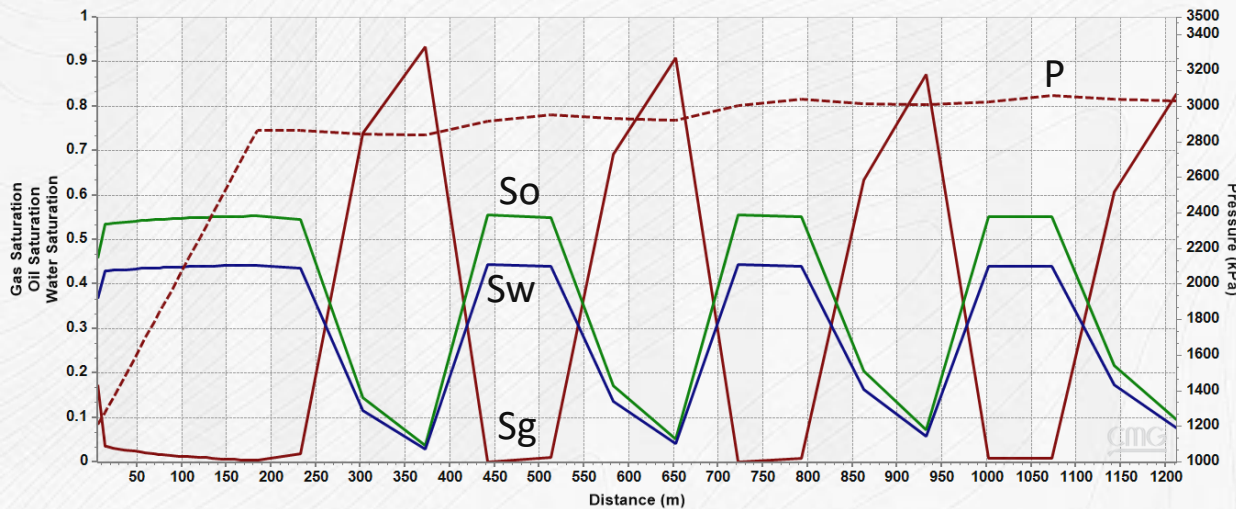
- Unified momentum Eqs.
- Smooth correlations for friction
- More robust for horizontal, undulated, vertical wells
- Pressure drop model used for NOSLIP and DRIFT-FLUX



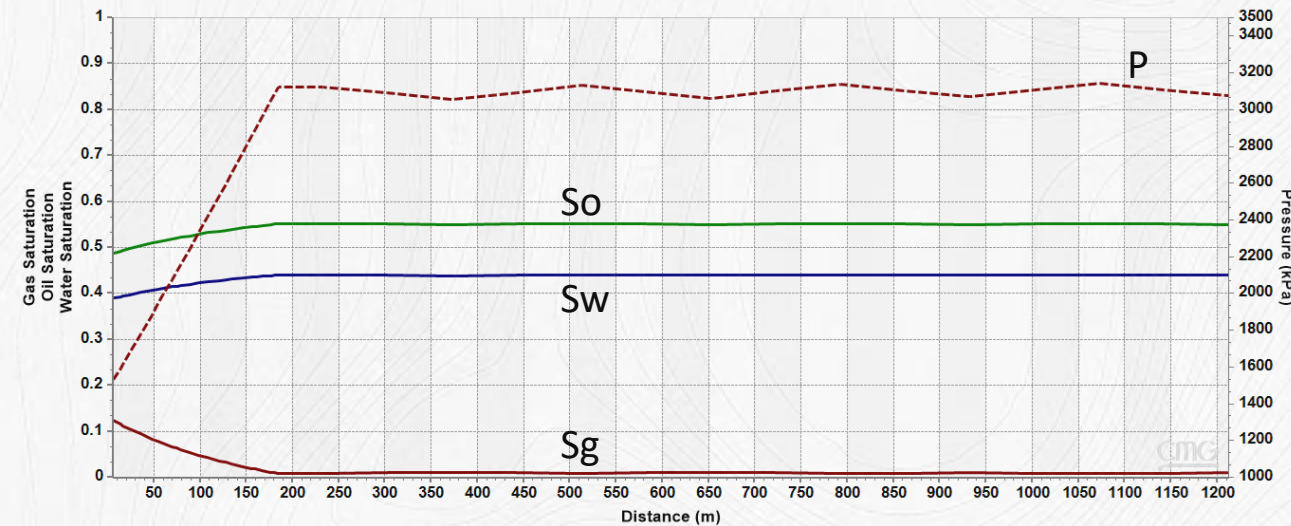


- A preliminary version of DF model has been completed in STARS
- Initial outcomes show that this model can capture the slippage between gas and liquid phases.
- Further tests will be conducted to improve the physics and efficiency of the DF model.

Drift Flux



NOSLIP





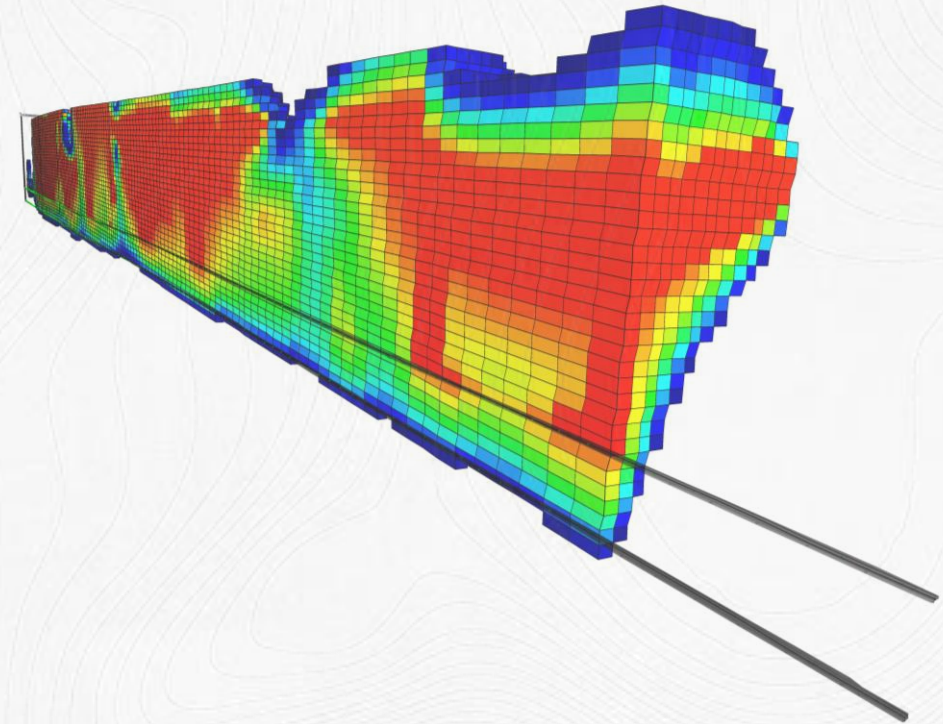
# Advanced Pump Model with All Segregation Options

## Pump Power Correlation

- Power is a function of volumetric flow rate of liquid and differential pressure across the pump
- Define a max power and pressure

## Table Input

- Rates versus Differential Pressure
- Similar to FCD Table setup
- Optional dependencies on Inlet:
  - Temperature
  - Pressure
  - Water Cut
  - Gas Fraction



# Released in 2024 - Greater Understanding of Complex Thermal Well Dynamics



## Improve accuracy of energy transfer with new Fluid Thermal Resistance calculation

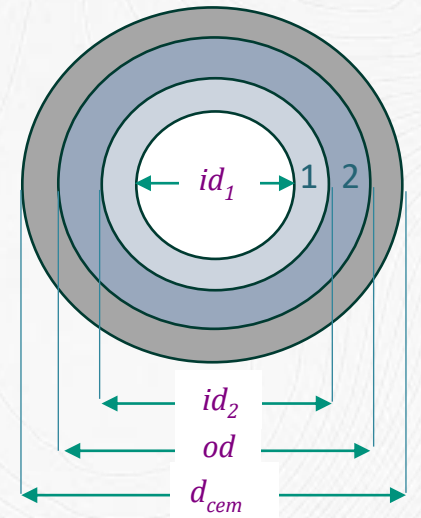
More accurate thermal resistance values through the fluid film near to the tubing and casing walls.

$$res(fl) = \frac{D_h}{\pi N_u K D_{to} L} = \frac{D_{ci} - D_{to}}{\pi N_u K D_{to} L}$$

## Better tuning to field operations with Radial Heat Transfer Outputs in FlexWell

Consider Intermediary temps and heat transfer changes **radially outwards**

- Define multiple concentric casing layers
- Output and plot the temperature along the intermediary points
- **Useful in matching thermal conduction with sensor data** (e.g. insulation or cemented zones)



# Better Fluid Modelling Accuracy for Thermal Runs



## Improved Default Enthalpy for gas-based enthalpy and Lee-Kesler thermal models

Default values of CPG, HVR, and EV are now based on library values of hydrocarbon components with the *nearest molecular mass* (C1 - C20)

- **Greater accuracy for thermal models**
- New defaults can be activated by the option **\*CPG\_LIB**



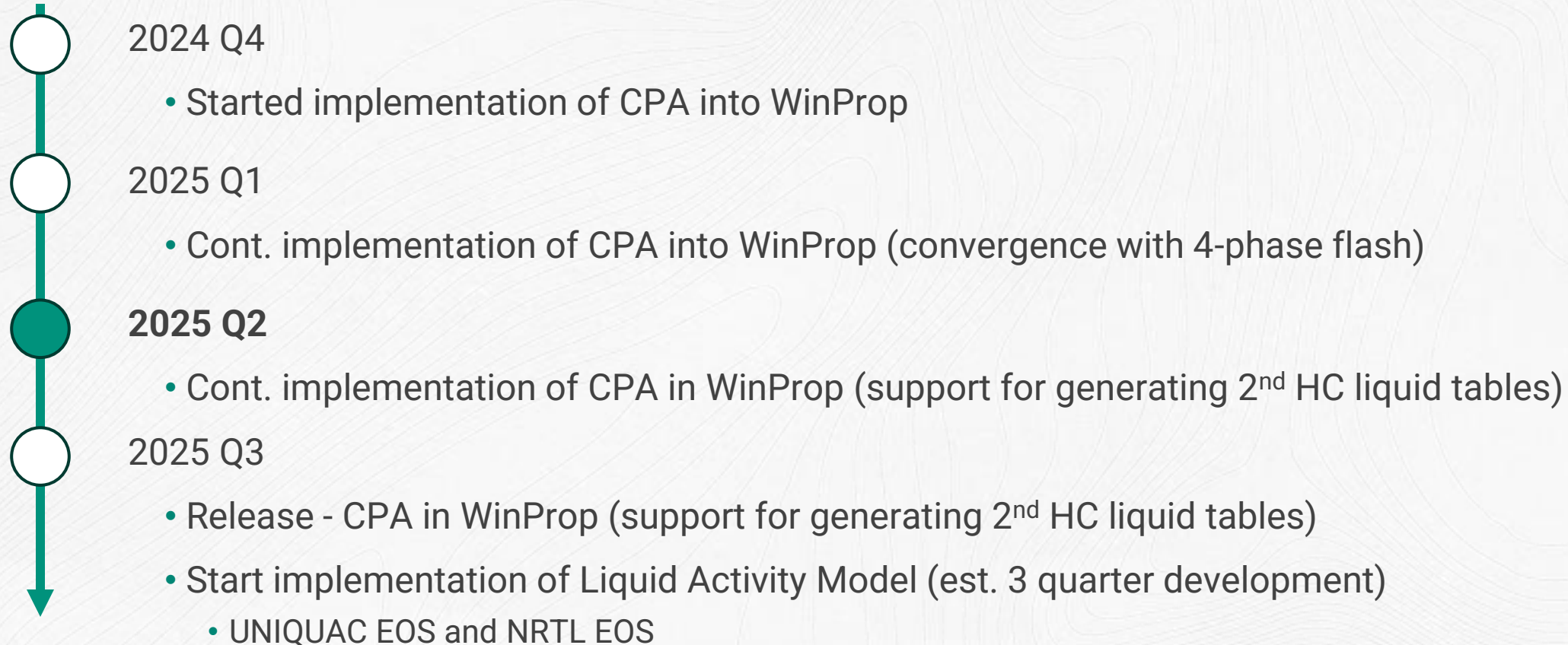
## Faster STARS runs with PVT K-Value Smoothing Option in WinProp

- Modified P-smoothing approach
- Best for Gas-Liquid K-value tables
- Also useful for liquid-liquid tables



# Modelling More Than 3-Phases – Phase 1

Providing ability to model LLV conditions with steam (2 oleic phases). Proposed development to support included comprehensive developments to both STARS and WinProp in 2-phases.





# Implementation of CPA

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1. Research conducted to understand the underlying formulae
2. Implemented:
  - a) Calculate molar-volume for one associated component (H<sub>2</sub>O) with EOS'
  - b) CPA Logic for two associated components
  - c) Ability to read in association energy and volume for association components
3. Current work:
  - a) Enhance general EOS calculations to work with CPA
  - b) Enhance multi-phase flash to work with CPA



# Modelling More Than 3-Phases – Phase 2

Providing ability to model LLV conditions with steam (2 oleic phases). Proposed development to support included comprehensive developments to both STARS and WinProp in 2-phases.



## 2025 Q2

- Common NIST data library
  - **Released (2025.20)** – CO<sub>2</sub> library properties and Visc vs Pressure option
- Start integration of second hydrocarbon phase table in STARS

## 2025 Q3

- Cont. integration of second hydrocarbon phase table in STARS

## 2025 Q4

- **Release** – second hydrocarbon phase table in STARS
- Combined liquid flow for hydrocarbons in STARS



# Accelerating Decisions

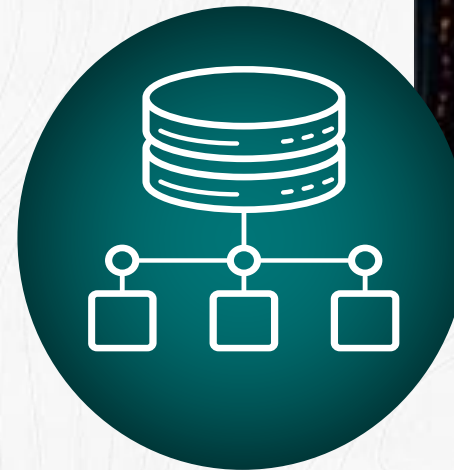
## Speed & Stability Enhancements

# PyControl Enablement with MPI



With the PyControl tool you can script simple-to-complex controls with Python that the simulators can use when running

With this latest extension to MPI, even the largest of models can leverage the external scripting and capabilities that PyControl workflows provide.



AWS Datacentre

<https://aws.amazon.com/compliance/data-center/data-centers/>

# Faster Distributed Runs with the Combinative Solver

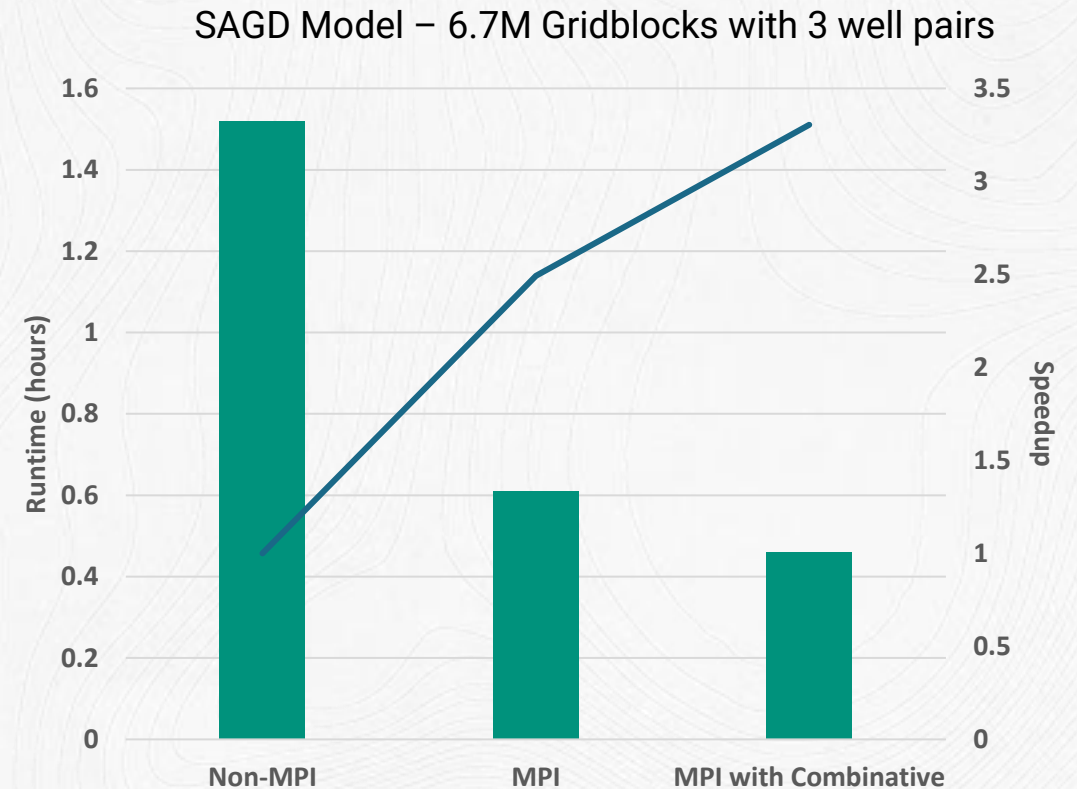



- Impactful with frontal-based processes
- Combinative previously available in IMEX, GEM, and STARS OpenMP (Shared Memory)
- Now Available in IMEX, GEM and STARS MPI (Distributed)
  - Combinative ILU currently deployed in MPI

## Coming Soon:

- Combinative AMG Option for STARS-MPI

**Average 40+% Speedup**  
*In benchmark models with Combinative Solver + MPI*





Thank you!  
Questions?