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***Flux and Power Mapping in
RFSP***

Benoit Arsenault, AECL

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Flux and Power Mapping in RFSP

- **Fuel Management Program RFSP has FLUX MAPPING and POWER MAPPING capability**
- **Alternative to unusual RFSP method of solving the finite-difference diffusion equation in 3 dimensions**
- **Mapping used to monitor core power distribution under nominal equilibrium core conditions at CANDU 6 sites (Gentilly 2 and Point Lepreau)**
- **Advantage over diffusion-type calculation is direct inclusion of in-core data in the power calculation**



Calculations of the Harmonics (*MONIC)

Steady State Diffusion Equation: $(R - P)\phi = 0$

where R is the Removal matrix $\begin{bmatrix} \nabla \cdot D_1 \nabla - (\Sigma_{a,1} + \Sigma_m) & \nabla \cdot D_2 \nabla - \Sigma_{a,2} \\ \Sigma_m & 0 \end{bmatrix}$

P is the Production matrix $\begin{bmatrix} 0 & v \Sigma_{f,2} \\ 0 & 0 \end{bmatrix}$

is the Flux vector $\begin{bmatrix} \phi_1 \\ \phi_2 \end{bmatrix}$

Objective: To find eigenfunctions of $R = \frac{1}{\lambda_i} P \phi_i$

At n^{th} iteration $\phi^n = R^{-1} P \phi^{n-1} = R^{-1} P \Sigma_i A_i \phi_i = \Sigma_i \lambda_i A_i \phi_i$

The eigenfunction with the largest λ_i will emerge and dominates.
Solution converges to this predominate mode.



Calculations of the Harmonics (con't)

The Adjoint Flux Vector $\phi^* = (\phi_1^* \ \phi_2^*)$ satisfy the $\phi^*(R-P) = 0$

Bi-Orthogonality of the natural modes:

$$\int (\phi_{1M}^*(\vec{r}) \phi_{2M}^*(\vec{r})) P \begin{bmatrix} \phi_{1N}(\vec{r}) \\ \phi_{2N}(\vec{r}) \end{bmatrix} d\vec{r} = 0$$

for any two different harmonics M .NE. N

For a pseudo-one-group flux $\phi_T = \phi_1 + \phi_2$, it is

self-adjoint:

$$\int \phi_{TM}(\vec{r}) v \Sigma_f(\vec{r}) \phi_{TN}(\vec{r}) d\vec{r} = 0 \quad M \neq N$$



Calculations of the Harmonics (con't)

Calculation of the Nth harmonic mode involves subtracting off from the unconverged flux the components of the previous harmonics:

$$\phi'_{uc} = \phi_{uc} - \sum_{I=1}^{N-1} A_I \phi_I$$

The Amplitude A_I of the component of the Ith harmonic determined using the approximate orthogonal property of the total flux:

$$A_I = \frac{\int \phi_{TI}(\vec{r}) \cdot \nabla \Sigma_f(\vec{r}) \phi_{Tuc}(\vec{r}) d\vec{r}}{\int \phi_{TI}(\vec{r}) \cdot \nabla \Sigma_f(\vec{r}) \phi_{TI}(\vec{r}) d\vec{r}} \quad I=1, \dots, N-1$$



Calculations of the Harmonics (con't)

- **A repetitive “Iterate - Subtraction” procedure forces convergence to the next higher harmonic. Harmonics generated are orthogonal (Gram-Scmidt Orthogonalization Procedure).**

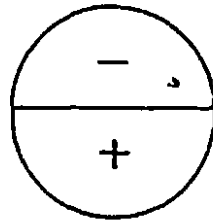


Selection of Mode Set

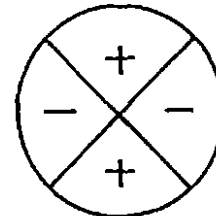
During normal full-power operation the set of modes consists of a fundamental mode (based on a recent core diffusion calculation) and the first 10-14 harmonic modes.

For the normal simulation e.g., a derating during which adjuster banks are withdrawn, a set of 22 flux modes is used:

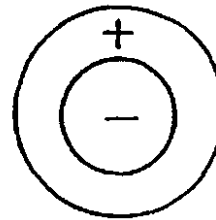
- a. fundamental based on diffusion calculation of core state before derating**
- b. 14 harmonic modes**
- c. 7 power recovery modes with 1 through 7 adjuster banks removed from core**



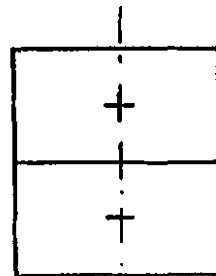
FIRST
AZIMUTHAL



SECONO
AZIMUTHAL



FIRST RADIAL



FIRST AXIAL



FLUX HARMONICS

MODE NUMBER	DESCRIPTION	ORDER/ALTY or	MODE SCHEMATIC (IDEALIZED)
0	FUNDAMENTAL	0	
1	FIRST AZIMUTHAL-A	01.1	
2	FIRST AZIMUTHAL-B	01.2	
3	FIRST AXIAL	02.1	
4	SECOND AZIMUTHAL-A	04.0	
5	SECOND AZIMUTHAL-B	04.0	
6	FIRST AZIMUTHAL-A X FIRST AXIAL	06.0	
7	FIRST AZIMUTHAL-B X FIRST AXIAL	06.0	
8	FIRST RADIAL X SECOND AXIAL-A	08.1	
9	FIRST RADIAL X SECOND AXIAL-B	08.1	

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Harmonics for CANDU (Cylindrical Reactor)

- For homogeneous bare cylindrical reactor, flux shape given by (in r, θ, z co-ordinate):

$$\phi(r, \theta, z) = J_M(\alpha_{ML} \cdot r / R_0) \cdot \cos(M \cdot \theta) \cdot \sin(N \cdot \pi \cdot z / H) \quad M \text{ even}$$

$$\phi(r, \theta, z) = J_M(\alpha_{ML} \cdot r / R_0) \cdot \sin(M \cdot \theta) \cdot \sin(N \cdot \pi \cdot z / H) \quad M \text{ odd}$$

where J_M is the M^{th} order Bessel function,

α_{ML} is the L^{th} zero of J_M

R_0 is the radius of the reactor

H is the height of the reactor

- Various combinations of M and N give the Harmonics. Flux shape used as initial guess in *MONIC



Harmonics - Natural Modes (Example)

1-D Problem Slab reactor, thickness from $x = -a/2$ to $+a/2$

$$\frac{d^2\phi}{dx^2} + B^2\phi = 0$$

Ignore flux extrapolation beyond slab surface,

i.e. assume $\phi = 0$ at $x = a/2$ and $-a/2$

Note also symmetry $\phi(x) = \phi(-x)$ and $\frac{d\phi}{dx} = 0$ at $x = 0$

Solution: $\phi(x) = A \cos Bx + C \sin Bx$

$$\frac{d\phi}{dx} = 0 \text{ at } x = 0 \quad \text{forces } C = 0$$

$$\phi\left(\frac{a}{2}\right) = 0 \quad \text{forces } \cos\left(\frac{Ba}{2}\right) = 0$$

Therefore, $\phi(x) = A \cos(B_n x) = A \cos\left(\frac{n\pi}{a} x\right) \quad n = 1, 3, 5, \dots$

B_n are the eigenvalues, $\cos(B_n x)$ are the eigenfunctions

(harmonics) $B_{1,2}$ is the buckling of the fundamental mode $= \left(\frac{\pi}{a}\right)^2$



Examples of Higher Harmonics - Natural Modes of a Slab Reactor

Steady State One-Group Diffusion Equation

$$D\nabla^2\phi - \Sigma_a\phi + s = 0$$

Define $L^2 = \frac{D}{\Sigma_a}$ (Unit cm²)

Since $s = \eta \Sigma_{aF} \phi$ and $f = \Sigma_{aF} / \Sigma_a$

then $s = \eta f \Sigma_a \phi = k^\infty \Sigma_a \phi$

$$\nabla^2\phi + \frac{k^\infty - 1}{L^2}\phi = 0$$

Define $B^2 = \frac{k^\infty - 1}{L^2}$

then $\nabla^2\phi + B^2\phi = 0$



Auxiliary Calculation Modules (Con't)

- ***RIPPLE**
Creates a new fundamental mode from current fluxes, and stores data in "FLUX MODES" "LATESTFUND" for subsequent use by *FLUXMAP

- ***MAPMATRIX**
Creates flux-mapping matrices for a specified flux mode N, and stores data in "FLUX MODES" "MODE N" for subsequent use by *FLUXMAP; If N=1, data stored in "FLUX MODES" "LATESTFUND"



Auxiliary Calculation Modules (con't)

- ***READAMODE**
Reads specific mode previously created, and stores data in “FLUX MODES” “MODE N” for subsequent use by *FLUXMAP; If N=1, data stored in “FLUX MODES” “LASTESTFUND”

- ***ONLINEMAT**
Calculates and copies to a specified file all mapping matrices for subsequent use by the on-line flux-mapping program at Point Lepreau



Sample *MONIC Input

```
*START      ALAN GRAY
480  SEU    0.9  %   U235  FUEL      MONIC  THIRD  HARMONIC  -  1ST  AZIMUTHAL  B

*MODEL      480  SEU    0.9  %   U235  FUEL      MONIC  THIRD  HARMONIC  -  1ST  AZIMUTHAL  B

*READ TAPESEUMONIC02
*MONIC
A          10  1      0.05      20  100  30
E  2  15 3243600.0 0.95470  0.00001  1.5      0.99  0.05      1  10 600
GUESS      1  1      1.0  422.9  3.832  270.0
LABEL      MONIC THIRD HARMONIC - 1ST AZIMUTHAL B
*RITE TAPESEUMONIC03*
TITLE      480  SEU    0.9  %   U235  FUEL      MONIC  THIRD  HARMONIC  -  1ST  AZIMUTHAL  B

*STORE
FROM      HARMONICS HARMONIC 3SLOW FLUX CELL PHI
TO        FLUX/POWERSLOW FLUX CELL PHI
*PRINT    CELL PHI
*RITE CARD
BLOCK     FLUX/POWERSLOW FLUX CELL PHI
FORMAT    (12E12.5)
WRITE     CELLPHI03      1      5760
*PRNT MASS
*CLOSE    NORMAL TERMINATION
```



Flux Mapping

The aim of flux mapping is to determine the amplitudes A_n to obtain the best fit of the mapped fluxes to the measured fluxes F_d .

There are many more detectors than modes, i.e., $D > N$. In the CANDU 6 there are 102 in-core detectors, i.e., $D=102$, and the number of modes used in the flux-mapping expansion, N , ranges between 15 and 28.

Since it is impossible to obtain a perfect fit to D detector fluxes using a smaller number N of unknowns A_n , the flux-mapping method obtains a least-squares fit of the mapped fluxes to the measured fluxes F_d .



Solving for the Mode Amplitudes

Working through the algebra in Matrix Notation:

$$A = \begin{pmatrix} A_1 \\ \vdots \\ A_N \end{pmatrix} \quad (7)$$

$$F = \begin{pmatrix} F_1 \\ F_2 \\ \vdots \\ F_D \end{pmatrix} \quad (8)$$

$$M = \begin{pmatrix} M_{11} & M_{12} & \dots & M_{1N} \\ & M_{21} & & \\ & \vdots & & \\ M_{D1} & M_{D2} & \dots & M_{DN} \end{pmatrix} \quad \text{where } M_{dn} = \psi_n(\bar{r}_d) \quad (9)$$

$$W = \begin{pmatrix} W_1 \\ \vdots \\ W_D \end{pmatrix} \quad (10)$$



Solving for the Mode Amplitudes (con't)

The detector measurements (readings) are electric currents generated by the in-core detectors and the lead cables. The readings converted to effective fluxes by dividing by sensitivity factors:

$$F_d = \frac{E_d}{S_d} \equiv K_d E_d \quad d = 1, \dots, D \quad (3)$$

where E_d is the reading detector d .

$K_d \equiv \frac{1}{S_d}$ is the inverse sensitivity of detector d and F_d is the derived “measured flux” (also sometimes called the “calibrated flux”) for detector d .



Solving for the Mode Amplitudes (con't)

Define a sum of squares of differences between the mapped and measured fluxes:

$$\epsilon = \sum_{d=1}^D W_d^2 \{ \phi_d - F_d \}^2 \quad (4)$$

where W_d is the weighting assigned to the Detector d .

Using Equation (1) for the mapped fluxes,

$$\epsilon = \sum_{d=1}^D W_d^2 \left\{ \sum_{n=1}^N \psi_n(\bar{r}_d) A_n - F_d \right\}^2 \quad (5)$$

ϵ is minimized by imposing the condition for an extremum:

$$\frac{\partial \epsilon}{\partial A_n} = 0 \quad n = 1, \dots, N \quad (6)$$



Solving for the Mode Amplitudes (con't)

Condition imposed by Eq. 6 leads to:

$$M^T \cdot (W \cdot W^T) \cdot M \cdot A = M^T \cdot (W \cdot W^T) \cdot F \quad (11)$$

Inverting this equation, the amplitude vector is obtained as

$$A = H \cdot F \quad (12)$$

where the NxN “pseudo-inverse” matrix H is given by:

$$H = \{(M^T \cdot (W \cdot W^T)) \cdot M\}^{-1} \cdot (M^T \cdot (W \cdot W^T)) \quad (13)$$

Once the modes-at-detectors matrix M has been computed and the weight vector W has been chosen, the matrix H can be calculated by inversion (Equation 13), and the amplitudes A_n can be determined by a simple matrix multiplication, Equation (12).

If $w = 1$, then Equation 13 reduces to : $H = (M^T \cdot M)^{-1} \cdot M^T$ (14)



Choice of Weighting

- **Choice arbitrary in principle**
- **Uniform Weighting - $W=1$**
 - **Equal absolute errors**
 - **Matrix H can be pre-calculated**
 - **Computation of A by a single matrix multiplication**
 - **Used in on-line flux mapping**
- **Relative Weighting - $W_d = 1 / F_d$**
 - **Sum of relative errors (percentage errors) minimized**
 - **High flux reading carries relatively smaller percentage error**
 - **Matrix H re-computed every time W changes**
 - **Used in off-line mapping**



Inclusion of Lead-Cable Effects

Detector current is the sum of the currents generated by the detector proper and its lead cable. The Ψ in Eq. (1) should then be replaced by the effective quantities.

$$\Psi_n^{\text{total}}(\vec{r}_d) = \Psi_n^{\text{detector}}(\vec{r}_d) + \alpha_d \cdot \Psi_n^{\text{(lead-cable d)}}$$

**where $\Psi_n^{\text{detector}}(\vec{r}_d)$ is the average flux in mode n at detector d
 $\Psi_n^{\text{(lead-cable d)}}$ is the lead-cable flux, summed over lengths of cable equal to the modelled length of the detector, α_d is the sensitivity of unit length lead cable relative to the detector d sensitivity.**

The remainder of the flux-mapping methodology is unchanged.



Three Dimensional Flux Distribution

Using the Modal Amplitudes A, 3-D flux distributions can be obtained:

$$\phi(\vec{r}) = \sum_{n=1}^N A_n \Psi_n(\vec{r})$$

With Lattice-Cell Modes (Bundle-Flux Modes Matrix B,

$$B_{kn} = \Psi_n(\vec{r}_k) \quad \begin{array}{l} k = 1, \dots, N_B \\ n = 1, \dots, N \end{array}$$

Mapped Lattice-Cell Flux is given by:

$$\begin{aligned} \phi_k \equiv \phi(\vec{r}_k) &= \sum_{n=1}^N B_{kn} A_n \\ &= \{B \cdot A\}_k \end{aligned}$$

Fuel Flux is deduced from Mapped Cell Flux by:

$$\phi_{k, \text{fuel}} = \phi_k \cdot F_k(\omega_k)$$



Three Dimensional Flux Distribution (con't)

With Channel-Flux-Modes Matrix C,

$$C_{jn} = \sum_{k \text{ in channel } j} B_{kn} \quad \begin{array}{l} j = 1, \dots, N_c \\ n = 1, \dots, N \end{array}$$

Mapped Channel Flux is given by:

$$\begin{aligned} \phi_j &= \sum_{n=1}^N C_{jn} A_n & j &= 1, \dots, N_c \\ &= \{C \cdot A\}_j \end{aligned}$$



Zone-Average Thermal Fluxes

With Zone-Average-Modes Matrix Z,

$$Z_{in} = \frac{\sum_{k \text{ in zone } i} B_{kn}}{\sum_{k \text{ in zone } i} 1} \quad \begin{array}{l} i = 1, \dots, N_z \\ n = 1, \dots, N \end{array}$$

Mapped Zone-Average Flux is given by:

$$\begin{aligned} \phi_i &= \frac{1}{N_{BZ_i}} \sum_{n=1}^N Z_{in} A_n \\ &= \frac{1}{N_{BZ_i}} \{Z \cdot A\}_i \quad i = 1, \dots, N_z \end{aligned}$$

where N_{BZ_i} is a number of bundles in Zone i



Mapped Powers

- **The mapped bundle power is obtained from the mapped cell flux by:**

$$P_k = \phi_k \cdot H_k(\omega_k)$$

where $H_k(\omega_k)$ is the H-factor of bundle k at instantaneous irradiation ω_k , ϕ_k is the mapped thermal flux for bundle k

Absolute normalization of the fluxes and powers is then imposed from the assumed the total reactor thermal power P_{th} , i.e., by renormalizing all fluxes such that:

$$\sum_{k=1}^{N_B} P_k = P_{th}$$



Mapped Powers (con't)

- **Note: Only F-Factors and H-Factors are required for deducing the power distribution and for fuel irradiation and burnup increments. Device incremental F and H need to be included.**



Failed Detectors

- **Detector Reading deemed irrational if less than 0.05 or greater than 1.5 of average value**
 - **Exclude irrational detectors in the mapping calculation: smaller matrix dimension, re-calculate the pseudo-inverse H Matrix**
 - **Replace failed detector reading by best-estimate, given by the fundamental mode detector coupling coefficient M_{d1} properly normalized**
- **(Absolute) difference between mapped and measured detector fluxes larger than an acceptable range (e.g. 3-sigma)**
 - **Multiple Passes: 1st pass with best-estimate, subsequent passes with mapped value, check for failed detectors after each pass (up to 4 passes)**



Comparison to Vanadium Fluxes (1992 Lepreau Restart Power Run-up Transient)

<u>Simulation Method</u>	<u>RMS Difference</u>
Pre-Simulation (Diffusion)	3.0%
Standard Diffusion	2.4%
History-Based Diffusion	2.2%
Mapping with Standard Diffusion Solution as Fundamental	1.7%
Mapping with History-Based Diffusion Solution as Fundamental	1.7%



Auxiliary Calculation Modules

- ***DLSENSIT**
 - **Calculates lead cable relative sensitivity factors for a given FPD**
 - **Detector sensitivity is a function of accumulated irradiation, same for lead cable**
 - **Irradiation by time-average flux for FPD full power days assumed**

- ***MONIC**

Computes higher harmonics of the diffusion equation. Modal Mesh fluxes for Mode N stored in Index “HARMONICS” “FUNDAMENTAL” or “HARMONICS” “HARMONIC N”



Auxiliary Calculation Modules (con't)

- ***ORTHOG**
Orthogonalizes a set of flux modes, using *MONIC orthogonalization procedure



Core-Tracking Application

- ***FLUXMAP Calculation Requires**
 - Updated fundamental Flux Mode (if desired)
 - Harmonic modes specification
 - Raw Vanadium detector readings
 - Detector sensitivity factors
- ***SIMULATE (POWERMAP Option) Calculation Requires**
 - Updated F-Factors and H-Factors at each bundle position
 - * Updated fuel irradiation (), fuelling bundle movements, and current lattice cell conditions (*POWDERPUF calculations may be required).



Core-Tracking Application (con't)

- **Incremental F-Factor and H-Factor due to devices
(current Zone fills required)**
- **Mapped Channel Powers and Bundle Powers used in
Power Limit Compliance statistics**