

Improving the Fluid Property Evaluation for RELAP5-3D

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Background and Purpose

- Fluid properties tabulated by many organizations, including:
 - National Bureau of Standards (NBS)
 - National Institute of Standards and Technology (NIST)
 - American Society of Mechanical Engineers (ASME)
 - Nuclear Regulatory Commission (NRC)
- Besides tables generators, complex mathematical functions of many variables are generated.
 - The generators were computationally intensive compared with table lookup for older computers
- RELAP5 codes read and store needed tables of fluid property values at input then interpolate to obtain properties at a given (P,T) or (P,U)
- Purpose of this study
 - Examine different calculations of the table fluid properties and their derivatives can improve RELAP5-3D calculations

Background

- RELAP5-3D has 33 different fluid table files.
 - 9 are variation on water properties: 8 for light water, 1 for heavy
 - 1967, 1984, 1995 in original and revised form
 - Liquid metals, molten salts, supercritical CO₂
 - Gases: hydrogen, helium, nitrogen, xenon
 - Refrigerants, glycerin, human blood
- All the property files start with the 3 letters “tpf” for Tabulated Properties of Fluids, or Fluid Property Tables for short.
- RELAP5-3D tables are generated in two formats:
 - ASCII and Machine independent binary XDR,

Water Property Accuracy & Mass Error

- The mathematical function, its evaluation, and interpolation all involve errors:
 - Approximation of reality, approximation of transcendental functions, numerical roundoff, and interpolation error
 - Interpolation error is the largest of these and it leads to mass error.
- In 2010, Cliff Davis studied the interpolation grid and found the regions of greatest difference between the function and interpolated values.
 - He created a new grid with more interpolation points in the areas of largest difference.
 - New grid reduced interpolation error
- A second study tested a methodology to improve fluid property derivative calculations

Mass Error and Consistency

- In RELAP5 codes, the derivatives are calculated from the nonlinear Clausius-Clapyron Equations

$$\begin{aligned} \bullet \left(\frac{\partial \rho_j}{\partial U_j} \right)_P &= \frac{v_j \beta_j}{(C_{pj} - v_j \beta_j P) v_j^2} & \left(\frac{\partial T_j}{\partial U_j} \right)_P &= \frac{1}{C_{pj} - v_j \beta_j P} \\ \bullet \left(\frac{\partial \rho_j}{\partial P} \right)_{U_j} &= \frac{C_{pj} v_j \kappa_j - T_j (v_j \beta_j)^2}{(C_{pj} - v_j \beta_j P) v_j^2} & \left(\frac{\partial T_j}{\partial P} \right)_{U_j} &= \frac{P v_j \kappa_j - T_j v_j \beta_j}{C_{pj} - v_j \beta_j P} \end{aligned}$$

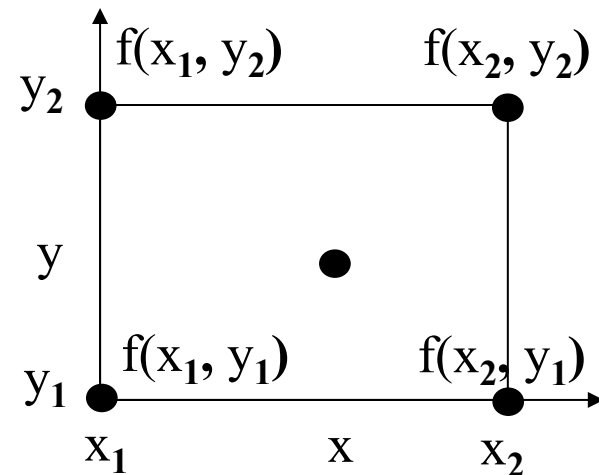
- $j = f$ or g (for liquid or gas)
- specific volume v
- specific heat capacity C_p
- isobaric coefficient of thermal expansion β
- isothermal coefficient of compressibility κ

Goal: Less Mass Error via Better Approximation

- The numerical approximations to Governing Partial Differential Equations arise from LINEARIZATION at every step.
 - No nonlinear terms of primary variables, namely P , U_f , U_g , α_g , and X_n , are allowed
 - Linearization by putting primary variables factors at old time and leaving only one factor at new time
 - Non-primary variables are replaced by functions of primary variables. These are linearized by
 - Using old time values
 - Or replacing nonlinear terms by first order Taylor polynomials
- Since the numerical system is linear, would linear derivatives be more CONSISTENT than default Clausius-Clapyron approximation?
- Would some other form that combines the two reduce mass error?

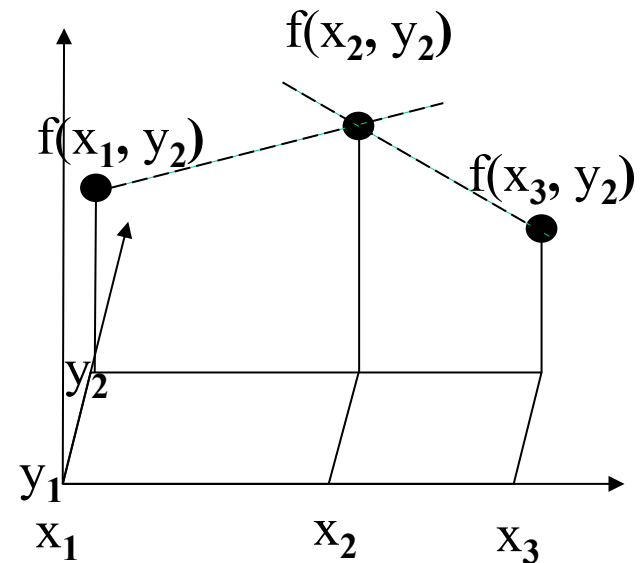
Second Study & Theory

- A second study showed promise for linear approximation
 - 1984 water properties with the detailed (U, P) mesh
 - A simple 2-vol test case showed the linear interpolant produced less mass error than the normal RELAP5-3D calculation
- Coding was written to carry out a more complete study.
- Calculate linear interpolants of the 8 derivatives based on the four corners of the bounding rectangle
- $\alpha(x) = \frac{x-x_1}{x_2-x_1}, \beta(y) = \frac{y-y_1}{y_2-y_1}$
- $\left(\frac{\partial f}{\partial x}\right)_{\text{lin}} = \frac{(1-\beta)[f(x_2,y_1)-f(x_1,y_1)]+\beta[f(x_2,y_2)-f(x_1,y_2)]}{x_2-x_1}$
- $\left(\frac{\partial f}{\partial y}\right)_{\text{lin}} = \frac{(1-\alpha)[f(x_1,y_2)-f(x_1,y_1)]+\alpha[f(x_2,y_2)-f(x_2,y_1)]}{y_2-y_1}$
- In example, $\alpha=0.6, \beta=0.4$



Theory – Jump Discontinuities

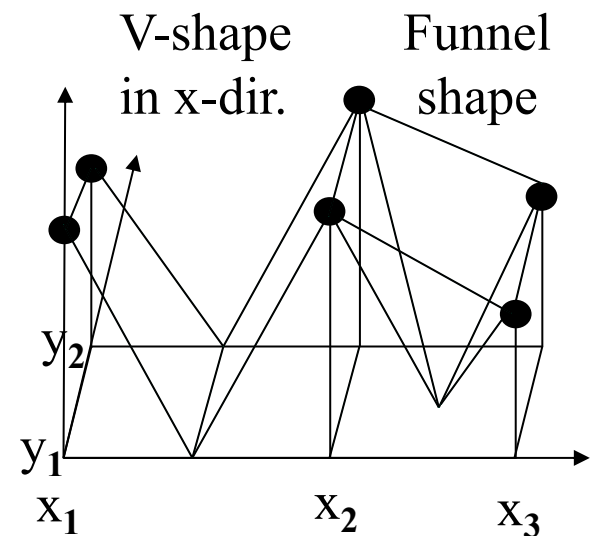
- reduced mass error for 2-vol test case, but not for Edward’s Pipe
 - Reason: Discontinuity of the derivatives at rectangle edges
- Consider three points of the energy grid: x_1, x_2, x_3 .
 - In left rectangle at $y = y_2$,
 - $\alpha(x_2) = (x_2 - x_1)/(x_2 - x_1) = 1$
 - In right rectangle at $y = y_2$,
 - $\alpha(x_2) = (x_2 - x_2)/(x_3 - x_2) = 0$
- The left and right derivative at y_2 do not match:
 - $\lim_{\epsilon \rightarrow 0} \left(\frac{\partial f}{\partial x} \right)_{\text{lin}} (x_2 - \epsilon, y_2) = \frac{f(x_2, y_2) - f(x_1, y_2)}{x_2 - x_1}$
 - $\lim_{\epsilon \rightarrow 0} \left(\frac{\partial f}{\partial x} \right)_{\text{lin}} (x_2 + \epsilon, y_2) = \frac{f(x_3, y_2) - f(x_2, y_2)}{x_3 - x_2}$
- Linear interpolant has **jump discontinuity**
- Default Clausius-Clapyron derivative does not



Theory – Hybrid Interpolation

- Combine the best of both - calculate a weighted average of the default Clausius-Clapyron derivatives and the linear interpolant.
- Weighted Average = $\omega * \left(\frac{\partial f}{\partial x}\right)_{\text{def}} + (1-\omega) * \left(\frac{\partial f}{\partial x}\right)_{\text{lin}}$
 - 1.0 produces fixed Default derivatives
 - 0.0 produces fixed Linear
 - 0.5 produces fixed Hybrid
 - -1.0 triggers variable V-shaped weighting in y-dir.
 - -2.0 triggers Funnel weighting
- Other shapes were considered, but not programmed

similar for y-direction



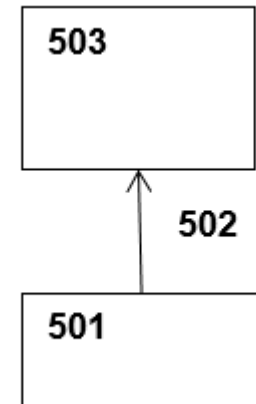
Coding Implementation

- Subroutine RMFLDS: 120-129 hydrodynamic system card adds weight as final value
 - The fluid can change from system to system, so the weighting must also.
- Subroutine LINPOLATE: new F90 routine that calculates:
 - Default or linear-interpolant derivatives
 - Matching fluid properties (Consistency)
 - T, rho, kappa, beta, Cp, k, mu, s, specific volume
 - Weighted average of linear and default
- Other modifications
 - SVPUPU, ISTATE, STPUPU – Calls to LINPOLATE for H2ON only
 - STATEP – uses weighted averages for H2ON only
- Test Cases
 - Two volume “Box” insurge and outsurge
 - Edwards Pipe Blowdown, Typical PWR, Moby Dick, Mixubub

Testing with Improved P,T Fluid Grid

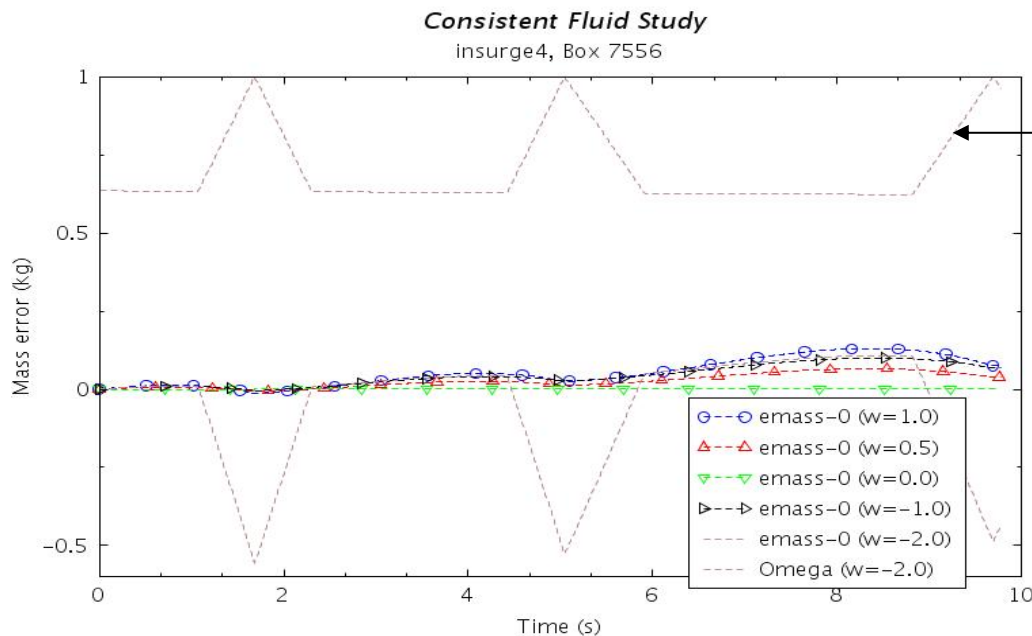
- Two volume insurge “Box” problem
 - P,T set at lowest values inside the bounding rectangle
 - P,T increased to maximum for box at 10s
 - Table shows ormalized mass errors during simple filling calculations

| Box | Midpoint values | | Comments |
|------|-----------------|-----------|---|
| | Pressure (MPa) | Temp. (K) | |
| 113 | 1.10E-03 | 275.33 | Liquid (worst box in Region 7) ¹ |
| 341 | 1.50E-03 | 286.25 | Vapor (worst box in Region 8) |
| 1936 | 9.00E-02 | 410.0 | Vapor (worst box in Region 10) |
| 5629 | 7.50 | 377.5 | Liquid (average box in Region 1) |
| 5651 | 7.50 | 617.50 | Vapor (average box in Region 2) |
| 7400 | 16.00 | 293.75 | Liquid (worst box in Region 9) ¹ |
| 7445 | 16.00 | 627.50 | Vapor (worst box in Region 2) |
| 7446 | 16.00 | 632.50 | Vapor (worst box in Region 4) |
| 7495 | 16.00 | 936.575 | Vapor (high temperature) |
| 7556 | 17.25 | 622.5 | Liquid (worst box in Region 1) |
| 7669 | 18.25 | 627.5 | Liquid (worst box in Region 3) |



Box 7556 Mass error, INSURGE Case

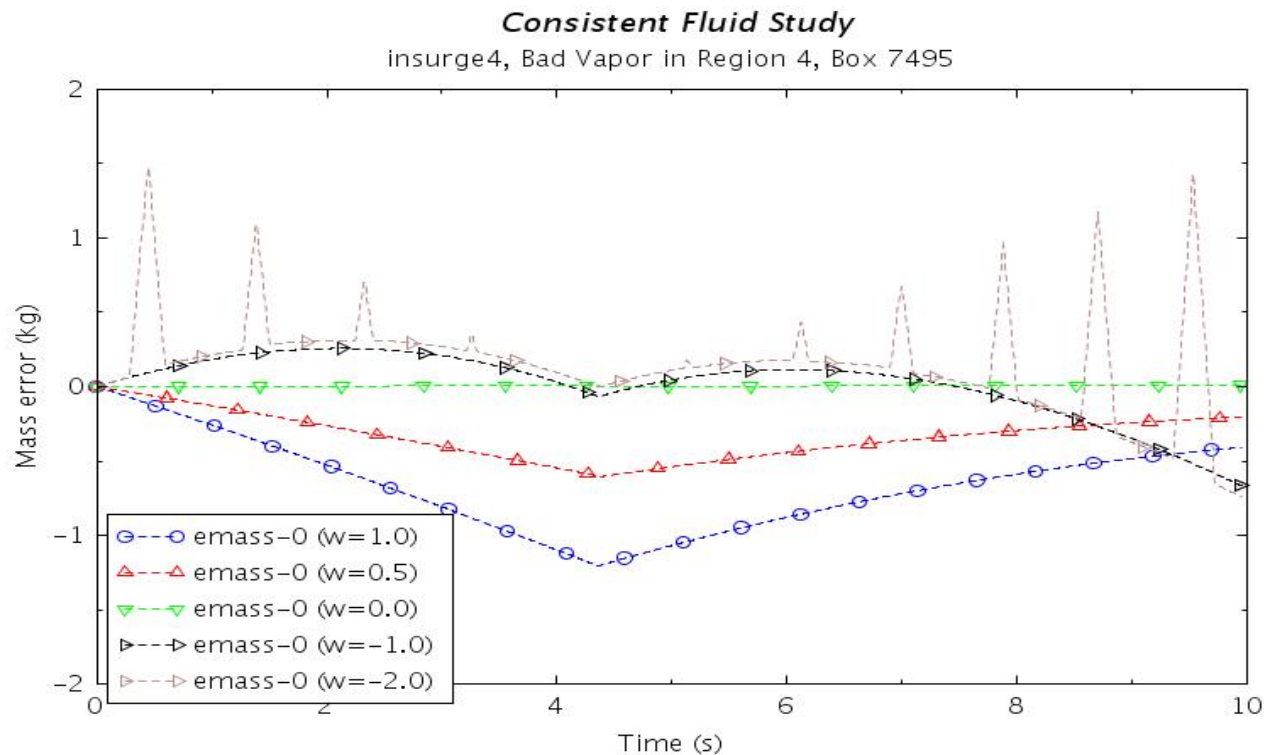
- Upper curve shows the variable ω -values generated when user selects the funnel shape (input flag $\omega = -2$).
 - The lower curve shows the mass error generate with these values
- Values $-1 \leq \omega < 1$ outperform default RELAP5-3D ($\omega = 1.0$).



ω -values from
flag $\omega = -2$

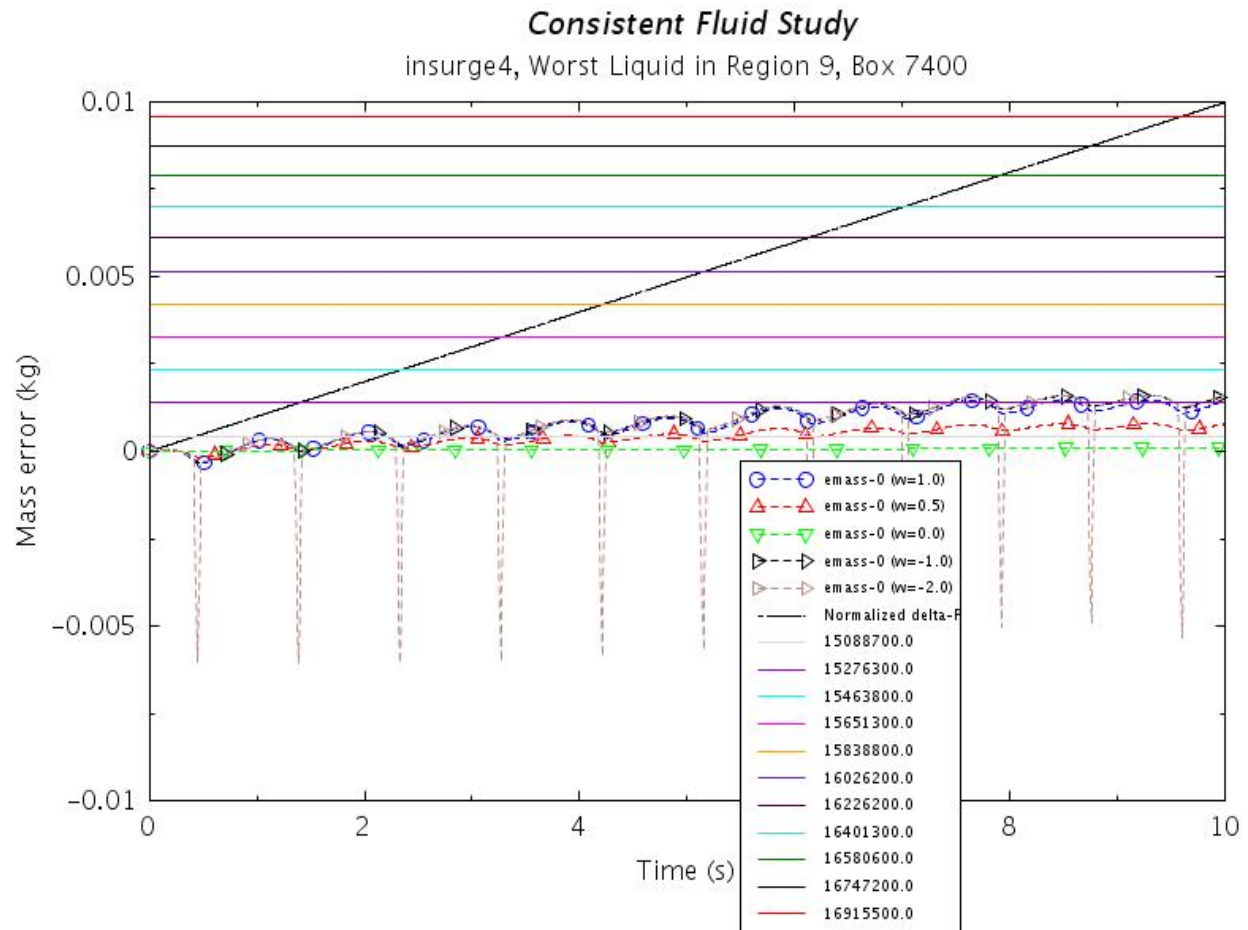
Box 7556 Mass error, INSURGE Case

- Values $-1 \leq \omega < 1$ outperform default RELAP5-3D ($\omega = 1.0$)
- What causes the spikes for funnel-shaped ω ?



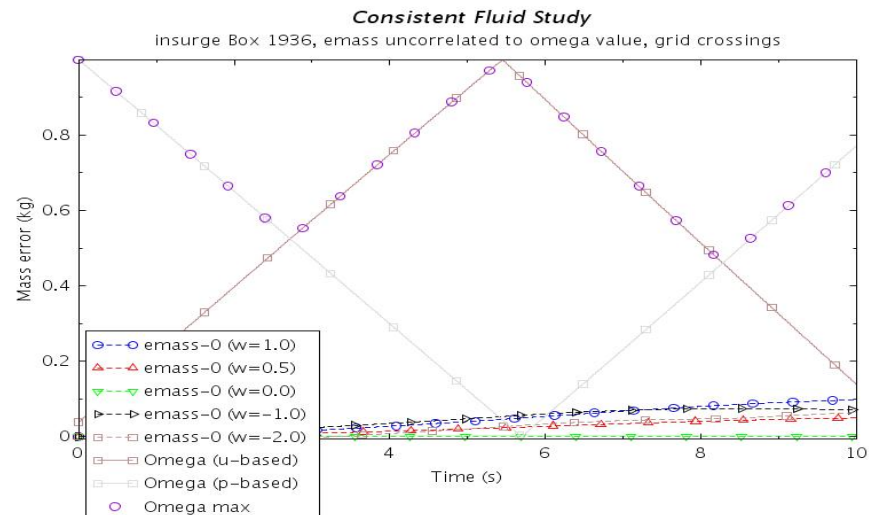
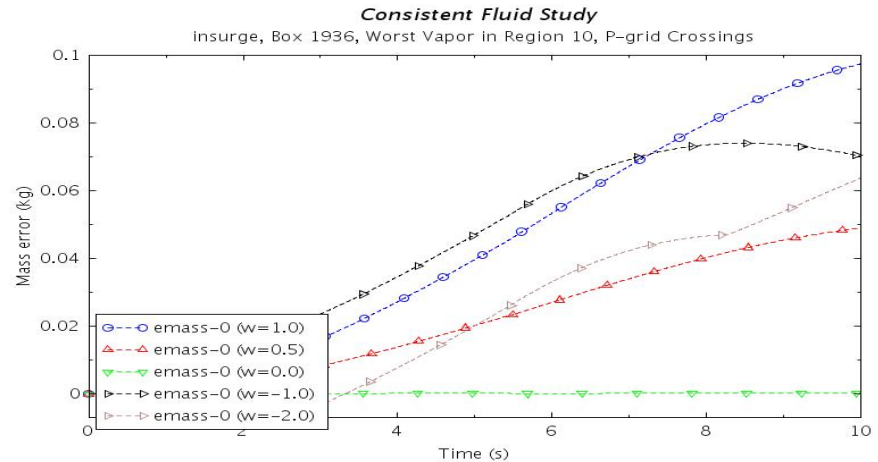
Box 7400 Mass error, INSURGE Case

- Black and horiz. lines are **P normalized**
- $-1 \leq \omega < 1$ default R5-3D ($\omega = 1.0$)
- Local *minima* for ALL mass error graphs at **pressure grid values**.
 - Also true in prior plots
- What about U grid values?



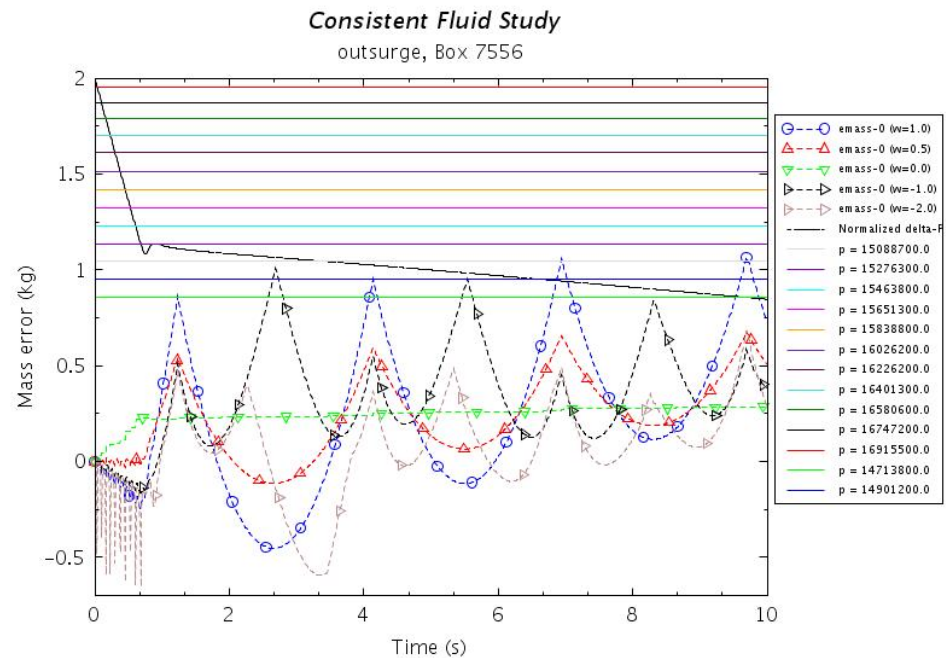
Box 1936 Mass error, INSURGE Case

- V-line gives P-based ω
- Inverted V-line shows U-based ω .
- Funnel ($\omega = -2$) uses $\omega = \max$ of V and inverted-V.
- **Inflection point at the U grid point** for $\omega = -2, -1, 0.5, 1$
- First time $\omega = 0.5$ is worst.
- All INSURGE cases show $\omega = 0$ has almost no mass error.



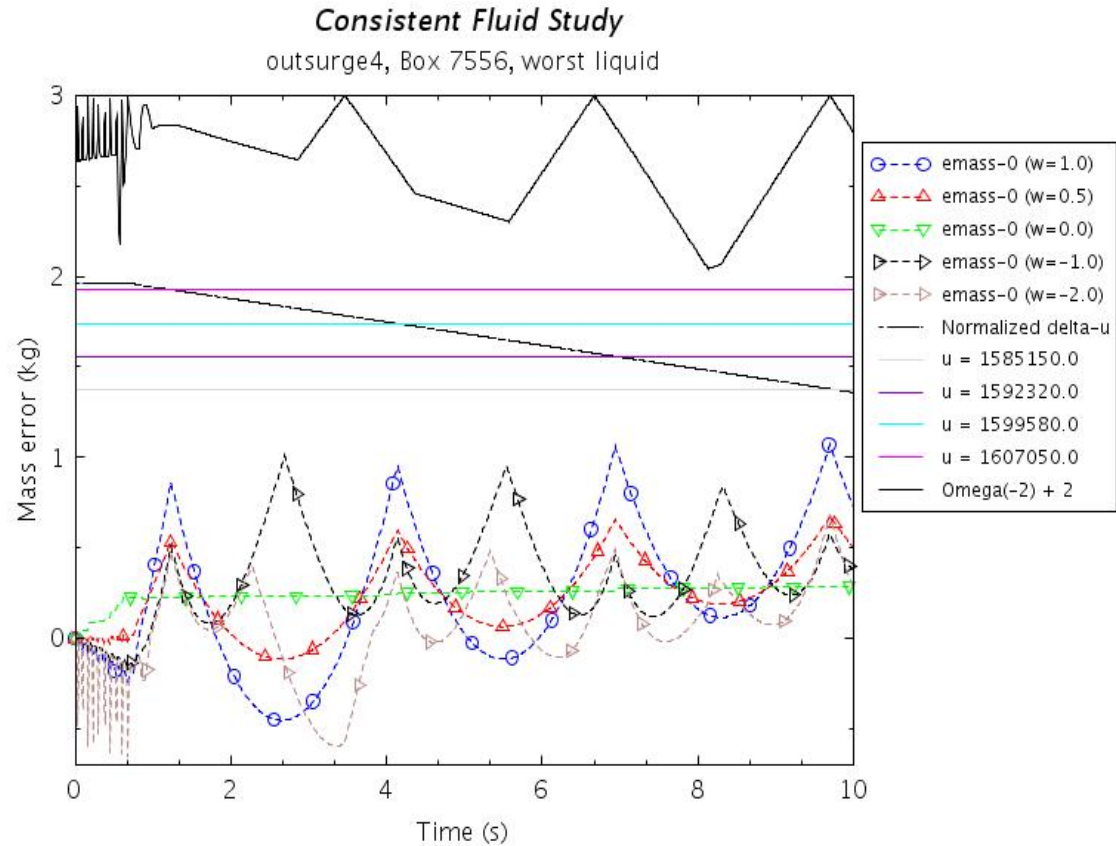
OUTSURGE Case, Box 7556 Mass error

- Identical noding diagram of the insurge and outsurge problems
- Outsurge case purpose: put the pressurizer volume through a blowdown, causing a phase boundary crossing
- Flow is reversed
- Horizontal lines show normalized U grid points
- Funnel ($\omega = -2$) values graphed at $\omega + 2$ for clarity
- Crossing normalized P-lines causes mass err oscillations
- Local extrema at or near P-line crossings for all ω
- $\omega=0$ mass error is **nonzero**



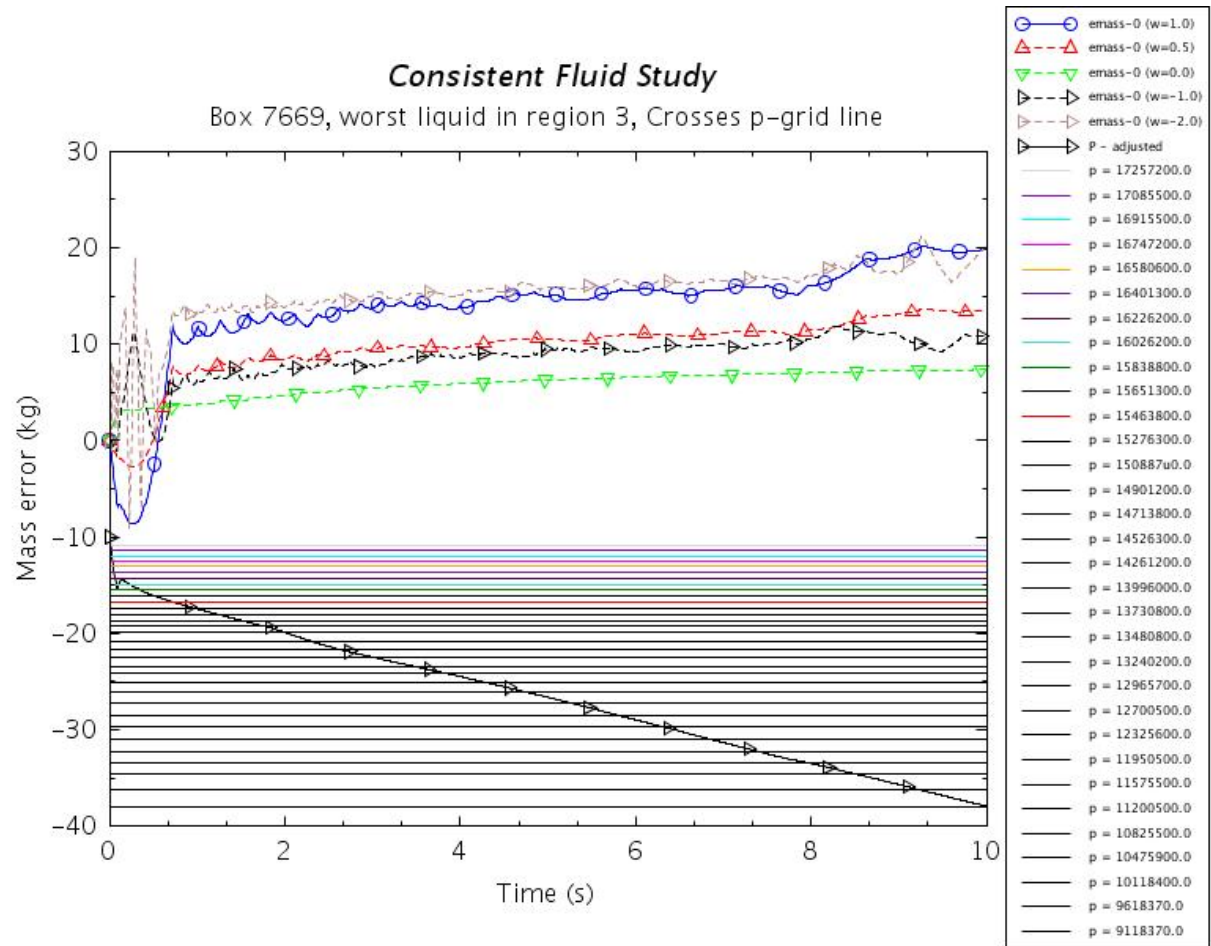
OUTSURGE Case, Box 7556 Mass error

- Local maxima at U-line crossings for all ω
- Inflection points in all graphs where funnel- ω either equals 1 or has a kink.



OUTSURGE Case, Box 7669 Mass error

- Oscillations at P-line crossings
- Again, $\omega=0.0$ is best but has nonzero mass error

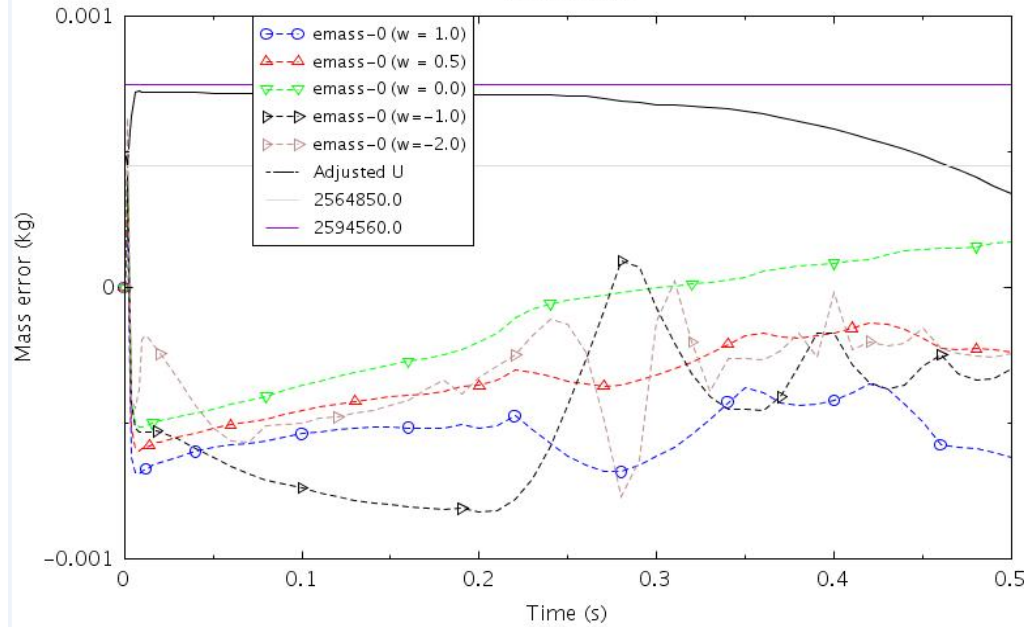
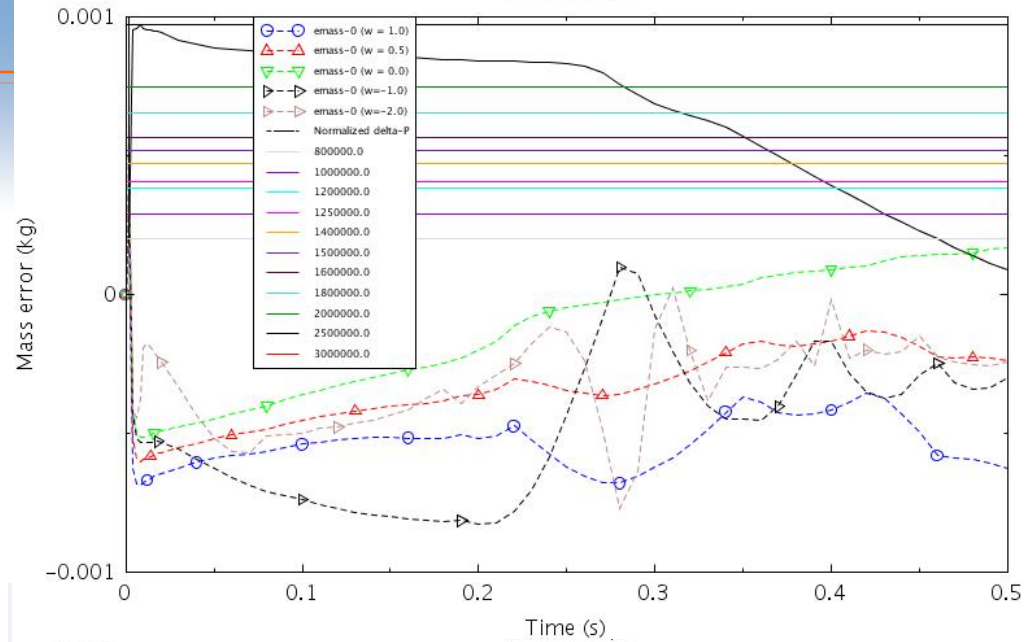


OUTSURGE Case, Box 7556 Mass error

- As with insurge, the plot of $\omega(-2)$ exhibits oscillatory behavior and has local extrema at pressure gridline crossings
- Plots of $\omega=1.0$ and $\omega=0.5$ have local maxima where the specific internal energy crosses gridlines
- Smallest mass error is created when $\omega=0.0$. However, unlike with insurge, the mass error is positive in outsurge problems
- Second lowest mass error for Box 7556 comes when $\omega=-2$
- Second lowest mass error for Box 7669 comes when $\omega=-1$

Edward's Mass Error

- Many local maxima and minima.
- Pressure grid point crossings and internal energy grid point crossing corresponds to an inflection point on the mass error curves
- From about 0.02s to 0.2s, where neither pressure nor energy grid points are crossed, the mass error plots are relatively smooth
- Smallest mass error for $\omega = 0$



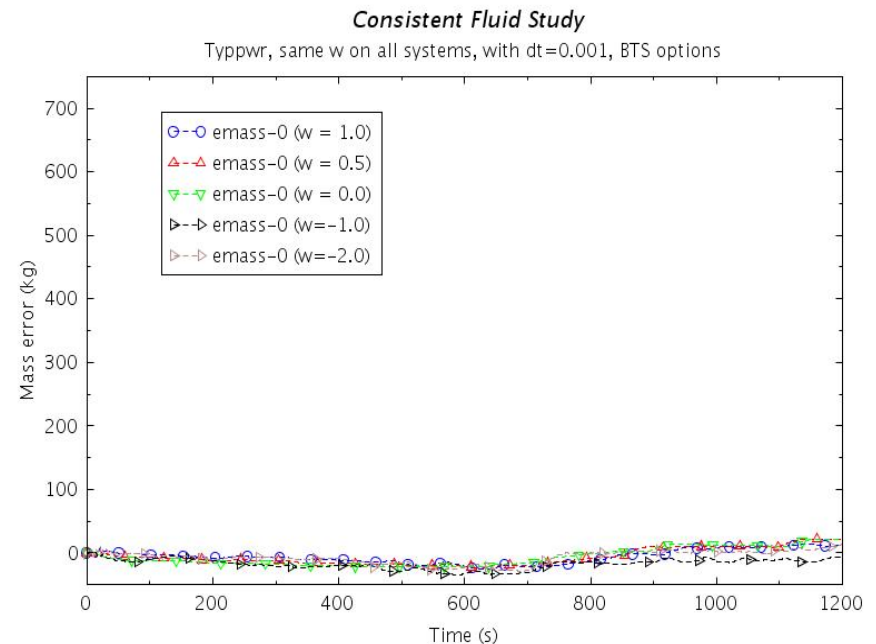
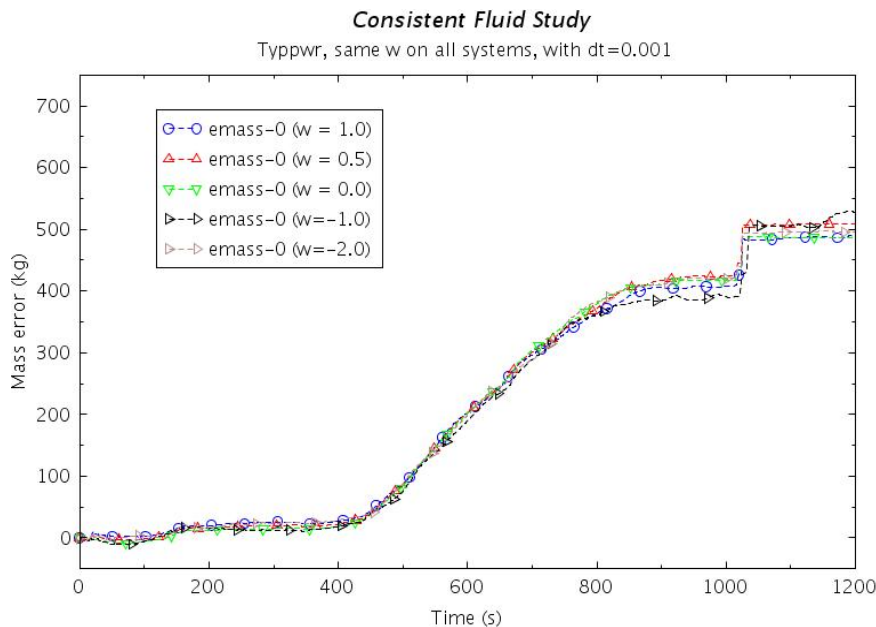
TYPWR Test Cases

- The typical PWR model has three systems
 - Same values of ω applied to each system
- Fifty code runs: Combinations of 5 ω -values (1.0, 0.5, 0.0, -1, -2) timestep values ($\Delta t = 0.1, 0.05, 0.01, 0.005, 0.001$) and 2 sets of card-one options: default and “Better Test Set” (BTS)
- Default and BTS produced similar results on earlier calculations

| Card-1 | “Better Test Set” of Options |
|--------|---|
| 29 | Accurately solves the momentum equations at low velocities. |
| 41 | Includes energy dissipation due to form loss (code calculated abrupt area change loss and user-specified loss) |
| 54 | Changes the 2-phase to 1-phase gas transition truncation limit in EQFINL for the semi-implicit |
| 55 | Model improvements to minimize numerical sources of oscillations for low pressure 2-phase flow: <ul style="list-style-type: none"> • Interfacial heat transfer for annular mist, Mist pre-CHF, Mist post-CHF flow regimes, More physical Hif and Hig |

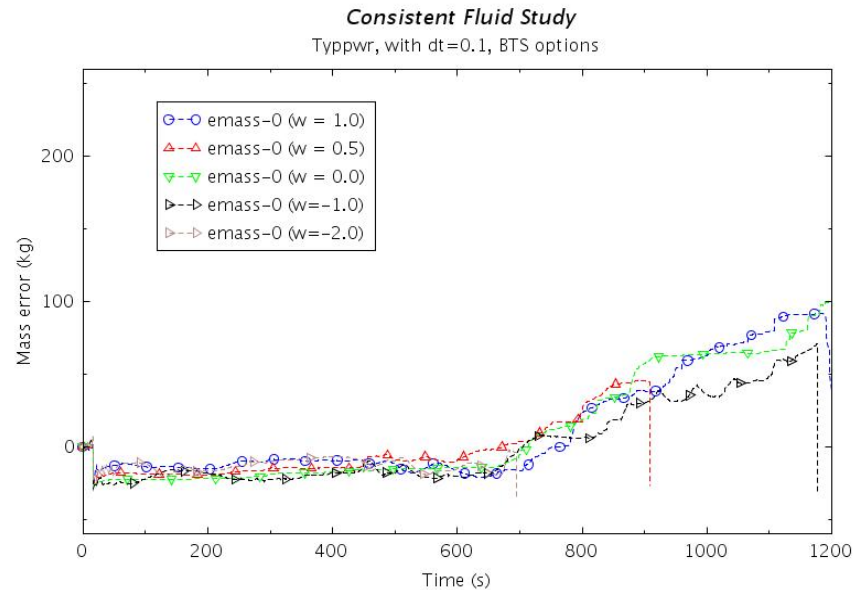
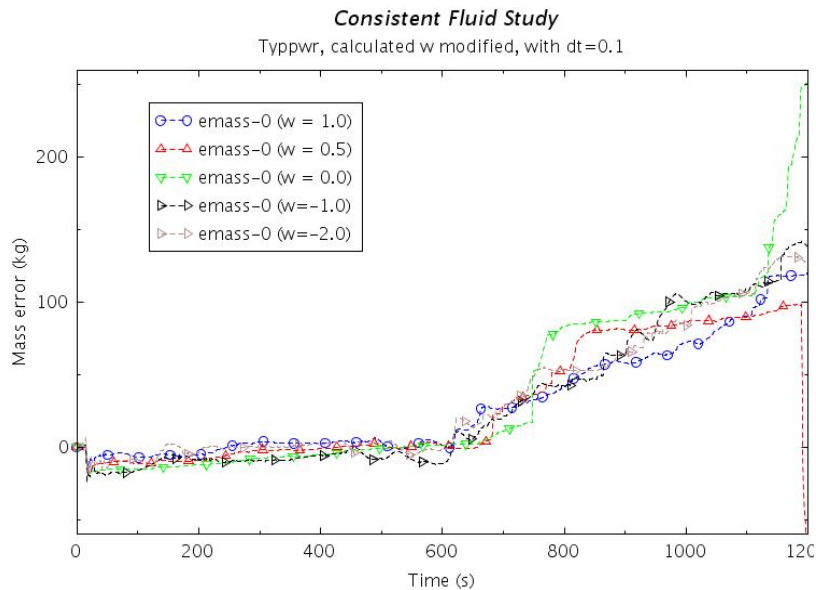
TYPWPW Mass Error, $\Delta t = 0.001$

- $\omega = -1$ has smallest mass error with BTS. $\omega = 0.0$ is best with default
- At smaller time step, BTS mass error is vastly superior to the default values of the card- options
- Why not make BTS the default?



TYPWPWR Mass Error, $\Delta t = 0.1$

- Except $\omega=1.0$ and $\omega=0.0$, all ω -values **FAIL** with **BTS** by 1200s
- All ω -values run with default options
- $\omega=0.0$ has the worst mass error, more than double any other, with the default options.



TYPWR Test Cases

- Runtime comparison
 - Orange is slowest
 - Yellow is fastest
 - Def = default
 - Ratio = failed advancements divided by total for Default options
- Ratio and runtime should correlate, but not well. Single run
- No option is best or worst for TYPWR

| ω | dt=0.5 | | | dt=0.01 | | |
|----------|--------|--------|--------|---------|--------|--------|
| | Ratio | Def | BTS | Ratio | Def | BTS |
| 1.0 | 0.235 | 38.148 | 36.552 | 0.035 | 163.78 | 166.70 |
| 0.5 | 0.194 | 37.089 | 37.490 | 0.028 | 161.27 | 163.87 |
| 0.0 | 0.230 | 37.791 | 38.465 | 0.042 | 164.99 | 165.38 |
| -1 | 0.174 | 37.033 | 38.344 | 0.045 | 164.85 | 162.29 |
| -2 | 0.196 | 37.679 | 37.148 | 0.046 | 164.26 | 165.19 |

| ω | dt=0.05 | | | dt=0.001 | | |
|----------|---------|--------|--------|----------|--------|--------|
| | Ratio | Def | BTS | Ratio | Def | BTS |
| 1.0 | 0.0262 | 322.62 | 323.07 | 0.00871 | 1586.4 | 1586.4 |
| 0.5 | 0.0253 | 319.88 | 321.44 | 0.00743 | 1573.0 | 1573.2 |
| 0.0 | 0.0242 | 324.12 | 321.31 | 0.00804 | 1569.7 | 1564.3 |
| -1 | 0.0192 | 320.21 | 322.82 | 0.00816 | 1572.2 | 1573.0 |
| -2 | 0.0205 | 321.49 | 321.54 | 0.00833 | 1575.7 | 1578.1 |

Moby Dick Problem Conclusions

- Four types of advancement
 - 1 = semi-implicit, explicit coupling with heat transfer
 - 2 = semi-implicit, implicit coupling with heat transfer
 - 3 = nearly-implicit, explicit coupling with heat transfer
 - 4 = nearly-implicit, implicit coupling with heat transfer
- Mass error ratio
 - Default options: constant for some, not for others
 - Default options: smaller mass error for case 1.
 - BTS options: lower mass error ratios than default for other cases
- The DA graph of pressure vs. elevation shows no visible differences between the five values of omega.
- *The mass error with H2ON was typically 1.5 times lower than the mass error ratio produced with default H2O.*

Conclusions

- No particular value of ω produced consistently better results in terms of mass error, though:
 - $\omega = -1.0$ somewhat outperformed other choices
 - $\omega = 0.0$ performed well
- Based on these limited results, ω did not visibly affect engineering parameters such as pressure and void fraction.
- Crossing pressure and internal energy grid lines from one time advancement to the next affects mass error
 - In simple cases, pressure crossings cause local extrema and internal energy crossing inflect the curves.
 - The effect diminishes as ω tends toward zero, but is visible even for $\omega = 0.0$ for some problems.

Conclusions and Recommendations

- With the BTS options, TH property failures occur with Typical PWR for all omega values at the largest DTMAX
- The default value of $\omega = 1.0$ was seldom the best and sometimes the worst in terms of mass error and runtime
 - The user may replace it by another choice through input
- Value $\omega = 0.0$ was:
 - Uniformly better for the simple insurge problems
 - Was generally better for the outsurge problems
 - Was sometimes worst with more complicated models
- Value $\omega = -1.0$ seldom performed poorly and performed well or the best in a significant number of more complicated cases
- Further study is recommended
- Cubic splines to smooth crossing should improve performance