# Jacobian-Free Newton-Krylov (JFNK) Methods for Nonlinear Neutronics/Thermal-Hydraulic Equations

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### **Outline**

- 1 Introduction
- 2 Governing Equations
- 3 Solvers
- 4 Results
- 5 Conclusions



### Motivation

■ Eventually will be part of thesis work

■ JFNK method not currently used in nuclear reactor analysis

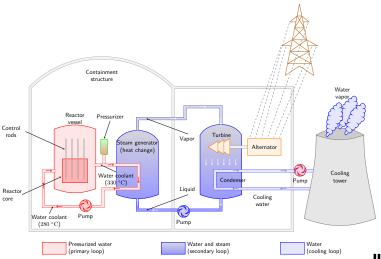
■ Incorporates a lot of ideas from 2.29 class

Coupled physics is fun!



Introduction Governing Equations Solvers Results Conclusions

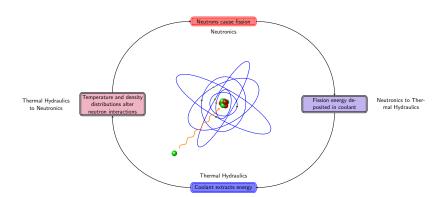
### **Nuclear Reactor Plant**





Introduction Governing Equations Solvers Results Conclusion

### **Nuclear Feedback**

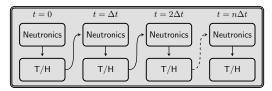


- lacktriangle Fuel Temperature Feedback  $T_f \uparrow$ , U-238 Capture  $\uparrow$ , Fission Rate  $\downarrow$ , Power  $\downarrow$
- Coolant Density Feedback  $\rho \downarrow$ ,  $E_n \uparrow$ , Fission Rate  $\downarrow$ , Power  $\downarrow$

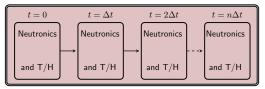


## Common Approach to Coupling: Operator Splitting

Solve physics independently and iterate between them



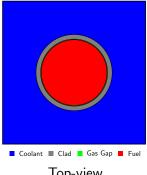
Fully coupled approach solves the nonlinear physics together





Introduction Governing Equations Solvers Results Conclusion

### 1-D Slab Reactor Geometry



Top-view
Fuel Rod Unit-Cell



Side-view (vertical)
1-D Model of Reactor



Basic Neutron Conservation:

$$Change + Leakage + Interactions = Scattering + Fission \\$$

**Neutron Transport Equation:** 

$$\underbrace{\frac{1}{v}\frac{\partial\varphi}{\partial t}}_{\text{time-dependent}} + \underbrace{\mathbf{\Omega}\cdot\nabla\varphi\left(\mathbf{r},E,\mathbf{\Omega},t\right)}_{\text{neutron leakage}} + \underbrace{\sum_{t}\left(\mathbf{r},E,t\right)\varphi\left(\mathbf{r},E,\mathbf{\Omega},t\right)}_{\text{interaction of neutrons with medium}} = \underbrace{Q\left(\mathbf{r},E,\mathbf{\Omega},t\right)}_{\text{neutron source}}$$

Neutron Diffusion Equation (1-D Energy Integrated)

$$\frac{1}{v}\frac{\partial\phi}{\partial t} - D\left(x,t\right)\frac{\partial^{2}\phi}{\partial^{2}x} + \Sigma_{a}\left(x,t\right)\phi\left(x,t\right) = \frac{1-\beta}{k_{eff}}\nu\Sigma_{f}\left(x,t\right)\phi\left(x,t\right) + \lambda_{d}c\left(x,t\right)$$

$$\frac{\partial c}{\partial t} = \frac{\beta}{k_{eff}}\nu\Sigma_{f}\left(x,t\right)\phi\left(x,t\right) - \lambda_{d}c\left(x,t\right)$$



# **Discretization of Neutronics Equations**

#### Assumptions:

- 1 One-dimensional finite volume spatial discretization, uniform  $\Delta x$
- 2 Central difference scheme for diffusion term
- 3 No incoming current of neutrons at boundaries
- 4 Implicit Euler time discretization
- Discretized neutronics equation for interior cell

$$\begin{split} \frac{1}{v} \frac{d\bar{\phi}_i}{dt} - \frac{2}{\Delta x^2} \frac{D_i D_{i-1}}{D_i + D_{i-1}} \bar{\phi}_{i-1} + \left( \frac{2}{\Delta x^2} \frac{D_{i+1} D_i}{D_{i+1} + D_i} + \frac{2}{\Delta x^2} \frac{D_i D_{i-1}}{D_i + D_{i-1}} + \Sigma_{a,i} \right) \bar{\phi}_i - \\ \frac{2}{\Delta x^2} \frac{D_{i+1} D_i}{D_{i+1} + D_i} \bar{\phi}_{i+1} = \frac{1 - \beta}{k_{eff}} \nu \Sigma_{f,i} \bar{\phi}_i + \lambda_d \bar{c}_i \end{split}$$

Matrix-form of neutronics equations

$$\bar{\mathbf{\Phi}}^{n+1} - \bar{\mathbf{\Phi}}^n + v\Delta t \left( \mathbb{M}\bar{\mathbf{\Phi}}^{n+1} - (1-\beta) \lambda \mathbb{F}\bar{\mathbf{\Phi}}^{n+1} - \lambda_d \bar{\mathbf{c}}^{n+1} \right) = 0$$

Matrix-form of precursors

$$\bar{\mathbf{c}}^{n+1} - \bar{\mathbf{c}}^n + \Delta t \left( \lambda_d \bar{\mathbf{c}}^{n+1} - \beta \lambda \mathbb{F} \bar{\mathbf{\Phi}}^{n+1} \right) = 0$$



Energy Equation - single phase fluid and inviscid fluid

$$\frac{\partial (\rho h)}{\partial t} + \nabla \cdot (\rho h \mathbf{u}) = -\nabla \cdot \mathbf{q}'' + q'''$$

Assuming fissions are a volumetic heat source in 1-D

$$\rho A \frac{\partial h}{\partial t} + \dot{m} \frac{\partial h}{\partial x} = q'$$

For an incompress2cmible fluid,  $dh=c_pdT$ 

$$\rho A c_p \frac{\partial T}{\partial t} + \dot{m} c_p \frac{\partial T}{\partial x} = q'$$



#### Assumptions:

- 1 One-dimensional finite volume spatial discretization, uniform  $\Delta x$
- 2 Upwind difference scheme for flux
- 3 Specify inlet conditions and mass flow rate
- 4 Implicit Euler time discretization
- Spatial discretization

$$\frac{\rho A \Delta x}{\dot{m}} \frac{d\bar{T}_i}{dt} + \bar{T}_i - \bar{T}_{i-1} = \frac{1}{2\dot{m}c_p} \left( Q_{i-1} + Q_i \right)$$

Matrix-form with time discretization

$$\bar{\mathbf{T}}^{n+1} - \bar{\mathbf{T}}^n + \frac{\dot{m}\Delta t}{\mathcal{P}^{n+1}A\Delta x} \left( \mathbb{S}\bar{\mathbf{T}}^{n+1} - \mathbb{R}\mathbf{Q}^{n+1} \right) = 0$$

Note:  $\mathcal{P}$  is a vector of cell-averaged coolant densities



## **Physics Coupling**

### Neutronics-Thermal Hydraulics

- Neutrons cause fission
- Large portion of fission energy deposited in coolant
- This is represented by

$$\mathbf{Q} = \tilde{c} \mathbb{E} \bar{\mathbf{\Phi}} \Delta x$$

where  $\mathbb{E}=\mathrm{diag}\left\{\kappa\Sigma_f\right\}$  characterizes energy per fission and  $\tilde{c}$  is the flux-to-power normalization constant

#### Thermal Hydraulics-Neutronics

- Diffusion theory parameters depend on coolant density
- This dependence is determined with a transport theory code
- D,  $\Sigma_a$ ,  $\nu\Sigma_f$ ,  $\kappa\Sigma_f$  are all affected by coolant density variations
- Data is correlated with a linear regression of the form:

$$\Sigma = \Sigma^{ref} + \frac{\partial \Sigma}{\partial \rho} \left( \mathcal{P} - \rho^{ref} \right)$$

achusett ute of nology

# The Steady State Eigenvalue Problem

- The steady state equations must be solved first
- Reducing the neutronics equation to steady state form:

$$\mathbb{M}ar{m{\Phi}} = \lambda \mathbb{F}ar{m{\Phi}}, \qquad \lambda = rac{1}{k_{eff}}$$

- lacktriangle Eigenvalue,  $\lambda$ , and eigenvector,  $\Phi$ , must be determined
- Flux-to-power normalization constant determined from reactor power:

$$Q_{R} = \tilde{c} \int_{0}^{L} dx \kappa \Sigma_{f}(x) \phi(x) = \tilde{c} \sum_{i} \kappa \Sigma_{f,i} \bar{\phi}_{i} \Delta x = \tilde{c} \kappa \Sigma_{f}^{\mathrm{T}} \bar{\Phi} \Delta x$$

 $m{\lambda}$  and  $\tilde{c}$  are specified as constants for time-dependent calculations



### **Newton's Method**

### Procedure:

9:

```
1: Goal: \mathbf{F}(\mathbf{x}) = \mathbf{0}

2: Guess \mathbf{x}

3: for n = 1, 2, 3, ... do

4: \mathbf{r} = \mathbf{F}(\mathbf{x})

5: if \|\mathbf{r}\| < ntol then

6: DONE!

7: end if

8: d\mathbf{x} = -\mathbb{J}^{-1}(\mathbf{x}) \mathbf{F}(\mathbf{x})
```

x = x + dx

Three major tasks:

10: end for

- 1 Evaluate residual vector in external function
- 2 Evalue Jacobian in external function (Do we have to?)
- 3 Calculate dx Direct or Iterative Solvers?



# **Krylov Subspace Methods**

- A class of iterative methods for sparse systems
- Solves  $A\mathbf{x} = \mathbf{b}$  by projecting a m dimensional problem into a lower dimensional Krylov subspace

$$\mathcal{K}_n(\mathbb{A}, \mathbf{v}) = \operatorname{span} \{ \mathbf{v}, \mathbb{A}\mathbf{v}, \mathbb{A}^2\mathbf{v}, ..., \mathbb{A}^{n-1}\mathbf{v} \}$$

- Here A is nonhermitian and we use the GMRES method
- GMRES uses the Arnoldi method to reduce the system to Hessenberg form

$$\mathbb{AQ} = \mathbb{QH}$$

$$\mathbb{H} = \begin{bmatrix} h_{11} & \cdots & h_{1n} \\ h_{21} & h_{22} & & \vdots \\ & \ddots & \ddots & \vdots \\ & & h_{n,n-1} & h_{n,n} \end{bmatrix}$$



### Generalized Minimal RESidual Method

- Goal:  $\mathbf{x}_* = \mathbb{A}^{-1}\mathbf{b}$
- A step n,  $\mathbf{x}_*$  is approximated by  $\mathbf{x}_n \in \mathcal{K}_n$  that minimizes the norm of the residual  $\mathbf{r}_n = \mathbf{b} - \mathbb{A}\mathbf{x}_n$

#### Procedure:

- 1:  $q_1 = b/\|b\|$
- 2: for n = 1, 2, 3, ... do
- 3: Perform step n of Arnoldi (Creates Hessenberg matrix)
- Find y to minimize  $\left\|\widetilde{\mathbb{H}}_n\mathbf{y} \|\mathbf{b}\|\,\mathbf{e}_1\right\|$ 4:
- 5: if  $\|\mathbf{r}\| < ltol$  then
- 6: DONE
- 7: end if
- 8: end for
- 9:  $\mathbf{x} = \mathbb{O}_n \mathbf{v}$
- Saad et al.  $^{\ddagger}$  defines a novel method to compute  $\widetilde{\mathbb{H}}_n$  from  $\widetilde{\mathbb{H}}_{n-1}$  from Givens rotations

<sup>&</sup>lt;sup>‡</sup>Youcef Saad and Martin H. Schultz. GMRES: A generalized minimal residual algorithm for solving nonsymmetric linear systems. Society for Industrial and Applied Mathematics, 7:856-859, 1986.



### **Inexact Newton's Method**

- Newton-Krylov is an inexact Newton method since the linear step is not determined exactly
- In Newton-Krylov framework, two tolerances were defined:
  - 1 Nonlinear tolerance for Newton iteration
  - 2 Linear tolerance for GMRES iteration
- Why have tight linear convergence when nonlinear residual is large?
- Instead, a relative residual tolerance,  $\eta$ , is used

$$\|\mathbb{J}(\mathbf{x}^n) d\mathbf{x}_m^n + \mathbf{F}(\mathbf{x}^n)\| < \eta \|\mathbf{F}(\mathbf{x}^n)\|$$

- At initial Newton iterations, GMRES will not be converged very tightly
- For the last couple of Newton iterations, convergence may be too tight ∴ limit how small linear tolerance can get



# **Jacobian-Free Approximation**

- $\blacksquare \quad \mathsf{Recall \ a \ Krylov \ subspace:} \ \mathcal{K}_n \left( \mathbb{A}, \mathbf{v} \right) = \mathrm{span} \left\{ \mathbf{v}, \mathbb{A}\mathbf{v}, \mathbb{A}^2\mathbf{v}, ..., \mathbb{A}^{n-1}\mathbf{v} \right\}$
- $\blacksquare$  Why create  $\mathbb A$  when it is only used to multiply a vector?
- Option 1: Perform Jacobian-vector product analytically

$$\mathbb{J}\mathbf{y} = \left[ \begin{array}{cc} \mathbb{M} - \lambda \mathbb{F} & -\mathbb{F}\bar{\mathbf{\Phi}} \\ -\bar{\mathbf{\Phi}}^\top & 0 \end{array} \right] \left[ \begin{array}{c} y_\phi \\ y_\lambda \end{array} \right] = \left[ \begin{array}{cc} (\mathbb{M} - \lambda \mathbb{F}) \, y_\phi - \mathbb{F}\bar{\mathbf{\Phi}} y_\lambda \\ -\bar{\mathbf{\Phi}}^\top y_\phi \end{array} \right]$$

Option 2: Approximate Jacobian-vector product with finite difference

$$\mathbb{J}\mathbf{y} \approx \frac{\mathbf{F}\left(\mathbf{x} + \epsilon\mathbf{y}\right) - \mathbf{F}\left(\mathbf{x}\right)}{\epsilon}$$

- Advantages: Saves memory and possibly computational time to form Jacobian
- lacktriangledown is the perturbation parameter and is somewhat arbitrary Mousseau  $\S$  recommends:

$$\epsilon = \frac{\sum_{i=1}^{N} bx_i}{N \|\mathbf{y}\|_2} \qquad b = 1 \times 10^{-8}$$

<sup>§</sup>V.A. Mousseau. Implicitly balanced solution of the two-phase flow equations couple to nonlinear heat IIII Massachusell conduction. Journal of Computational Physics, 200:104-132, 2004.

# Want to limit number of GMRES iterations

Before a calculation, a Jacobian matrix is formed analytically and a zero-fill Incomplete LU (ILU) is performed:

$$\mathbb{R} = \mathbb{L}\mathbb{U} - \mathbb{A}$$

- In ILU, residual matrix  $\mathbb R$  is constrained to certain conditions
- Zero-fill implies that the number and location of nonzeros is preserved
- Left preconditioning is used in this project:

$$\mathbb{U}^{-1}\mathbb{L}^{-1}\mathbb{A}\mathbf{x} = \mathbb{U}^{-1}\mathbb{L}^{-1}\mathbf{b}$$



# Steady State - Neutronics

Residual Equations

$$\mathbf{F} = \left[ egin{array}{c} \mathbb{M}\mathbf{\Phi} - \lambda \mathbb{F}\mathbf{\Phi} \ -rac{1}{2}ar{\mathbf{\Phi}}^{ op}ar{\mathbf{\Phi}} + rac{1}{2} \end{array} 
ight]$$

■ Resulting neutron flux distribution:

- A Newton method cannot guarantee that the fundamental eigenmode is calculated
- Any mode satisfies the nonlinear set of equations
- Here the ratio of the first two eigenvalues is close to unity (dominance ratio)
- Use a few power iterations to get gross flux shape



# Steady State - Coupled Neutronics/Thermal Hydraulics [1/2]

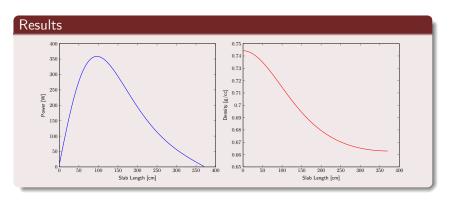
#### Residual Equations

$$\mathbf{F} = \begin{bmatrix} \mathbf{M}\boldsymbol{\Phi} - \lambda \mathbb{F}\boldsymbol{\Phi} \\ Q_R - \tilde{c}\kappa \boldsymbol{\Sigma}_f^{\mathrm{T}}\boldsymbol{\Phi}\Delta x. \\ \mathbf{Q} - \tilde{c}\mathbf{E}\boldsymbol{\Phi}\Delta x \\ \mathbf{S}\mathbf{T} - \mathbb{R}\mathbf{Q} \\ \mathcal{F} - \rho\left(\mathbf{T}, p\right) \\ \boldsymbol{\Sigma}_a - \boldsymbol{\Sigma}_a^{ref} - \frac{\partial \boldsymbol{\Sigma}_a}{\partial \rho} \left[\mathcal{P} - \rho^{ref}\right] \\ \nu \boldsymbol{\Sigma}_f - \nu \boldsymbol{\Sigma}_f^{ref} - \frac{\partial \nu \boldsymbol{\Sigma}_f}{\partial \rho} \left[\mathcal{P} - \rho^{ref}\right] \\ \mathbf{D} - D^{ref} - \frac{\partial \boldsymbol{D}}{\partial \rho} \left[\mathcal{P} - \rho^{ref}\right], \\ \kappa \boldsymbol{\Sigma}_f - \kappa \boldsymbol{\Sigma}_f^{ref} - \frac{\delta \kappa \boldsymbol{\Sigma}_f}{\partial \rho} \left[\mathcal{P} - \rho^{ref}\right] \\ - \frac{1}{2}\boldsymbol{\Phi}^{\mathrm{T}}\boldsymbol{\Phi} + \frac{1}{2} \end{bmatrix}$$

- Neutronics
- Flux normalization
- Energy deposition
- Temperature distribution
- Density distribution
- Absorption xs
- Fission xs
  - Diffusion
- Energy deposition xs
- Eigenvalue
- X-Steam MATLAB tables were used to look up density as a function of temperature
- An analytical Jacobian-vector product can be formed for everything except  $\frac{\partial \rho}{\partial T}$
- For this block a finite difference approximation must be used



# Steady State - Coupled Neutronics/Thermal Hydraulics [2/2]



- Timing results using MATLAB profiler:
  - 7 seconds to converged overall
  - 6+ out of 7 seconds in X-Steam!
  - Solution: fit density range with polynomial with linear fit < 0.5 seconds



# Transient - Coupled Neutronics/Thermal Hydraulics [1/2]

#### Residual Equations

$$\mathbf{F} = \begin{bmatrix} \mathbf{\Phi}^{n+1} - \mathbf{\Phi}^n + v\Delta t \left[ \mathbb{M}\mathbf{\Phi}^{n+1} - (1-\beta) \lambda \mathbb{F}\mathbf{\Phi}^{n+1} - \lambda_d \mathbf{c}^{n+1} \right] \\ \mathbf{c}^{n+1} - \mathbf{c}^n + \Delta t \left( \lambda_d \mathbf{c}^{n+1} - \beta \lambda \mathbb{F}\mathbf{\Phi}^{n+1} \right) \\ \mathbf{Q} - \tilde{c}\mathbb{E}\mathbf{\Phi}\Delta x \\ \mathbf{T}^{n+1} - \mathbf{T}^n + \frac{\tilde{m}\Delta t}{\mathcal{P}^{n+1}A\Delta x} \left( \mathbb{S}\mathbf{T}^{n+1} - \mathbb{R}\mathbf{Q}^{n+1} \right) \\ \mathcal{P} - \rho^{ref} - \frac{\partial \mathcal{P}}{\partial \mathcal{P}} \left( \mathbf{T} - T^{ref} \right) \\ \mathbf{\Sigma}_a - \mathbf{\Sigma}_a^{ref} - \frac{\partial \mathbf{\Sigma}_a}{\partial \rho} \left[ \mathcal{P} - \rho^{ref} \right] \\ \nu \mathbf{\Sigma}_f - \nu \mathbf{\Sigma}_f^{ref} - \frac{\partial \nu \mathbf{\Sigma}_f}{\partial \rho} \left[ \mathcal{P} - \rho^{ref} \right] \\ \mathbf{D} - D^{ref} - \frac{\partial \mathbf{D}}{\partial \rho} \left[ \mathcal{P} - \rho^{ref} \right] \\ \kappa \mathbf{\Sigma}_f - \kappa \mathbf{\Sigma}_f^{ref} - \frac{\partial \kappa \mathbf{\Sigma}_f}{\partial \rho} \left[ \mathcal{P} - \rho^{ref} \right] \end{bmatrix}$$

- Code time per loop
- Number of GMRES and Newton





### Conclusions

■ JFNK methods are well suited for solving coupled systems

■ Finite difference method for Jacobian-vector product works well

Write your own steam tables or fit data with polynomial

Watch out for scaling issues when coupling physics



- Rewrite code in a compiled language (Fortran)
- Use PETSc Nonlinear solvers
- Model other important nuclear feedback mechanisms
- Compare JFNK to Operator-splitting methods
- This project can be downloaded from GitHub:

www.github.com/bhermanmit/JFNK
git clone git@github.com:bhermanmit/JFNK

